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**The KEDAK Program  
Compendium  
Part VI  
Mechanized Transfer of  
Nuclear Data from ENDF/B  
to KEDAK and vice versa**

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Projekt Schneller Brüter

**Kernforschungszentrum Karlsruhe**

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KERNFORSCHUNGSZENTRUM KARLSRUHE GMBH

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Part VI

Mechanized Transfer of Nuclear Data  
from ENDF/B to KEDAK and vice versa

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## Abstract

Translation of nuclear data from one representation into another one permit programs, which are linked to a specified data base, to use also data out of another data base.

The programs BRIGITTE and COPEND, used for translating from ENDF/B to KEDAK, will be described in detail. The description includes source program listings. Translation from KEDAK to ENDF/B can be done with the program KTOE. The report contains a copy of the NEA (Nuclear Energy Agency) CPL (Computer Programme Library) abstract. The programs BRIGITTE (abstract no. NEA 438) and KTOE (abstract no. NEA 342) can be obtained from the NEA CPL.

## Das KEDAK Programm Compendium

### Teil VI

#### Mechanisierte Übertragung von Nuklearen Daten von ENDF/B nach KEDAK und umgekehrt

### Zusammenfassung

Die Übertragung von nuklearen Daten aus einer Darstellung in eine andere erlaubt es in Programmen, die an eine Datenbasis angeschlossen sind, auch Daten aus einer anderen Datenbasis zu verwenden.

Die zur Übertragung von ENDF/B nach KEDAK verwendeten Programme BRIGITTE und COPEND werden ausführlich beschrieben. Die Beschreibung enthält auch Quellprogrammlisten. Die Übertragung von KEDAK nach ENDF/B kann mit dem Programm KTOE erfolgen. Der Bericht enthält eine Kopie der NEA (Nuclear Energy Agency) CPL (Computer Programme Library) Kurzbeschreibung. Die Programme BRIGITTE (abstract no. NEA 438) und KTOE (abstract no. NEA 342) können von der NEA CPL bezogen werden.

### Acknowledgement

The authors of this report are grateful to Mrs. K. Mayer for typing this report with great carefulness and patience. Besides typing Mrs. K. Mayer established reproducible graphical representations out of sketches of flowcharts for programs.

The authors are also indebted to Dr. H. Küsters for his critical reading of this report.

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Section 1.

Some remarks about the  
exchange of nuclear data

E. Stein

Normally the representation of nuclear data, used to exchange these data, is different from that one used to work with these data. This must be done, because using the exchange data set, which is in nearly all cases sequentially written on a magnetic tape, is ineffective, if a lot of random accesses are made. It is evident, that the translation from the exchange representation of the data to the local library and vice versa, should be rather simple. E.g. the translation program must not be changed, if new data types are defined.

Unfortunately the described state is not at all realized. There are existing - due to historical reasons - different nuclear neutron data libraries. The data of these libraries can only be exchanged easily, if the receiving library is of the same type. The difficulties result from the different conventions used when storing the same physical data. One solution of this problem could be to use world wide one set of conventions when representing the same data, that means to use one library. But this can be done only with a very great effort of work, because all programs, which are working with the data of a specified library, are using the special conventions of this library. That means, that all these programs then had to be changed.

A second solution of the problem is to translate the data from one representation into another and vice versa. This does not only mean to represent the data in such a manner that they can be read by the retrieval routines of the other library, but also to use all conventions of the other library in the translation of the data.

So far we have made only general considerations about the problem. In the following we will restrict to the exchange of data between the libraries ENDF/B (see /2/ and /3/) and KEDAK (see /4/, /5/, and part II of this KEDAK Program Compendium). Because the data in KEDAK are mainly intended to be used for the calculation of fast reactors, only a subset of the ENDF/B data types must be exchanged. To enable the programs, which are working with KEDAK, an access to this subset of ENDF/B-data, the program BRIGITTE - described in section 2. - has been written. A change of all the programs would have been much more difficult.

If - in the future - KEDAK users will be interested in datatypes, which are not processed by BRIGITTE until now, the data can be used directly from

ENDF/B, or the program BRIGITTE has to be extended. An extension of the program BRIGITTE should be made, if the existing programs can work with the translated data after performing only minor program changes. If processing of the new data types requires new programs or extensive program changes, the data should be directly read from ENDF/B. In this case the data should not be read from the card image form of ENDF/B, but from a dataset to be created on a direct access device, which allows random access, without sequential processing.

Until now only one way - ENDF/B to KEDAK - of data exchange has been treated. The tools for the other direction are described in section 4. Until now, we have not done any translation from KEDAK to ENDF/B. But it is planned in the near future to offer KEDAK materials, which have been evaluated in Karlsruhe, also in an ENDF-representation. This will be done by using the program, which is described in section 4., or, if this is not possible, by changing this program, or by developing own methods.

But it should be kept in mind, that a card image representation of the KEDAK library exists. This card image representation is used to exchange the data between KEDAK libraries on different installations. But it can also be used to print the data, or if retrieval programs exist<sup>+)</sup> , to process the data. However, due to the sequential representation, this is not an optimal way.

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<sup>+)</sup> A retrieval program package for the KEDAK card image representation, which will have the same subroutine calls as the LDFPAC (see part III), will be developed in the near future.

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Section 2.

BRIGITTE

A program for translating nuclear data  
in ENDF/B-representation to nuclear data  
in KEDAK-representation

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---

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## 2.1 Introduction

The program BRIGITTE translates a selection (see 2.2.2) of ENDF/B (see /2/ and /3/) data types into the KEDAK-representation (see /4/, /5/ and II.1 of this compendium). The necessity to do this has been explained in section 1. The present version of the program solves the problem in a nearly automatic way. But because the translated data must be controlled and in few cases changes must be made, we prefer to use the term "mechanized transfer" instead of automatic transfer.

Several people have developed the program BRIGITTE. The first version was written by J.C. Schepers (see /1/) from CEN at MOL (Belgium). The present running version of the program at GfK in Karlsruhe has been developed out of the program written by J.C. Schepers by making corrections, improvements and extensions. This work has been influenced by some important proposals and ideas from P. Vandeplass, from the CEN at MOL. P. Vandeplass improved some parts of the program in the version of BRIGITTE, running at MOL.

This report describes the Karlsruhe version of BRIGITTE only. To understand this report, the user should be familiar to some extent with ENDF/B and KEDAK. For ENDF/B, one may read the reports /2/ or /3/; for KEDAK, the preceding parts of this Compendium (mainly II.1 and sections of III) or the reports /4/ and /5/ can be used.

The following information about the program BRIGITTE starts with a guide for the program user. The recommendations given in 2.2 should be obeyed, when reading this section. The user's guide is followed by a programmer's guide (2.3), where one can find all information, which is necessary for program changes or for the location of program errors. After this section one can find the reference listings (2.4) followed by the source program. Section 2.5 contains a list of the source programs together with some aids to find one's way in this listing.

At the end of this introduction, we would like to point to section 2.2.4, where the normal way of inserting data into a public KEDAK library is described. This section will explain the necessity of the program COPEND, which is described in 3.

## 2.2 User's guide

If one likes to run the program BRIGITTE, one may read the following sections until 2.2.10 and additionally 2.3.1 and 2.3.3, and then start the program. But if one wants to have a BRIGITTE output at once and are only reading 2.2.10, we must give a warning for using a KEDAK-library produced by or processed with KEMA (see part II of this Compendium or /4/) as direct acceptor (dataset linked to FTO1FOO1) for the converted data. Before running the program, one should in any case read 2.2.4.

### 2.2.1 Differences between the ENDF and the KEDAK representation

In this section the expense of work, which must be done by the program BRIGITTE, will be shown. This work results from the differences between the ENDF and the KEDAK representations of the same data.

(1) ENDF-conventions (see /2/ and /3/)

(α) ENDF is built up of mainly four record types: CONT, LIST, TAB1, and TAB2. To represent a specified cross section type these records are composed in a defined manner. But this composition can depend from the value of a quantity in a first record. The type and sequence of records can differ from datatype to datatype and must be given for all datatypes in a written description.

The data in ENDF can be retrieved in units of records. A TAB1-record can contain e.g. the pointwise representation of the total cross section in the whole energy range.

(β) ENDF uses five interpolation schemes. By this the number of stored data can be reduced, but problems occur when combining the ENDF-data (e.g. addition of two pointwise cross sections).

- ( $\gamma$ ) ENDF did not contain compound cross sections like the transport cross section .
  - ( $\delta$ ) Often a parametric representation is chosen in ENDF instead of a pointwise representation; e.g. in the resonance region only parameters plus a correction cross section is given.
- (2) KEDAK conventions (see /4/, /5/ and part II of this Compendium):
- ( $\alpha$ ) KEDAK is built up of data items. A data item can have  $N$  arguments ( $N \geq 0$ ) and  $M$  functional values ( $M \geq 0$ ). An example for a data item is with  $N = 1$  and  $M = 1$  an energy value together with a cross section value, or with  $N = 3$  and  $M = 8$  an item describing a resolved resonance. The data in KEDAK are retrieved in units of data items. A number of data items is composed to a data type, e.g. a cross section in the whole energy range or a set of all resonance information for a specified material. This datatype can have one or more so called "further names". In the case of the inelastic excitation cross sections, the excitation energy is a further name. To each further name or each combination of further names a set of data items, which have all the same number  $N$  of arguments and the same number  $M$  of functional values, is associated.
  - ( $\beta$ ) KEDAK uses only linear interpolation.
  - ( $\gamma$ ) Programs, which are using KEDAK, expect to find a number of compound cross sections, like the transport or the absorption cross section.
  - ( $\delta$ ) Some cross sections must not be given only in parametric form on KEDAK, but also in a pointwise representation, because users of KEDAK and above all, programs working with KEDAK are expecting these data.

By (1)( $\alpha$ ) and (2)( $\alpha$ ) we can see that each datatype of ENDF with a different structure must be processed with a different method. But this is not only a problem of the program BRIGITTE but also a problem in a normal retrieval of data from ENDF/B. Points (1)( $\beta$ ) and (2)( $\beta$ ) shows that converting from



ENDF/B to KEDAK means the introduction of only linear interpolation. This can be done by generating enough points between the existing ones.

Due to (1)( $\gamma$ ) and (2)( $\gamma$ ) the compound cross sections must be produced in BRIGITTE. This is only done for the transport and the absorption cross section.

The differences mentioned in (1)( $\delta$ ) and (2)( $\delta$ ) give rise, in the case of the resonance parameters, to a lot of computer work, because the calculation of pointwise cross sections out of resolved resonance parameters is very CPU-time consuming in the case of many resonances.

### 2.2.2 Datatypes converted by the program

In the following table all datatypes, that can be converted by the version of the program described here, are listed. The ENDF identification (see /2/, /3/) is followed by the associated KEDAK alphanumerical name (see /4/, /5/, II). Some data types, which are given in ENDF/B by different formulations, as in the case of energy distributions of secondary neutrons, can only be converted for some formulations. This is also mentioned in the following table. If the ENDF MT-number is enclosed in "(", that datatype is no longer used in ENDF/B. That means this datatype must be generated during conversion. A "-" sign in the MT-column means: this KEDAK-datatype did not exist in ENDF/B. Remember: In the case of SGT,SGN,SGF and SGG the resolved and unresolved resonance part - if existent - is added to the appropriate pointwise cross sections of ENDF/B MF = 3.

ENDF MF	MT	KEDAK alphanumerical name	remarks
1	452	NUE	average total number of neutrons per fission (If given in a polynomial representation, the pointwise representation is calculated in conversion.)
2	151	RANGRES	see /4/, /5/, II
		ISÔT1	" "
		ISOT2	" "

ENDF		KEDAK alphanumeric name	remarks
MF	MT		
2	151	ISOT3	generated in the case of more than 1 isotope. See /4/, /5/, II.
		RES	resolved resonance parameter set
		ST	unresolved resonance parameter set
		STGF	generated in the case of fissionable materials. Contents see /4/, /5/, II
			pointwise cross sections are produced out of the resonance parameters and added to the appropriate data of ENDF MF = 3. Conversion and generation of pointwise cross sections is only done for single and multilevel Breit-Wigner parameter sets
3	1	SGT	total cross section
3	2	SGN	elastic scattering cross section
3	3	SGX	non elastic cross section <sup>1)</sup>
3	4	SGI	total inelastic cross section
3	16	SG2N	(n,2n) cross section
3	17	SG3N	(n,3n) cross section
3	18	SGF	total fission cross section
3	22	SGIA	(n,n' $\alpha$ ) cross section
3	23	SGI3A	(n,n'3 $\alpha$ ) cross section
3	24	SG2NA	(n,2n $\alpha$ ) cross section
3	25	SG3NA	(n,3n $\alpha$ ) cross section
3	(27)	SGA	absorption cross section. Produced in the program. SGA = SGG + SGF + SGP + SGD + SGH3 + SGHE3 + SGALP + SG2HE
3	28	SGIP	(n,n'p) cross section
3	29	SGI2A	(n,n'2 $\alpha$ ) cross section

<sup>1)</sup> The translated cross section SGX does not contain the resonance parts.

ENDF		KEDAK alphanumeric name	remarks
MF	MT		
3	51 . . 90	SGIZ	inelastic cross section for excitation of rest nucleus. (If MT = 51 did not exist, but higher MT-numbers can be found, see 2.2.6, namelist INPUT, variable SALINE.)
3	91	SGIZC	inelastic scattering cross section to the continuum
3	102	SGG	(n, $\gamma$ ) radiative capture cross section
3	103	SGP	(n,p) cross section
3	104	SGD	(n,d) cross section
3	105	SGH3	(n,t) cross section
3	106	SGHE3	(n,He <sup>3</sup> ) cross section
3	107	SGALP	(n, $\alpha$ ) cross section
3	108	SG2HE	(n,2 $\alpha$ ) cross section
3	-	SGTR	transport cross section SGTR = SGT - MUEL * SGN Always calculated in conversion (if MUEL exists)
3	-	ETA	average number of fission neutrons per neutron absorption. ( <u>Not</u> produced in the running version of the program but foreseen in later versions.)
3	-	ALPHA	ratio of capture (SGG) to fission (SGF) cross section. (Foreseen to be produced in later program versions.)
3	251	MUEL	average cosine of the elastic scattering angle in the laboratory system
4	2	SGNC	angular distributions of secondary neutrons for elastic scattering. SGNC and SGNL are tabulated distributions in the center of mass resp. laboratory system. LEGNC and LEGNL are sets of Legendre coefficients. If the ENDF/B data are given in tabulated form, only SGNC resp. SGNL is generated. In the case of Legendre coefficients, the tabulated data are generated.
		SGNL	
		LEGNC	
		LEGNL	
5	16	SED2N	energy distributions of secondary neutrons emitted by the (n,2n) process
5	17	SED3N	energy distributions of secondary neutrons emitted by the (n,3n) process

ENDF		KEDAK alphanumeric name	remarks
MF	MT		
5	18	SEDF	energy distributions of secondary neutrons emitted in the fission process
5	91	SEDIC	energy distributions of secondary neutrons emitted in the inelastic scattering to the continuum
			Only the following ENDF energy distribution laws can now be converted: LF = 3 in K = 4 (excitation of discrete levels) LF = 7 in K = 2 (Maxwellian spectrum) LF = 9 in K = 1 (evaporation spectrum) LF = 10 in K = 2 (Watt spectrum)

### 2.2.3 General description of the program (options, region, time, restrictions)

As described in the preceding sections, the program BRIGITTE translates nuclear data from the ENDF- into the KEDAK-representation. Only a selection of the data types is processed (see 2.2.2).

Provided that a load module of the program exists, one can use the JCL-cards, described in 2.2.9, and the input control cards, described in 2.2.6, to start the program. The construction of a complete job will be much simplified, when using the examples given in 2.2.10.

An exact description of the program flow can be found in 2.3.1. For this section it is sufficient to know, that the translation of the ENDF/B data is made filewise. First some data of MF = 1 are processed and stored in an auxiliary file, then the resonance data (MF = 2) are read, stored and used for the calculation of pointwise cross sections, which are also stored in an auxiliary file. These pointwise data are then combined with the MF = 3 data. The result of this process is then adapted to KEDAK conventions and stored in an auxiliary file. After this some data of MF = 4 and afterwards MF = 5 are processed in a similar mode. In the last section of the program the data, stored on the auxiliary file, are included into a KEDAK-library<sup>\*)</sup>

<sup>\*)</sup> This KEDAK-library must not be created by or processed with KEMA (see 2.2.4)

There are some options in the program, which can be selected by using the quantities in the namelist control input. Processing of data of MF = 4 and/or MF = 5 can be suppressed, or processing only MF = 4 or MF = 5 data can be selected (see IF4 and IF5 in the namelist INPUT, 2.2.6). Only due to historical reasons the data translation can be made with respect to a not longer used KEDAK convention: all inelastic cross sections have the same energy mesh (see KEDAVE in the namelist INPUT, 2.2.6). Until now, only two options, controlled by quantities in the namelist INPUT, have described. To get a total knowledge of all options, one must read section 2.2.6.

Some important options are controlled by variables in the namelist LIB (also described in 2.2.6). One can select, whether the program has to built up a new KEDAK-library, or can use an existing one. It is also possible, to delete in an existing library a material, if it has the same name, as the converted one. But remember: all possible program options are described in 2.2.6.

The region used by the program itself is about 102 k-byte. Some region is needed by the system, for read-input and print-output buffers, and for buffers for other datasets, used by the program. As a mean value we can take about 150 k-byte necessary for the program, for system use, and for buffers. The rest of the region, which is given in the JOB-card, is used for data arrays. The size of the data field region mainly depends on the number of resolved resonance parameters. If there are no such or only a few parameters, 300 k-byte are normally sufficient to run the program. In the case of many resolved resonance parameters, 300 k-byte will not be enough. But we recommend, to start with 300 k-byte. If the program has not enough region, an error message will be produced. This error message will help to decide, how much additional region is necessary.

The CPU (central processing unit)-time needed by the program, also depends on the number of the resolved resonance parameters. This time will be given for an IBM/370 model 168. If there are no parameters, 1 minute is sufficient. For about 80 single level Breit Wigner resonances about 11 minutes are used. The CPU-times will increase up to hours, if there are many multilevel resolved Breit Wigner resonance parameters.

Although the data storage used in the program is variably dimensioned, there are some arrays of a fixed length, which causes some restrictions. These are:

- One (natural) element may have up to 10 isotopes.
- Five different l-states ( $l = 0, 1, 2, 3, 4$ ) are allowed in the resolved Breit Wigner resonance parameter set.
- Three different l-states and 6 different j-states for each l-state are allowed in the unresolved Breit Wigner resonance parameter set.
- One hundred points are allowed as primary energy grid for energy distributions of secondary neutrons.

#### 2.2.4 Insertion of the translated data into a local KEDAK library

The program BRIGITTE produces a KEDAK-library, which differs from the KEDAK-libraries produced by the program KEMA (see II, /4/ and /5/). The BRIGITTE created libraries contain in word 880 of the first record (the word is the last one) the number of the next available record in the library.

This word is used, when inserting new data in the subroutine SICKY (see 2.3.5.86). In the library created by BRIGITTE also the data of each material start on a new record. By this, deleting materials and compressing the library in the direct access form is much simpler than for the KEDAK-libraries created by KEMA (see /4/ resp. II). Because of the differences described here, it is not possible to use a KEDAK-library, which has been created by or processed with KEMA, as direct acceptor for the data converted by the program BRIGITTE. But there are no difficulties for the KEDAK retrieval codes, as described in III, to use the BRIGITTE created libraries. Because all programs, which are working with KEDAK, use these routines, all these programs can work with the "BRIGITTE-created" libraries.

In most cases using of a KEMA created (or processed) library will cause the program to do no insertion of the translated data. But because the present version of the program only tests word 880 of the first record, the decision whether the library is BRIGITTE- or KEMA-created (or processed) may be

incorrect. Therefore we recommend never to use a KEMA created (or processed) library as direct acceptor for the data translated by BRIGITTE.

The only way to insert the translated data in a public KEMA-created (or processed) library is to produce KEMA-input out of the BRIGITTE-created libraries. This problem is solved by the program COPEND, described in 3.

One may ask, why BRIGITTE itself does not produce this KEMA-input instead of a KEDAK-library. The answer is very simple: in many cases it is not necessary to insert the data into the public KEDAK-library. As stated before all retrieval routines, described in III, can work with the BRIGITTE-created libraries. Why then insert the data into one library? There are several reasons for this action:

- (1) If the data should be kept for a longer time interval, it is much simpler to manage one library (e.g. produce back-up copies), than a number of different libraries.
- (2) If a BRIGITTE-created library has been processed by KEMA (e.g. to correct some data), it can not be longer used by BRIGITTE.
- (3) There are programs, which are processing several materials in one run. This materials must be contained in one library, because these programs use the retrieval program III.3, which permits only one KEDAK-library.

A last question may be left. Why BRIGITTE does not write directly in public KEMA-created KEDAK-libraries. There are three answers to this question:

- (1) To have a control about changes in KEDAK, only one program should be allowed to update the library. This is the program KEMA.
- (2) The program KEMA performs some formal checks before inserting the data.
- (3) The program BRIGITTE can delete materials, with names identical to the translated ones, and compress the library after the deletion. This would be very difficult and expensive with KEMA-created (or processed) libraries.

The contents of this section can be summarized in the following statement.  
Never use a KEMA created (or processed) KEDAK-library in a BRIGITTE run as acceptor for the translated data. If the checks in BRIGITTE will fortuitously fail, the KEMA-created (or processed) library may be destroyed.

#### 2.2.5 External data sets

The programm BRIGITTE uses 7(8) data sets represented by the following ddnames (see /10/):

FT06FOO1  
FT05FOO1  
FT01FOO1  
FT09FOO1  
FT10FOO1  
FT11FOO1  
FT12FOO1  
(FT15FOO1) optional

The datasets, which the program expects to be allocated to this ddnames, will be described in the following.

##### FT06FOO1

###### Description:

Print output.

###### Actions taken:

The program makes only WRITE-statements on this dataset. This is done in the subroutines TITEL, PRIEIN, GSPACE, INITNL, INGRID, REGINE, MARY, MARY01, LUCY, RESERR, MILLI, MELA, ANNICK, ANNIO1, DENISE, MARION, SOPHIE, SUZY, ADDPNT, SUPNEG, ISABEL, FILE5, PRIZEI, ERROR, PRCONT, PRLIST, PRTAB1, PRTAB2, PRHOL, PRTPID, WDA, WRECS, SICKY, LISA, TABPRT, PAGE and A8FORM.

###### DD-parameters:

The printer must be able to produce lines with a maximum of 132 characters.

###### Possible errors:

Maximal line length of the printer is smaller than 132 characters.



FT05FO01

Description:

Input unit for (card) control input.

Actions taken:

The data are read from FT05FO01 in the subroutines INITNL, PRIEIN and SICKY. A REWIND is made in PRIEIN.

DD-parameters:

The record format (RECFM) must be fixed (F) or fixed blocked (FB) with a logical record length (LRECL) of 80 bytes.

Possible errors:

The control input does not fulfil the conventions necessary for namelist input (see 2.2.6).

FT01FO01

Description:

KEDAK-library \* (created in the job or old one).

Actions taken:

The converted material is added to the KEDAK-library. If the library does not exist, it is created. If the library exists, READ and WRITE is made. If it is created, only WRITE-statements are executed.

DD-parameters:

A SPACE parameter must also be given for datasets with DISP=OLD, because the values in this parameter are used in the program in the subroutine DEFI (see 2.3.5.93). Because the BLKSIZE must be equal 3520, the parameter can look like SPACE=(3520,(600)).

The DCB parameter must be

DCB=(RECFM=F,BLKSIZE=3520,LRECL=3520,DSORG=DA).

Possible errors:

No SPACE parameter given for a dataset with DISP=OLD.

BLKSIZE not equal 3520.

Number of blocks not sufficient to receive the converted material (use another or a new KEDAK-library \* where the material can be included).

---

\* The KEDAK-library must not be created by or processed with KEMA; it must be created by BRIGITTE itself (see 2.2.4).

FT09FO01

Description:

ENDF/B-data.

Actions taken:

The data are processed by READ-statements. No writing is done. First a REWIND is made in the subroutine INGRID and finally in the subroutine PART1. Reading is done - only in forward direction - in the subroutines SEARCH and PREC. BACKSPACE can be done in the subroutines REGINE, SOHPIE, ISABEL and FILE5.

DD-parameters:

Because normally ENDF/B is used in the MODE=3 representation (see /3/ chapter R-2.2) the record format (RECFM) must be fixed (F) or fixed blocked (FB) and the logical record length (LRECL) must be equal 80 bytes. The data must be in EBCDIC or BCD representation.

Possible errors:

Wrong tape used.

In the case of NL-types: incorrect DCB-specifications.

FT10FO01

Description:

Auxiliary dataset, where temporarily ENDF/B-data are stored (and read) in the MODE=2 (see /3/ Chapter R-2.2).

Actions taken:

The data set is used in the subroutines SOPHIE (only if KEDAVE=0. See 2.2.6 namelist INPUT), ISABEL and FILE5. The data are written by using the subroutine WREC and reread by using the subroutine RREC.

DD-parameters:

Because writing is done, using no FORMAT-statements, the record format (RECFM) must be variable (VS) or variable blocked (VBS). The block size can be chosen with respect to an optimal buffer size. We recommend the values BLKSIZE=1016 or BLKSIZE=2040 or BLKSIZE=3064. Normally a SPACE of about 50 blocks of the size 1016 is sufficient. The dataset can reside on a disk, a tape, a drum, or somethinglike that.

Possible errors:

- Incorrect record format (RECFM).
- Block size (BLKSIZE) too small when using RECFM=VS.
- Space in the SPACE-parameter not sufficient.

FT11FOO1

Description:

Auxiliary dataset. Contains all data to be inserted in a KEDAK-library \* by the subroutine SICKY. The structure of the dataset is described in the subroutine WDA (see 2.3.5.76).

Actions taken:

The data are written on the dataset by the subroutine WDA (see 2.3.5.76) and read by the subroutine RDAB (see 2.3.5.77). Only the subroutine ISABEL writes on FT11FOO1 without using these subroutines. To get the table of contents and addresses, the subroutines WDA and ISABEL are also reading from FT11FOO1 (if initialized). Writing of data on the dataset is done (directly or by the subroutine WDA) in the subroutines PART1 (initialisation), REGINE, MARYO1, ANNIO1, SUZY, ISABEL, and FILE5. WDA is also called by the subroutine PART1 to print the table of contents and addresses of file 11. Reading and processing (insertion in a KEDAK-library \*) the data of file 11 is only done by the subroutine SICKY.

DD-parameters:

The allocated dataset must reside on a direct access storage (e.g. a disk). The blocksize (BLKSIZE) must be 1600 bytes, the record format (RECFM) must be fixed (F). You can give the DCB-subparameter DSORG=DA. In any case the space must be given in units of the blocksize 1600, e.g. SPACE=(1600,(400),CONTIG) (CONTIG means, that the space must be contiguous.) The contents of the SPACE-parameter are used in the subroutine DEFI (see 2.3.5.93).

Possible errors:

- BLKSIZE not equal 1600 bytes.
- SPACE not given in units of 1600 bytes.
- Not enough SPACE given (Error produced is of the form "Record number out of range for .....").

---

\* The KEDAK library must not be processed with or created by KEMA; it must be created by BRIGITTE itself (see 2.2.4).

FT12F001

Description:

Auxiliary dataset. Used to store ENDF/B TAB1- and LIST-records, which have been read from the ENDF/B file, but cannot be processed at the moment. Also data calculated in the program, which will be used later, are stored on FT12F001. The structure of the dataset is described in 2.3.5.78.

Actions taken:

The data are written on the dataset by the subroutine WRECS (see 2.3.5.78) and read from it by the subroutine RRECS (see 2.3.5.79). Writing of data is done in the subroutines PART1 (initialisation), INGRID, MELA, ANNIO1, DENISE and SOPHIE.

The data are read in the subroutines MYRIAM, MARION, SOPHIE and SUZY.

DD-parameters:

The allocated dataset must reside on a direct access storage (e.g. a disk). The blocksize (BLKSIZE) must be 1600 bytes, the record format (RECFM) must be fixed (F). The DCB-subparameter DSORG=DA can be given. Because the contents of the SPACE-parameter are used in the subroutine DEFI (see 2.3.5.93), the space must be given in units of the block-size 1600, e.g.

SPACE=(1600,(400),,CONTIG).

(CONTIG means, that the space must be contiguous)

Possible errors:

BLKSIZE not equal 1600 bytes.

SPACE not given in units of 1600 bytes.

Not enough SPACE given (Error produced is of the form "Record number out of range for ...").

If ITAPE (see 2.2.6 namelist INPUT) is equal 1, the following ddname must exist:

FT15FOO1

Description:

Datatypes of ENDF/B MF=3 with added resonance parts are written in ENDF representation on the dataset.

Actions taken:

Writing (including HEAD- and SEND-records) is done in the subroutine SUZY (see 2.3.5.37) by using the subroutines WCONT (see 2.3.5.50) and WREC (see 2.3.5.51).

DD-parameters:

Because writing is done with MODE=3, the record format (RECFM) must be fixed (F) or fixed blocked (FB), the logical record length (LRECL) must be equal 80 bytes. The dataset can reside on a disk, a tape, a drum, or something like that. The data can also be punched.

Possible errors:

Completion code D37, because not enough SPACE has been given

2.2.6 Input description

The program BRIGITTE is controlled by the contents of three (plus one) namelists. All three namelists must be given in the input, but some can be empty. If more than one material is to be converted in one job, these three namelists must be given for each material.

The namelist input is read from unit FT05FOO1 respectively SYSIN.

The sequence of the namelists is as following:

```
&SPACE  data-list or empty  &END
&INPUT  data-list           &END
{ &ISO   data-list           &END }
  &LIB   data-list or empty  &END
```

In the description of the variables of all namelists, the abbreviations, described in the following, are used.

R4 means REAL\*4 quantities  
R8 means REAL\*8 quantities  
I4 means INTEGER\*4 quantities  
L4 means LOGICAL\*4 quantities

In the "dimension column" either the maximal dimension for arrays or the word scalar for scalar quantities is given.

The values in the "default value column" are used, if the variable is not found in the namelist input. If no default value exists a "-" sign is made.

There are some syntax rules for the namelist input data, which can be found in FORTRAN manuals (see e.g. /9/). The important rules will be listed in the following.

- (1) The first character in each record to be read must be a blank.
- (2) The second character must be an &, immediately followed by the namelist name, which must not contain any embedded blanks.
- (3) The namelist name must be followed by at least one blank.
- (4) After this blank the data items separated by commas are listed. (A comma after the last item is optional.) There may be blanks between the comma and the next data item symbolic (resp. array) name.
- (5) The end of a specified namelist input is signaled by &END.
- (6) The input for one namelist can be given on more than one card. The data item list should end with a comma on one card and can be continued in each column of the next card.

(7) A data item is of the form

symbolic name = constant

or

array name = set of constants  
(separated by commas)

The symbolic name can be an array element name (e.g. A(21))

or a variable name (e.g. V).

(Valid symbolic or array names are given in the following tables).

The constants may be of the type integer (e.g. 5), real (e.g. 3.0E+3), literal (e.g. 'ALPHA'), complex, or logical (e.g. .TRUE., or F).

(8) In the case of a set of constants with the same value, you can write  $k * \text{constant}$ , where  $k$  gives the number of times the value constant is to occur.

(9) If the input for one namelist is distributed over several cards, the cards must not contain a numeration in columns 73 to 80.

Examples for the namelist input are given in 2.2.10.

In the following tables a description of all variables of all namelists is given. Normally only input for the namelist INPUT is necessary, whereas the namelist input for SPACE and LIB can be of the form

&SPACE &END

respectively

&LIB &END

In the case of the namelist INPUT a survey of all necessary quantities is given at the end of the table.

namelist name: SPACE				
name	type	dimension	default value	meaning
NBKX	I4	scalar	12 000	maximal dimension of the array BK (see 2.3.3)
NIX	I4	scalar	500	maximal dimensions of the arrays NBT and JNT (see 2.3.3)
N2X	I4	scalar	5 000	maximal dimensions of the arrays X and Y (see 2.3.3)
NBX	I4	scalar	3 000	maximal dimensions of the array B (see 2.3.3)
IADIM	I4	scalar	20 000	maximal dimensions of the array A (see 2.3.3)
ITDIM	I4	scalar	32 560	maximal dimension of the array TAB (see 2.3.3)

Caution: The default values of the quantities in the namelist SPACE should be only changed by the input, if the user has a sufficient knowledge of the organization of the data storage in the program as described in section 2.3.3. If there are difficulties with the maximal dimension of an array, we recommend to enlarge the total region and to give only "&SPACE &END" with an empty list.



namelist name: INPUT

name	type	dimension	default value	meaning
MODE	I4	scalar	3	input format of ENDF/B-data (see /3/ chapter R-2.2). (MODE=3 means card image tape with standard arrangement.)
LABEL	I4	scalar	-1	<p>&lt; 0 TPID-record (see /3/ chapter R-2.3.6) is not read. The tape identification number is not checked.</p> <p>= 0 TPID-record is read. The tape identification number is printed out but not checked.</p> <p>&gt; 0 TPID-record is read. The tape identification number is printed out and must be equal LABEL.</p> <p>We recommend to give always the correct tape identification number and not to use LABEL-values <math>\leq 0</math>.</p>
MATP	I4	scalar	-	ENDF/B-material number
NAMKDK	R8	scalar	-	alphanumerical KEDAK-name of the material MATP
MATKDK	I4	scalar	-	numerical KEDAK-name of the material MATP
EB	R4	scalar	-	binding energy of the last neutron (in eV)

namelist name: INPUT (continuation)

name	type	dimension	default value	meaning
NOTTRL	R8	(25)	'blank'	<p>data types contained in this array (as KEDAK datatype names), are not translated from ENDF/B to KEDAK. These data types - all sections of ENDF/B MF=3 - can be:</p> <p>SGT, SGN, SGX, SGI, SG2N, SG3N, SGF,            SGIA, SGI3A, SG2NA, SG3NA, SGA, SGIP,            SGI2A, SGIZ, SGIZC, SGG, SGP, SGD,            SGH3, SGHE3, SGALP, SG2HE, SGTR,            ETA, ALPHA, MUEL.</p>
KEDAVE	I4	scalar	1	<p>0 an old not longer used KEDAK version will be created. (All SGIZ and SGIZG will have the same energy grid.)</p> <p>1 the translated data will fulfil the valid KEDAK conventions.</p>
NISO	I4	scalar	0	<p>number of isotopes in the ENDF/B material, for which resonance (and some other) information is given. (NISO=0 and NISO=1 have the same meaning: resonance information is given only for one isotope.)</p> <p>If <u>NISO &gt; 1</u> the <u>namelist ISO must be given</u> in the input</p>

namelist name: INPUT (continuation)

name	type	dimension	default value	meaning
LNER	I4	scalar	0	<p>used in the subroutine MARY (see 2.3.5.12). Maximum number of points per resonance used in calculating the energy grid in the resolved resonance region with the subroutine EPT (see 2.3.5.17).</p> <p>0 51 points per resonance are used</p> <p>≥1 LNER points per resonance - with a maximum of 200 - are used</p>
MLSLSW	I4	scalar	0	<p>≥1 single level Breit Wigner formalism is used in the resolved resonance region to calculate pointwise cross sections independently from the ENDF/B recommended formalism.</p> <p>=0 the ENDF/B recommended formalism (single or multilevel Breit Wigner) is used</p>
ELINEG	I4	scalar	0	<p>Deletion of negative SGT and SGN cross section values.</p> <p>≠1 negative SGT and SGN cross section values are printed out</p> <p>=1 the values are printed out and deleted</p>

namelist name: INPUT (continuation)

name	type	dimension	default value	meaning
IF4	I4	scalar	0	<p>control of calling the subroutine ISABEL (Processing of ENDF/B MF=4 data; angular distributions of secondary neutrons. See 2.3.5.42)</p> <p>= 1 only the subroutine ISABEL is called in BRIGITTE (only angular distributions of secondary neutrons are converted)</p> <p>=-1 the subroutine ISABEL is not called in BRIGITTE (angular distributions of secondary neutrons are <u>not</u> converted)</p> <p>≠ 1 and ≠ -1 normal processing of BRIGITTE (ENDF/B MF=4 data are included in the conversion process)</p>
IF5	I4	scalar	0	<p>control of calling the subroutine FILE5 (Processing of ENDF/B MF=5 data: energy distributions of secondary neutrons. See 2.3.5.44).</p> <p>= 1 and IF4≠1 only the subroutine FILE5 is called (only energy distributions are converted)</p> <p>=-1 and IF4≠1 the subroutine FILE5 is not called in BRIGITTE (energy distributions of secondary neutrons are not converted)</p> <p>≠ 1 and ≠ -1 and IF4≠1 normal processing of BRIGITTE (ENDF/B MF=5 data are included in the conversion process)</p>

namelist name: INPUT (continuation)

name	type	dimension	default value	meaning
ICOMB	I4	scalar	0	<p>used in the subroutine COMBP (see 2.3.5.67)</p> <p>≠ 1 the optimal interpolation scheme is chosen by the routine</p> <p>= 1 only a linear-linear interpolation scheme is used</p> <p><u>Caution:</u> This variable should have always the value 0 in production runs</p>
EPSRES	R4	scalar	1.OE-2	<p>accuracy used in the following cases:</p> <ul style="list-style-type: none"> <li>- energy grid used in the calculation of pointwise cross sections of resolved resonance parameters. It is guaranteed that a cross section value between two energy points can be predicted by linear interpolation with a relative error less or equal EPSRES</li> <li>- in the calculation of pointwise cross sections out of unresolved resonance parameters the energy grid created guarantees that each of the statistical resonance parameters can be linearly interpolated with an error less or equal EPSRES</li> <li>- in changing the interpolation methods used in ENDF/B to linear-linear. It is guaranteed that functional values between two argumental values can be predicted by linear interpolation with</li> </ul>

namelist name: INPUT (continuation)

name	type	dimension	default value	meaning
				<p>an error less or equal EPSRES. In some cases the program increases the given value of EPSRES temporarily up to <math>2.0 \cdot 10^{-2}</math></p> <ul style="list-style-type: none"> <li>- in the combination of two pointwise cross sections with non linear interpolation rule. It is guaranteed, that cross section values between two energy points can be predicted - using the associate interpolation rule - with an error less or equal EPSRES</li> <li>- in cropping unnecessary points, before inserting the data in KEDAK, functional values, which can be predicted by linear interpolation with an error less or equal <math>\epsilon</math>, are cropped. Normally <math>\epsilon</math> has the value SCREPS (see next variable). But if two adjacent points fulfil the condition <math>\left  \frac{x_2 - x_1}{x_1} \right  &lt; 5.0 \cdot 10^{-6}</math> <math>\epsilon</math> is increased to the value EPSRES.</li> </ul> <p>(EPSRES should always be greater than SCREPS). (see 2.3.5.75)</p>

namelist name: INPUT (continuation)

name	type	dimension	default value	meaning
				<p>- in the construction of scratch classes (see 2.3.5.75) in the subroutine DZMSCH (see 2.3.5.40). In this case 10 classes are built up with the boundaries 0, EPSRES, <math>EPSRES + \frac{1.-EPSRES}{9.0}</math>, <math>EPSRES + 2 \cdot \frac{1.-EPSRES}{9.0}</math>, ..... , <math>EPSRES + 8 \cdot \frac{1.-EPSRES}{9.0}</math>, 1.0.</p> <p>Points, which are replaced by DZMSCH, are checked, with which accuracy they could have been interpolated. This accuracy defines the membership in a specified scratch class. The number of points in the scratch classes is printed out. EPSRES is also used in the construction of scratch classes in the subroutine CROP (see 2.3.5.75). In this subroutine it has the name EPSMAX.</p>
SCREPS	R4	scalar	$5.0 \cdot 10^{-3}$	<p>Before inserting the pointwise cross section into KEDAK and after combination of two cross sections it is checked, whether there are points, which can be predicted by interpolation between adjacent points to relative accuracy of <math>\epsilon</math>.</p>

namelist name: INPUT (continuation)

name	type	dimension	default value	meaning
				<p>Points with an <math>\epsilon</math> better than or equal to SCREPS (resp. EPSRES, see description of this quantity) are deleted. SCREPS and EPSRES (see variable before) are also used in the subroutine CROP to built up - for printing purposes - so called scratch classes (see 2.3.5.75 "problem solved").</p>
FAICOD	R4	scalar	10.0	<p>used in combining two pointwise cross sections in the subroutine COMBP (see 2.3.5.67 "problem solved" point (5)).</p> <p>The variable FAICOD should <u>not</u> be changed by the user. Only program users, who have a very well knowledge of the program BRIGITTE, can try to work with this variable.</p>



namelist name: INPUT (continuation)

name	type	dimension	default value	meaning
LEDIT1	I4	scalar	1	control of print output = 1 descriptive data and dictionary of ENDF/B (MF=1, MT=451) are <u>printed</u> . ≠ 1 these data are <u>not</u> printed.
LEDIT2	I4	scalar	1	control of print output ≠ 1 the data, listed in the following, are <u>not</u> printed. = 1 the data, listed in the following, are <u>printed</u> . - resolved resonance parameters in ENDF and KEDAK representation - unresolved resonance parameters in ENDF representation - the KEDAK data types ST and STGF (for fissionable materials)
LEDITS	I4	scalar	0	Control of print output ≠ 1 the data, listed in the following, are <u>not</u> printed. = 1 the data, listed in the following, are <u>printed</u> . - pointwise cross sections calculated out of resolved and unresolved resonance parameters - pointwise total cross sections calculated out of resolved and unresolved resonance parameters - ENDF/B (pointwise) reaction types of MF=3

namelist name: INPUT (continuation)

name	type	dimension	default value	meaning
BIB12	I4	scalar	1	number of records of the length 400 4-byte words used for the table of contents and addresses of the direct access data set NF=12 (see 2.3.5.78). The maximal value for BIB12 is 10.
ITAPE	I4	scalar	0	control of writing the converted data of ENDF/B MF=3 on unit 15. The data written in the subroutine SUZY (see 2.3.5.37) are neutron cross sections of ENDF/B file 3 with added resolved and unresolved resonance contributions. The interpolation scheme is linear-linear and all unnecessary points have been deleted (all conversion work of this type has been done). But writing is done in ENDF mode=3 (card image). ≠ 1 <u>no</u> writing is done = 1 writing is done
SALINE	L4	scalar	.FALSE.	used in the subroutine SOPHIE (see 2.3.5.36). Normally the program assumes, if in ENDF/B MF=3 is no section MT=51, that there are also no sections 52,53,...,90. But there are some materials with special data types stored for example under MT=54 with no MT=51,52,53 existing. To be able to convert such cross section types, SALINE is used. = .FALSE. if MT=51 is not found, no search for MT=52,53,...,90 is made. = .TRUE. if MT=51 is not found, a search is started to find higher MT-values.

namelist name: INPUT (continuation)

In the namelist INPUT there are quantities, which must be given, should be given, must be given only for special cases, and should not be given. These quantities are:

(1) quantities, which must be always given in the input:

MATP,NAMKDK,MATKDK,EB

(2) quantities, which should be given in the input:

LABEL

(3) quantities, which must be given in the input only for special cases:

MODE (ENDF/B-data not in card image)

NOTTRL (suppress translation of some datatypes)

KEDAVE (old KEDAK conventions wished)

NISO (information for more than one isotope given in the resonance region)

MLSLSW (always single level Breit Wigner desired)

ELINEG (suppress negative  $\sigma_t$  and  $\sigma_e$  cross section values)

IF4 (control of processing angular distributions of secondary neutrons)

IF5 (control of processing energy distributions of secondary neutrons)

EPSRES (change of the accuracy)

SCREPS (" " " " )

LEDIT1

LEDIT2 } (more or less print output desired)

LEDITS }

BIB12 (table of contents too small. In this case BRIGITTE produces an error message and a STOP)

SALINE (no MF=3 MT=51 but higher MT occurs)

(4) quantities which should not be given in the input:

LNER (if too small, not enough or too much points may be produced)

ICOMB } (should only be used after a very good knowledge of the program

FAICOD }

ITAPE (there is normally no use for the produced dataset)

namelist name: ISO (Only given for NISO > 1. See namelist INPUT)

name	type	dimension	default value	meaning
NAMISO	R8	(10)	-	alphamerical KEDAK-names for the additional isotopes of a material (NISO-names) (see 2.2.10)
MATISO	I4	(10)	-	numerical KEDAK-names - associated to the alphamerical ones - for the additional isotopes of a material (NISO-names) (see 2.2.10)
EBISO	R4	(10)	-	binding energies of the last neutron for the NISO additional isotopes (NISO-values). (EB, in the namelist INPUT, means for NISO > 1 the binding energy of the last neutron of the natural element.)

namelist name: LIB

name	type	dimension	default value	meaning
IFIRST	I4	scalar	0 <sup>1)</sup>	<p><math>\leq 0</math> the KEDAK library <sup>2)</sup> is existing. The converted material has to be added to this library.</p> <p><math>\geq 1</math> it is assumed that the direct access data set contains no KEDAK library. The KEDAK library is initialized.</p> <p><u>Caution:</u> Existing libraries are destroyed by IFIRST <math>\geq 1</math></p>
IMOD	I4	scalar	0 <sup>1)</sup>	<p>= 0 the converted material is not yet present in the library. It is to be inserted. (If it is present no insertion is done.)</p> <p><math>\neq 0</math> material yet present in the library. It is to be replaced. (If it is not present, a warning message is printed and the converted material is inserted in the KEDAK-library.)</p> <p>The quantity IMOD is only meaningful for IFIRST <math>\leq 0</math></p>
IDAT	I4	scalar	date <sup>1)</sup> fetched by the routine DATUM (see 2.3.5.98)	<p>creation (or change) date of the KEDAK-library in the form ddmmyy specifying day, month, year (see /4/ or II). In nearly all cases the date fetched by the subroutine DATUM should be used.</p>

1) Before each reading of the LIB namelist input these default values are set.

2) This library must not be created by or processed with KEMA; it must have been created by BRIGITTE in a previous run (see 2.2.4).

### 2.2.7 Output description

The printed output, produced by the program, is almost self explaining. Therefore only a general survey together with some helps for some details, where questions may arise, will be given. The production of some parts of the printed output is controlled by the values of the variables LEDIT1, LEDIT2, and LEDITS. These variables have default values but may also be given in the input (see 2.2.6 namelist INPUT).

The printed output of the program BRIGITTE starts with the program name in "big" letters. After this a list of the control input is printed. This list is followed by the results of the dynamic working space allocation (see 2.3.3). On the next page a printout of the contents of the namelist INPUT is given. After this the program name together with the material name and some other informations are printed.

Under the heading "ENDF/B FILE 1" the results of the conversion of the ENDF/B MF=1 data are printed. In this section first the printout of the not translated section MT=451 is given. This may be followed by a listing of the  $\bar{\mu}$ -data.

The next heading is "ENDF/B File 2". If desired, resolved and unresolved parameters in ENDF/B- and KEDAK-representation are printed together with the calculated pointwise cross sections. All pointwise cross sections are listed in the ENDF/B MODE=4 (see /3/ chapter R 2.2.4). At the end of the printout of the pointwise total cross sections, calculated from resolved resonance parameters by adding the partial cross sections, the accuracy EPS used in combining the partial cross sections is printed out.

After this the text "FILE 3 - SMOOTH CROSS SECTIONS" is printed starting on a new page. Under this heading the cross sections read from ENDF/B MF=3 are listed. If there is a resonance part, which has been calculated from the data of MF=2, this part has been added before printing. The printout of all these cross sections is controlled by the quantity LEDITS (see 2.2.6 namelist INPUT). The printout also includes the calculated cross sections as e.g.  $\sigma_a$  and  $\sigma_{tr}$ . In the case of  $\sigma_a$  the partial cross section names, used in calculating  $\sigma_a$ , are listed.

After changing the interpolation rule to linear-linear, deletion of points, which can be predicted by linear interpolation, combination of points with nearly equal energy values, and suppressing of negative cross section values, all cross sections read from MF=3 (including the calculated ones) are printed once more. These data will be inserted into the KEDAK-library. Each listing of a cross section type is headed with some information. In the first line the ENDF/B MT-number is listed together with the used accuracies. The "PRECISION EPS" is used in changing the interpolation rule, "EPSMIN" in scratching points, which can be predicted by linear interpolation, and "EPSMAX" in scratching points, which can be predicted by linear interpolation and are having nearly equal energy values. This first line is followed by two lines, which will show the number of points, which have been scratched due to the possibility of linear interpolation. These points are classified in so called "scratchclasses". If e.g. 38 points are in the scratchclass 0.2 % to 0.3 % (1 % means 1 percent), this means, that these scratched points could have been predicted with an accuracy in the range 0.2 % to 0.3 %.

If there are some points with nearly equal energy values, which have endured the scratching process, these points will be combined. Some information about this combination will be printed. The accuracy classes have the same meaning as the described scratchclasses.

If there are negative  $\sigma_t$  or  $\sigma_e$  cross section values, the confounded energy ranges will be printed. If desired (see ELINEG in the namelist INPUT in 2.2.6), the values will be suppressed. All this information is followed by the KEDAK cross section type name printed in "big" letters and finally the data itself. In order to know where the computing time is spent, an asterisk surrounded text is printed. This text includes the date, the time of the day, and the consumed CPU-time (CPU abbreviates Central Processing Unit). This information may be found very often in the program printout.

The next heading is "FILE 4 - ANGULAR DISTRIBUTIONS OF SECONDARY NEUTRONS". This part ends with "REACTION TYPE TABLES". The first column in these tables means the energy. In the "polynomial" table, which is the first one, if a polynomial representation exists, the second and the third column means internal addresses. In the "pointwise" table, which is normally the second one, but may also be, if no polynomial representation exists, the first one, the second column means the number of cosine values. The third and fourth column are internal addresses.

The next heading is "FILE 5 - ENERGY DISTRIBUTIONS OF SECONDARY NEUTRONS". If a secondary energy distribution law (used in ENDF/B) can not be processed by the program, an one line text is printed, which may be overlooked, if the printoutput is inspected only very carelessly. We recommend, at least in the beginning, to read the program printoutput very carefully. All data type names, which will be included into KEDAK, are listed under the heading "CONTENTS OF FILE 11". In the first column the numerical KEDAK material name can be found. The second column means the numeration of the isotope (normally 1). In the third column the alphameric KEDAK datatype name is listed. The fourth column means the record number in file 11, where the data starts. The number of these data can be found in column five, which is the last one.

The section headed with "INSERTION OF TRANSLATED DATA INTO KEDAK" starts with a printout of the namelist LIB, followed by some messages, resp. warnings. After this the first words of record 1 of the KEDAK-library are listed, followed by the material conversion table, the type conversion table and the material address table (see II.1). After this the type address table for the inserted material is listed. The last line of this section is

NEXT AVAILABLE RECORD IN THE KEDAK LIBRARY

followed by a number. This number must not be greater than the extents (plus 1) of the library measured in 3520 byte records. The last section is a "PRINTOUT OF THE KEDAKTABLES". The general data of the KEDAK-library, the material conversion table, the type conversion table, the material address table, and the type address table are listed in an edited form. The last line of the printoutput of the program is

"\*\*\*\*\* NORMAL END OF TABPRT \*\*\*\*\*"

We recommend, to check whether all addresses in the KEDAK-tables are within the library extent.

#### 2.2.8 Listing of possible errors

In this section a listing of all possible not self explaining error messages, which are printed by the program, is given.



Messages generated by the operating system or input/output routines will not be listed here, because these messages must be self explaining. Messages relating to auxiliary datasets see section 2.2.5.

The error messages produced by the program have all the same format. First the text

ERROR STOP      lmn

with lmn to be found in the following tables, is printed. Then the values of MAT, MF, and MT in the Common RECS are printed out followed by the text

\*\*\*DUMP OF THE ARRAYS NBT,JNT,X,AND Y.\*\*\*

These arrays, which are a part of the data storage (see 2.3.3), are normally used to store a TAB1-record (see /3/ chapter R-2.3.3). After the printout of the contents of the arrays a STOP 1 is made.

In the list on the following pages first the error number lmn is tabulated. Then the names of the subroutines, where this error can happen, are specified. Afterwards the reason for the occurrence of the error number lmn is given. In the last column proposals for possible corrections are made.

In nearly all cases the numbered error stops are caused by exceeding the maximal dimensions of an array. That means, increasing the total region or the maximal dimension of an array, will correct the error. But there are also other errors. Therefore it is better, before increasing the region, to look at the following table.

The table of the numbered error messages is followed by the table "Possible corrections for the error type program or ENDF/B-data error". In the last part of this section some hints for other error situations are given. The remarks (1) and (2) will be found at the end of the numbered error messages.

error number	happened in subroutine (s)	reason	possible correction(s)
99	RREC	ENDF/B-tape number NT less or equal zero	Program or ENDF/B-data error. See special table at the end of this section
100	RREC	JT out of range 1-6	
101	RREC	MODE out of range 1-3	
102	RREC	Temperature T not in range given in data	
103	RRECS FETCH	Maximal dimensions of the arrays NBT and JNT too small.	Increase total region or N1X in the namelist SPACE (see 2.2.6). See also remarks ① .
	RREC	Maximal dimensions of the arrays NBT and JNT too small or incorrect ENDF/B data (N1 less or equal zero).	In the first case, which is normally the reason for this error, increase the total region or N1X in the namelist SPACE (see 2.2.6). See also remarks ① . In the second case, correct the ENDF/B-data, which are read by the subroutine.
104	RRECS FETCH	Maximal dimension of the array B too small	Increase the total region or NBX in the namelist SPACE (see 2.2.6)
	RREC	Maximal dimension of the array B too small or incorrect ENDF/B-data (N1 less or equal zero)	In the first case, which is normally the reason for this error, increase the total region or NBX in the namelist SPACE (see 2.2.6). See also remarks ② . In the second case, correct the ENDF/B-data, which are read by the subroutine.
105	MELA ANNICK ADDPNT RRECS FETCH	Maximal dimension of arrays X and Y too small	Increase total region or N2X in the namelist SPACE (see 2.2.6). See also remarks ① .

error number	happened in subroutine(s)	reason	possible correction(s)
105 (continuation)	RREC	Maximal dimension of the arrays X and Y too small or incorrect ENDF/B-data (N2 less or equal zero).	In the first case, which is normally the reason of the error, increase the total region or N2X in the namelist SPACE (see 2.2.6). See also remarks ①. In the second case correct the ENDF/B-data read by the subroutine.
106	RREC	Improper temperature dependence	Program or ENDF/B-data error. See special table at the end of this section
107	RREC	Incorrect MAT,MF,MT on the ENDF/B-tape	Correct the ENDF/B data
108	WREC	JT out of range 1-6	Program or ENDF/B-data error. See special table at the end of this section
109	WREC	MODE out of range 1-4	
133	TERP1	Interpolation code is less than 1 or greater than 5	Program or ENDF/B-data error. See special table at the end of this section
134	DANI TERP1	-	
135	TERP1	Lower and upper end of the interpolation interval have the same value	
201	INGRID	Material MATP not found in specified ENDF/B library	Correct MATP in the namelist INPUT (see 2.2.6) or give correct ENDF/B library
202	INGRID	$\bar{\nu}$ -data (number of neutrons per fission, MF=1,MT=452) not found for the specified material MATP although the material is flagged as fissionable and MT=452 is listed in the table of contents.	Correct the ENDF/B data

error number	happened in subroutine(s)	reason	possible correction(s)
203	LUCY	I out of range 2-3	Program or ENDF/B-data error. See special table at the end of this section
204	ANNICK ANNIO1 REGINE	Maximal dimension of array BK too small	Increase total region or NBKX in the namelist SPACE (see 2.2.6)
205	INGRID	Maximal dimension of array A too small	Increase total region or IADIM in the namelist SPACE (see 2.2.6)
206	MARY MARYO1	Maximal dimension of array BK too small	Increase total region or NBKX in the namelist SPACE (see 2.2.6)
207	MARY	Number of l-states, for which resolved resonance parameter sets are given, is greater than 5.	Can only be corrected by program changes if checking of ENDF/B-data leads to NLS > 5. (Due to ENDF conventions NLS must be less or equal 3).
208	MARY MILLI	Maximal dimension of array A too small	Increase total region or IADIM in the namelist SPACE (see 2.2.6)
209	MARCEL	Negative l-state number	Check ENDF/B resolved resonance data (quantity L). If correct, inform program author. See also special table "Possible corrections for the error type program or ENDF/B-data error".
210	REGINE	Number of isotopes given in ENDF/B MF=2 MT=151 data greater 10	Inform program author to increase the appropriate dimensions in the program
219	REGINE	Number of isotopes given in ENDF/B MF=2 M=151 not consistent with the number given in the control input	Correct NISO in the namelist INPUT and rerun the job
220	ANNICK	Maximal dimension of array A too small	Increase total region or IADIM in the namelist SPACE (see 2.2.6).

error number	happened in subroutine(s)	reason	possible correction(s)
221	ANNICK	Maximal dimension of arrays X and Y too small	Increase total region or N2X in the namelist SPACE (see 2.2.6). See also remarks ① .
222	ANNICK	Maximal dimension of array A too small	Increase total region or IADIM in the namelist SPACE (see 2.2.6)
224	ANNIO1	-	Check unresolved resonance parameter set on consistence. If correct, see special table "Possible corrections for the error type program or ENDF/B-data error".
225	ANNIO1	-	Check unresolved resonance parameter set on consistence. If correct, see special table "Possible corrections for the error type program or ENDF/B-data error".
226	ANNIO1	Maximal dimension of array BK too small	Increase total region or NBKX in namelist INPUT (see 2.2.6).
300	STORE	JT out of range 1-6	Program or ENDF/B-data error. See special table at the end of this section
301	STORE	Identification number MA is equal zero	
310	COMB	Lower energy limit greater or equal upper energy limit	
311	COMB	Identification of the records to be combined is zero	
312	CROP	Incorrect interpolation tables NBT and JNT.	
313	CROP	Maximal dimensions of the arrays NBT and JNT too small	Increase total region or NIX in the namelist SPACE (see 2.2.6). See also remarks ① .
314	IPDS	Incorrect interpolation tables NBT and JNT	Program or ENDF/B-data error. See special table at the end of this section

error number	happened in subroutine(s)	reason	possible correction(s)
317	SEARCH	MODE less 1 or greater 3	Correct MODE in the namelist INPUT (see 2.2.6). If correct, see special table "Possible corrections for the error type program or ENDF/B-data errors"
400	INGRID	MODE out of range	MODE must be 1,2, or 3 (see 2.2.6 namelist INPUT variable MODE)
401	INGRID	ENDF/B type identification number does not match with variable LABEL	Correct variable LABEL in the namelist INPUT (see 2.2.6)
403	MYRIAM	Maximal dimension of the arrays X and Y too small	Increase total region or N2X in the namelist SPACE (see 2.2.6). See also remarks ① .
404	MYRIAM	-	Program or ENDF/B-data error. See special table at the end of this section
407	MYRIAM	Maximal dimension of array A too small	Increase total region or IADIM in namelist SPACE (see 2.2.6)
408	MARION	Maximal dimension of array A too small	Increase total region or IADIM in the namelist SPACE (see 2.2.6)
409	MARION	Maximal dimensions of the arrays X and Y too small	Increase total region or N2X in the namelist SPACE (see 2.2.6). See also remarks ① .
410	MARION	-	Program or ENDF/B-data error. See special table at the end of this section
500	SOPHIE	Maximal dimension of the array A too small	Increase total region or IADIM in the namelist SPACE (see 2.2.6)
501	SOPHIE	Maximal dimension of the arrays X and Y too small	Increase total region or N2X in the namelist SPACE (see 2.2.6). See also remarks ① .

error number	happened in subroutine(s)	reason	possible correction(s)
502	SOPHIE	-	Program or ENDF/B-data error. See special table at the end of this section
505	SOPHIE	Maximal dimension of the arrays X and Y too small	Increase total region or N2X in the namelist SPACE (see 2.2.6). See also remarks ① .
506	SOPHIE	Maximal dimension of the array BK too small	Increase total region or NBKX in the namelist INPUT (see 2.2.6).
510	SOPHIE	Maximal dimension of the array BK too small	Increase total region or NBKX in the namelist SPACE (see 2.2.6).
515	SUZY	Maximal dimension of the array BK too small	Increase total region or NBKX in the namelist SPACE (see 2.2.6).
516	WDA	Maximal dimension of array BK too small	Increase total region or NBKX in the namelist SPACE (see 2.2.6).
550	WDA	Too many entries in the table of contents and addresses of file 11	Probably program error. Rerun the job. See also special table "Possible corrections for the error type program or ENDF/B-data error. If the error persists, reduce the number of datatypes converted by using NOTTRL in the namelist INPUT (see 2.2.6). In any case inform the author of the program
600	ANNICK	Number of j-states for a particular l-state for statistical resonance parameter sets is greater than 6.	Check ENDF/B data (unresolved resonance parameter set, quantity NJS). If NJS is less than or equal to 6 inform program author, else program changes must be done due to changes in ENDF conventions.
601	ANNICK	Number of l-states in a statistical resonance parameter set > 3.	Check ENDF/B data (unresolved resonance parameter set, quantity NLS). If NLS is $\leq 3$ inform program author, else program changes must be made due to changes in ENDF conventions.

error number	happened in subroutine(s)	reason	possible correction(s)
620	DANT	Changing interpolation to linear-linear, 300 points had not been enough to represent a non linear interval	Check whether this is correct. If true increase the arrays in the common RE and the number 300 in the subroutine DANI. Inform the program author
621	CHGINT	Interpolation code less 1 or greater 5 found in the array JNT	Program or ENDF/B-data error. See special table at the end of this section
700	ISABEL	Maximal dimension of array BK too small	Increase total region or NBKX in the namelist SPACE (see 2.2.6).
701	ISABEL	Too many entries in the table of contents and addresses of file 11	See error number 550
703	ISABEL	Maximal dimension of array A too small	Increase total region or IADIM in the namelist SPACE (see 2.2.6).
750	FILE5	Maximal dimension of the arrays X and Y too small	Increase total region or N2X in the namelist SPACE (see 2.2.6). See also remarks ① .
751	FILE5	Maximal dimension of array ENERG in FILE5 too small	Check ENDF/B data of MF=5, which are to be converted, on correctness. If no errors found, increase NCMAX and dimension of the array ENERG in the subroutine. In any case inform the author of the program.
752	FILE5	Maximal dimension of the array BK too small	Increase total region or NBKX in the namelist SPACE (see 2.2.6).
777	MILLI MARION MYRIAM ELIMZ SUZY DZMSCH ADDPNT SUPNEG CHGINT CROP		Program or ENDF/B-data error. See special table at the end of this section



error number	happened in subroutine(s)	reason	possible correction(s)
799	ISABEL	Maximal dimension of array A too small	Increase total region or IADIM in the namelist SPACE (see 2.2.6)
800	ITABI	Incorrect interpolation tables (arrays NBT and JNT).	Program or ENDF/B-data error. See special table at the end of this section
850	SUZY	Maximal dimension of the arrays X and Y too small	Increase total region or N2X in the namelist SPACE (see 2.2.6). See also remarks ① .
851	SUZY	Maximal dimension of the arrays X and Y too small	Increase total region or N2X in the namelist SPACE (see 2.2.6). See also remarks ① .

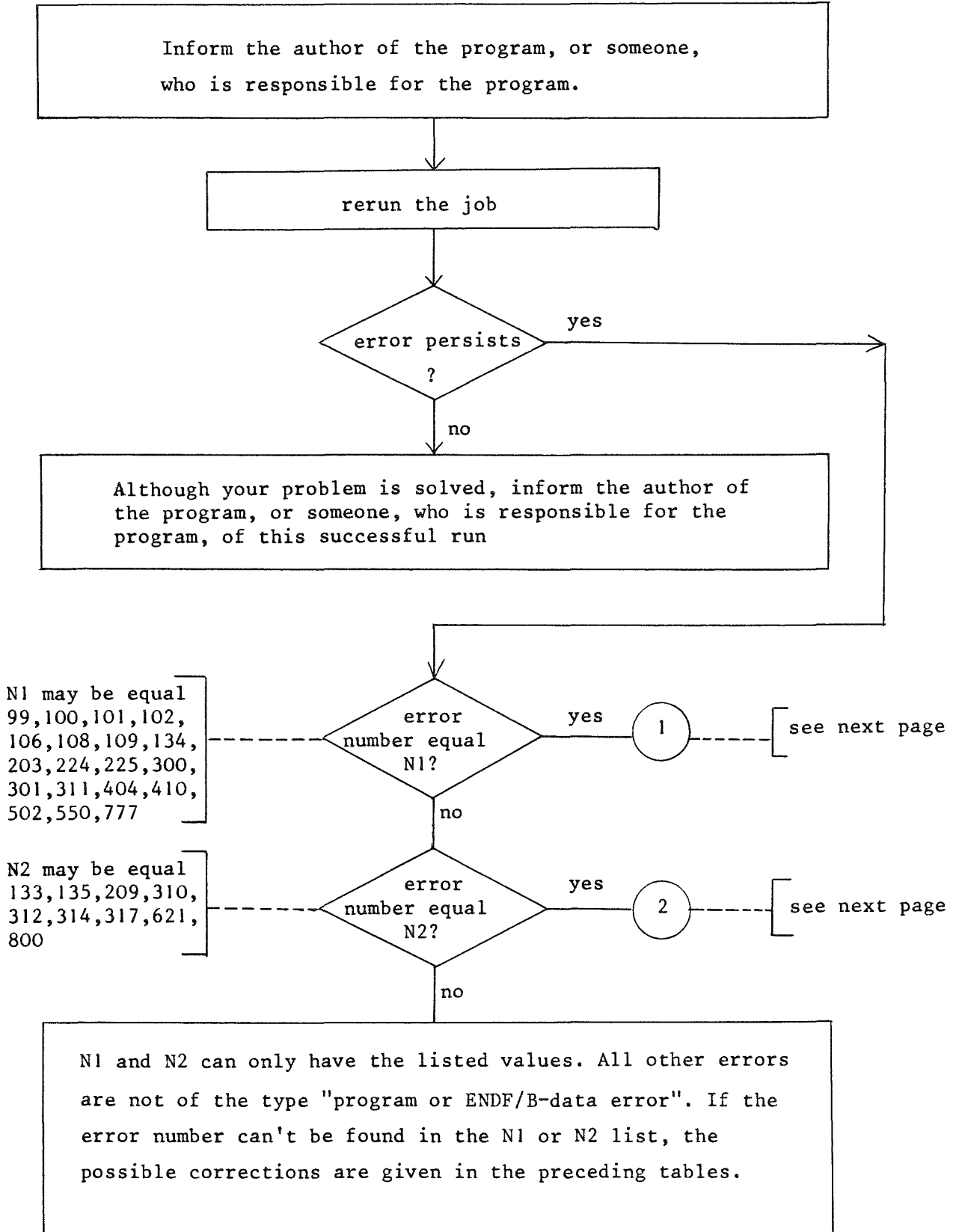
①

Before increasing the total region or maximal dimensions of arrays, look at the printout of the error message. There is a part headed with "\*\*\*DUMP OF THE ARRAYS NBT,JNT,X, and Y\*\*\*". In this part the values of N1 and N2 can be found. Are these values correct? If not, correct the ENDF/B-data.

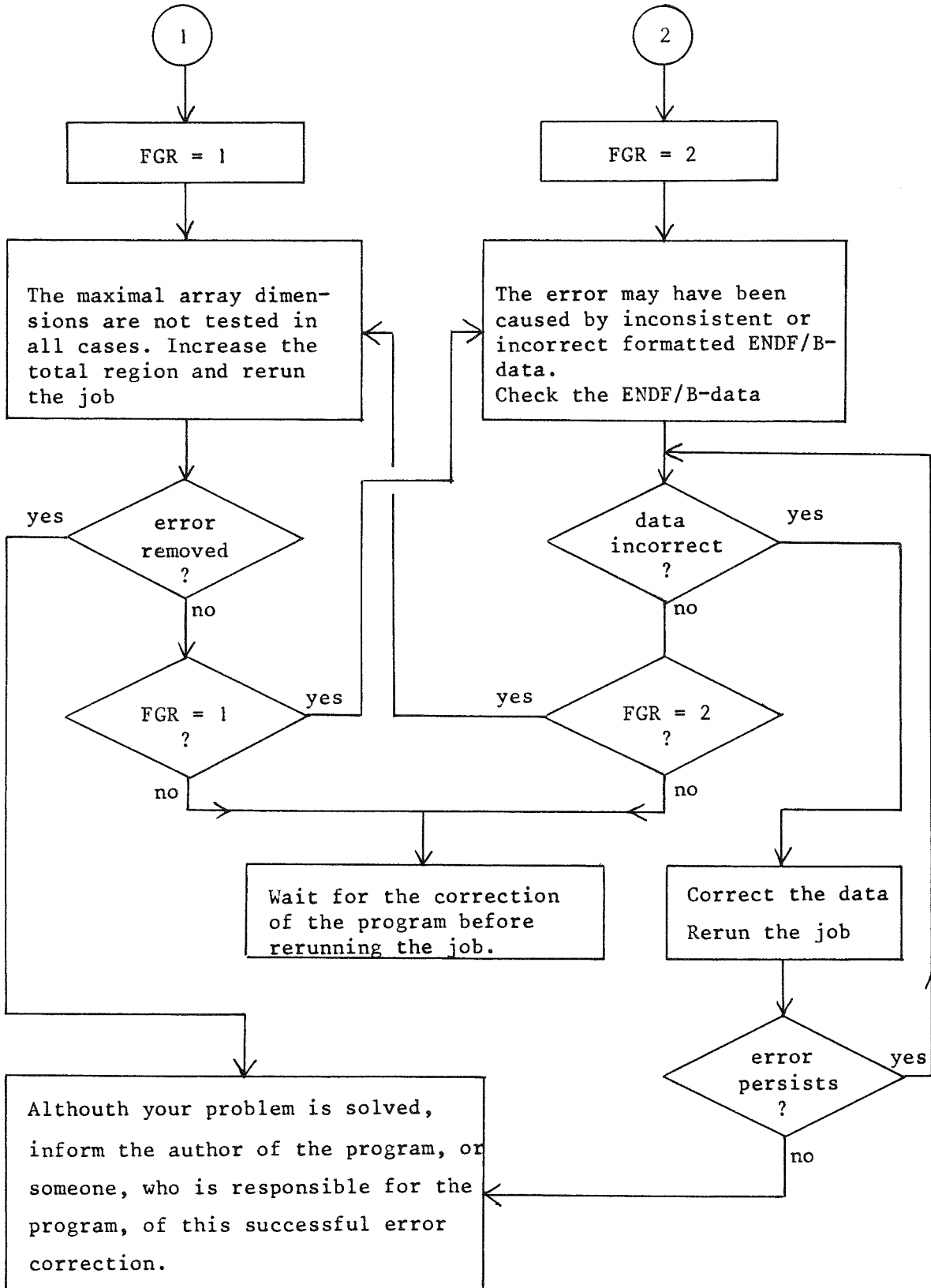
②

Before increasing the total region or maximal dimensions of arrays, check - if possible -, whether the ENDF/B-LIST-record, which is processed, is correct. If not, correct the ENDF/B-data.

Possible corrections for the error type  
"program or ENDF/B-data error".



Possible corrections for the error type "program or ENDF/B-data error" (continuation of the preceding page).



Unnumbered error messages

In the case of an error message produced by system input output routines, which can not be explained, check the following:

- SPACE-parameter given for FT01FO01 (see 2.2.5)?
- SPACE-parameter for FT01FO01 in units of 3520 bytes?
- SPACE-parameter for FT11FO01 and FT12FO01 in units of 1600 bytes?

In case of the error message

RECORD NUMBER N OUT OF RANGE ON UNIT M

with M equal 11 or 12, increase the appropriate SPACE. If M is equal 1, give another or a new KEDAK-library.

### 2.2.9 JCL-cards for an IBM 05370

To run the program BRIGITTE on an IBM computer of the series 360 or 370, which are controlled by an operating system OS360 respectively OS370, the following JCL-cards are necessary:

- (1) // (job card information),
- (2) // REGION=nnnn K, TIME=(minutes)
- (3) // EXEC PGM=progname
- (4) //STEPLIB DD DSN=dsname, UNIT=type,
- (5) // VOL=SER=volname, DISP=SHR
- (6) //SYSABEND DD DUMMY
- (7) //FT05FOO1 DD DDNAME=SYSIN
- (8) //FT06FOO1 DD UNIT=(CTC,,DEFER), LABEL=(.,NL),
- (9) // DCB=(LRECL=133, BLKSIZE=1995, RECFM=FBA)
- (10) //FT01FOO1 DD UNIT=typek, VOL=SER=volnamek,
- (11) // DISP=(dispka, dispkn) DSN=dsnamek,
- (12) // DCB=(RECFM=F, BLKSIZE=3520, LRECL=3520, DSORG=DA),
- (13) // SPACE=(3520, (pq))
- (14) //FT09FOO1 DD UNIT=typee, VOL=SER=volnamee,
- (15) // DISP=(dispea, dispen),
- (16) { // LABEL=(ln, la, , IN), }
- (17) { // DCB=(RECFM=recfm, BLKSIZE=blksize, LRECL=lrecl, DEN=den), }
- (18) // DSN=dsnamee
- (19) //FT10FOO1 DD UNIT=types,
- (20) // SPACE=(blk10, (pq10, sq10)),
- (21) // DISP=(NEW, DELETE, DELETE),
- (22) // DCB=(RECFM=recfm10, BLKSIZE=blk10)
- (23) //FT11FOO1 DD UNIT=types,
- (24) // SPACE=(1600, (pq11), , CONTIG),
- (25) // DISP=(NEW, DELETE, DELETE),

```
(26) // DCB=(DSORG=DA,RECFM=F,BLKSIZE=1600)
(27) //FT12FOO1 DD UNIT=types,
(28) // SPACE=(1600,(pg12),,CONTIG),
(29) // DISP=(NEW,DELETE,DELETE),
(30) // DCB=(DSORG=DA,RECFM=F,BLKSIZE=1600)
(31) { //FT15FOO1 DD parameter }
(32) //G.SYSIN DD *
           control input(see 2.2.6)
(33) //
```

Explanation of the parameters in the listing of the JCL-cards:

We have used the rule, that information, which must be given by the user, is written with underlined lowercase characters. The other characters must be given as typed.

card (1)

(job card information) all necessary information to **start** a job  
(see /10/).

card (2)

nnnn region size. You should start with a value  
of 300.

minutes estimated time value in minutes. You should start  
with values of 1 to 10 minutes.

card (3)

progname name of the program in the partitioned dataset  
dsname (see card (4)). For GFK-users progname =  
BRIGITTE.

card (4)

dsname name of the partitioned dataset containing the  
program named  
progname (see card (3)). For GFK-users  
dsname = STEINPDS

type unit containing dsname. For GFK-users type=3330.

card (5)  
volname name of the volume containing dsname (see card (4)).  
For GFK-users volname=KAPROS.

card (10)  
typek unit containing the KEDAK-dataset<sup>+)</sup> dsnamek (see  
card (11)). Must be a direct access device.

volnamek name of the volume containing the KEDAK-dataset<sup>+)</sup>   
dsnamek (see card (11)). Must be a direct access  
device.

card (11)  
dispka Must be NEW for new datasets, or OLD for old  
datasets.

dispkn Must be KEEP if the KEDAK-library<sup>+)</sup>  should be kept.  
If it should be deleted, it can be DELETE.

dsnamek dataset name of the KEDAK-library<sup>+)</sup> .

card (13)  
pq extension of the KEDAK-library measured in units  
of the BLKSIZE=3520. This card must also be given  
for old datasets.

card (14)  
typee unit containing the ENDF/B-data. (Normally a tape  
unit).

volnamee name of the volume containing the ENDF/B-data  
(Normally a tape).

---

<sup>+)</sup> Never use a KEDAK-library created by or processed with KEMA to receive the translated data. This library must have been created by BRIGITTE itself. (see 2.2.4).

card (15)

dispea In the case of ENDF/B-data stored on a tape  
dispea=OLD. In other cases dispea=SHR may be  
better.

dispen In the case of ENDF/B-data stored on a tape  
dispen=PASS. In other cases the subparameter  
dispen can be omitted.

card (16)

This card is only given for tapes.

ln number of the label.

la SL for standard label tapes.  
NL for non label tapes.

card (17)

This card must be normally only given for non label tapes.

recfm record format of the dataset

blksize block size of the dataset

lrecl logical record length of the dataset

den density, with which the data have been written.

card (18)

dsname dataset name of the ENDF/B-dataset.  
(Not used for non label tapes).

card (19)

types unit, which will contain, the auxiliary dataset.  
For GFK-users types=SYSDA.

card (20)

blk10 block size for the auxiliary dataset.  
Recommended values see 2.2.5.

pq10 primary space quantity for the auxiliary dataset.  
Recommended values see 2.2.5.



sq10 secondary space quantity for the auxiliary dataset. Can be one tenth of pq10.

card (22)

recfm10 Must be VS or VBS.

blk10 see card (20)

card (23)

types see card (19).

card (24)

pq11 primary space quantity for the auxiliary dataset. Recommended values see 2.2.5.

card (27)

types see card (19).

card (28)

pq12 primary space quantity for the auxiliary dataset. Recommended values see 2.2.5.

card (31)

These cards must only be given, if ITAPE is equal 1 (see 2.2.6 namelist INPUT). If the data written on FT15F001 should be punched, the correct parameters for the puncher must be given. If the data should be written on a tape or a disk or something like that, also the correct parameters must be given. See 2.2.5 for DCB-parameters.

card (32)

control input see 2.2.6

GFK-users can use the local procedures and must use additionally ASP control cards. Complete jobs to run BRIGITTE at the GFK computers are listed in 2.2.10.

### 2.2.10 Examples (input and JCL) for the use of the program

In this chapter three examples will be given for the use of the program. All these examples contain the JCL-cards and ASP-cards, which are necessary to run the jobs on the GFK computer installation. Because the installation and the systems can change, the cards are only valid for the existing installation and the operating systems at the time of writing this report.

In all three examples, it is assumed, that the ENDF/B-data can be found on a magnetic tape and are written with MODE=3 (card image standard arrangement see /3/ chapter R-2.2.3). The specifications of the tape, where the ENDF/B-data to be converted can be found, must be asked at the author of this report or the local ENDF/B-manager.

#### Example 1

We assume, that Cm-244 is to be translated. The KEDAK-library is created in the job. A maximal output is wished, and negative  $\sigma_t$  and  $\sigma_e$  cross section values in the resolved resonance region will be deleted.

```
(1) //      (job card information),
(2) //      REGION=300K,TIME=3
(3) /*MAIN  LINES=2
(4) /*SETUP  DEVICE=TAPE9,ID=(ididid,NORING,,SL)
(5) /*FORMAT PR,DDNAME=FTO6FOO1,OVFL=ON
(6) //      EXEC  FHG,NAME=BRIGITTE
(7) //G.STEPLIB DD  UNIT=3330,VOL=SER=KAPROS,DSN=STEINPDS,DISP=SHR
(8) //G.FTO1FOO1 DD  UNIT=uuuu,VOL=SER=volvol,
(9) //      SPACE=(3520,(200)),DISP=(NEW,KEEP),
(10) //      DSN=dsndsn,
(11) //      DCB=(RECFM=F,BLKSIZE=3520,DSORG=DA,LRECL=3520)
(12) //G.FTO9FOO1 DD  UNIT=TAPE9,VOL=SER=ididid,
```

```
(13) // DSN=dsname,DISP=(OLD,PASS),
(14) // LABEL=(n,SL,,IN),
(15) // DCB=(RECFM=FB,LRECL=80,BLKSIZE=3520,DEN=2)
(16) //G.FT10FOO1 DD UNIT=SYSDA,SPACE=(3520,(30,5)),
(17) // DISP=(NEW,DELETE,DELETE),
(18) // DCB=(RECFM=VS,BLKSIZE=3520)
(19) //G.FT11FOO1 DD UNIT=SYSDA,SPACE=(1600,(400),,CONTIG),
(20) // DCB=(DSORG=DA,RECFM=F,BLKSIZE=1600)
(21) //G.FT12FOO1 DD UNIT=SYSDA,SPACE=(1600,(400),,CONTIG),
(22) // DCB=(DSORG=DA,RECFM=F,BLKSIZE=1600)
(23) //G.SYSIN DD *
(24) &SPACE &END
(25) &INPUT MATP=mmmm,NAMKDK='CM244EN4',
(26) MATKDK=o960244,EB=6.799E+6,
(27) LEDITS=1,LABEL=111,ELINEG=1,
(28) &END
(29) &LIB IFIRST=1,&END
(30) /*
(31) //
```

Remarks to the cards:

card (1)

(job card information) These data are jobname, JOB-card identification, accounting information, and name of the user.

card (4)

ididid the tape name must be asked at the local ENDF/B-manager.

card (5)

By this card the maximal printed lines for a page is limited to 66.  
The next line(s) is (are) printed on a new page.

card (8)

uuuu } unit type and volume name of a  
volvol } disk storage.

card (10)

dsndsn this name can be chosen by the user. The translated data are written on this data set. (Never use a KEDAK-library, which has been created by or processed with KEMA. The library must have been created by BRIGITTE itself. See 2.2.4)

card (12)

ididid see card (4)

card (13)

dsname this name must be asked at the local ENDF/B-manager.

card (14)

n the label number must be asked at the local ENDF/B-manager.

card (25)

mmmm the ENDF/B material number must be asked at the local ENDF/B-manager.

card (27)

l11 the label identification, where the material mmmm can be found, must be asked at the local ENDF/B-manager.

When writing the namelist input cards, remember the rules listed in 2.2.6.

Example 2

Np-237 is to be translated. The result is to be inserted in an existing KEDAK library, which must have been created by BRIGITTE in a preceding job (see 2.2.4). If this library contains Np-237-data, these data are to be deleted. A maximal output is wished; negative  $\sigma_t$  and  $\sigma_e$  cross section values in the resolved resonance region are to be deleted.

```
(1) //      (job card information),
(2) //      REGION=700K,TIME=17
(3) /*MAIN  LINES=4
(4) /*SETUP  DEVICE=TAPE9,ID=(ididid,NORING,,SL)
(5) /*FORMAT  PR,DDNAME=FTO6FOO1,OVFL=ON
(6) //      EXEC  FHG,NAME=BRIGITTE
(7) //G.STEPLIB DD  UNIT=3330,VOL=SER=KAPROS,DSN=STEINPDS,DISP=SHR
(8) //G.FTO1FOO1 DD  UNIT=uuuu,VOL=SER=volvol,
(9) //      SPACE=(3520,(200)),DISP=(OLD,KEEP),
(10) //      DSN=dsndsn,
(11) //      DCB=(RECFM=F,BLKSIZE=3520,DSORG=DA,LRECL=3520)
(12) //G.FTO9FOO1 DD  UNIT=TAPE9,VOL=SER=ididid,
(13) //      DSN=dsname,DISP=(OLD,PASS),
(14) //      LABEL=(n,SL,,IN),
(15) //      DCB=(RECFM=FB,LRECL=80,BLKSIZE=3520,DEN=2)
(16) //G.FTO10FOO1 DD  UNIT=SYSDA,SPACE=(3520,(30,5)),
(17) //      DISP=(NEW,DELETE,DELETE),
(18) //      DCB=(RECFM=VS,BLKSIZE=3520)
(19) //G.FT11FOO1 DD  UNIT=SYSDA,SPACE=(1600,(800),,CONTIG),
(20) //      DCB=(DSORG=DA,RECFM=F,BLKSIZE=1600)
(21) //G.FT12FOO1 DD  UNIT=SYSDA,SPACE=(1600,(800),,CONTIG),
(22) //      DCB=(DSORG=DA,RECFM=F,BLKSIZE=1600)
```

```
(23) //G.SYSIN DD *
(24) &SPACE &END
(25) &INPUT MATP=mmmm,NAMKDK='NP237EN4',
(26) MATKDK=0920237,EB=6.619E+6,
(27) LEDITS=1,LABEL=111,ELINEG=1,
(28) &END
(29) &LIB IFIRST=0,IMOD=1,&END
(30) /*
(31) //
```

Remarks to the cards:

See remarks for example 1.

### Example 3

Mo (Molybdenum) is to be translated. The result is to be inserted in an existing BRIGITTE created KEDAK-library (see 2.2.4). If this library contains Mo-data, the converted data must not be inserted in the library. A maximal output is wished; negative  $\sigma_t$  and  $\sigma_e$  cross section values in the resolved resonance region are to be deleted. The ENDF/B-material contains in MF=2 (resonance parameters) information for 7 different isotopes. The namelist &ISO must be present. The default value of NIX (500) is to be increased on 600.

```
(1) // (job card information),
(2) // REGION=300K,TIME=2
(3) /*MAIN LINES=2
(4) /*SETUP DEVICE=TAPE9,ID=(ididid,NORING,,SL)
(5) /*FORMAT PR,DDNAME=FTO6FOO1,OVFL=ON
(6) // EXEC FHG,NAME=BRIGITTE
(7) //G.STEPLIB DD UNIT=3330,VOL=SER=KAPROS,DSN=STEINPDS,DISP=SHR
```

```
(8) //G.FTO1FOO1 DD UNIT=uuuuu,VOL=SER=volvol,
(9) // SPACE=(3520,(200)),DISP=(OLD,KEEP),
(10) // DSN=dsndsn,
(11) // DCB=(RECFM=F,BLKSIZE=3520,DSORG=DA,LRECL=3520)
(12) //G.FTO9FOO1 DD UNIT=TAPE9,VOL=SER=ididid,
(13) // DSN=dsname,DISP=(OLD,PASS),
(14) // LABEL=(n,SL,,IN),
(15) // DCB=(RECFM=FB,LRECL=80,BLKSIZE=3520,DEN=2)
(16) //G.FT10FOO1 DD UNIT=SYSDA,SPACE=(3520,(30,5)),
(17) // DISP=(NEW,DELETE,DELETE),
(18) // DCB=(RECFM=VS,BLKSIZE=3520)
(19) //G.FT11FOO1 DD UNIT=SYSDA,SPACE=(1600,(400),,CONTIG),
(20) // DCB=(DSORG=DA,RECFM=F,BLKSIZE=1600)
(21) //G.FT12FOO1 DD UNIT=SYSDA,SPACE=(1600,(400),,CONTIG),
(22) // DCB=(DSORG=DA,RECFM=F,BLKSIZE=1600)
(23) //G.SYSIN DD *
(24) &SPACE N1X=600,&END
(25) &INPUT MATP=mmmm,NAMKDK='MO EN4',
(26) MATKDK=0420000,EB=8.6424E+6,NISO=7
(27) LEDITS=1,LABEL=111,ELINEG=1,
(28) &END
(29) &ISO NAMISO='MO 92EN4','MO 94EN4',
(30) 'MO 95EN4','MO 96EN4',
(31) 'MO 97EN4','MO 98EN4',
(32) 'MO100EN4',
(33) MATISO=0420092,0420094,0420095
(34) 0420096,0420097,0420098,
(35) 0420100,
```

```
(36) EBISO=12.692.E+6,9.6722E+6,7.3751E+6,  
(37)          9.1542E+6,6.8161E+6,8.6424E+6,  
(38)          8.301E+6,&END  
(39) &LIB IFIRST=0,IMOD=0,&END  
(40) /*  
(41) //
```

Remarks to the cards:

See example 1.



## 2.3 Programmer's guide

The following sections have mainly been written for programmers. But the section 2.3.1 is very useful for all those users, desiring to get a general survey about the program. It is also possible, that problems, which may arise have with the program, may be solved by reading the corresponding section in this part of the description.

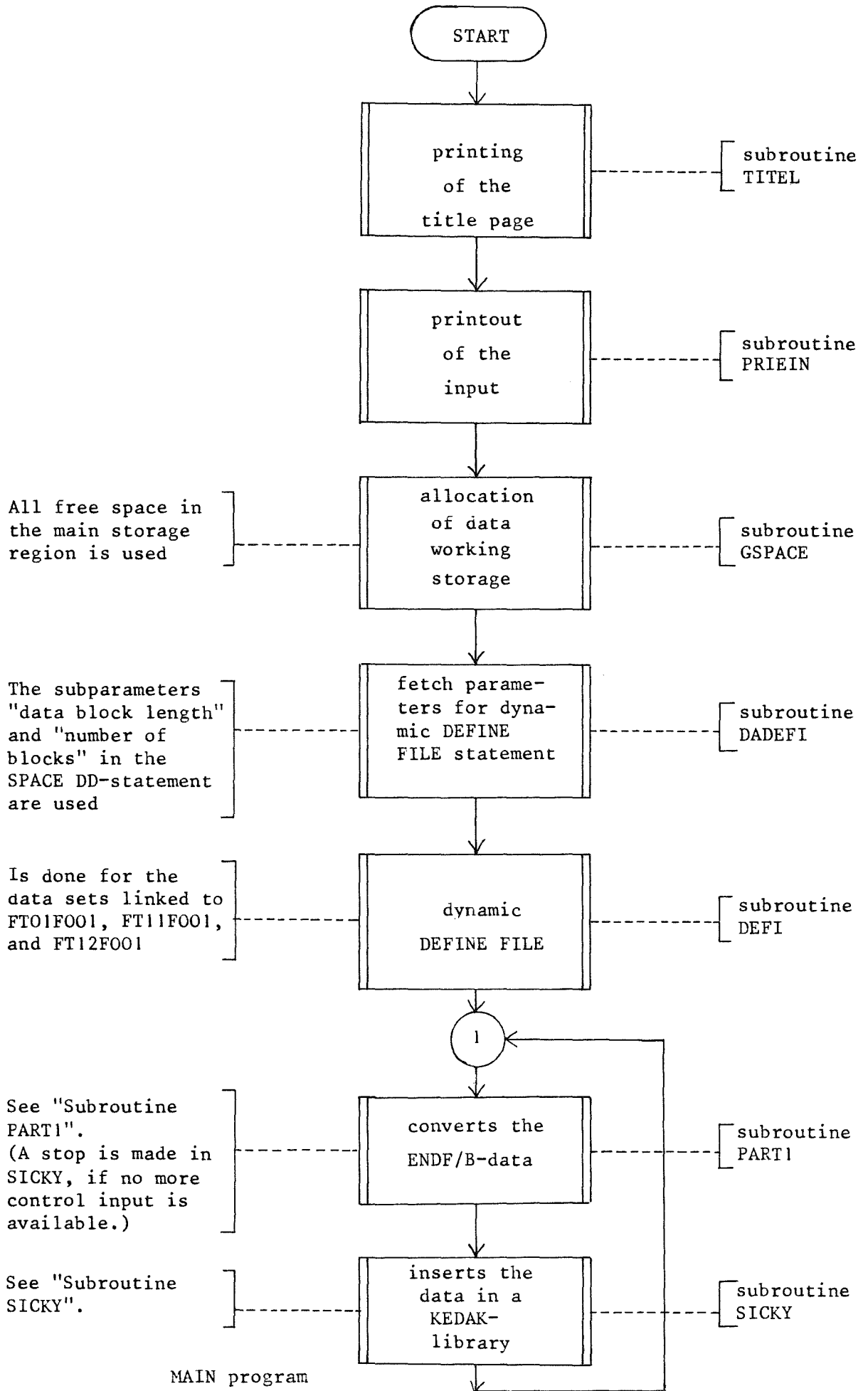
The contents of the following pages can best be seen in the table of contents. The sections containing general aspects of the program (2.3.1 until 2.3.4) are followed by a detailed description of each subroutine resp. function (2.3.5.1 until 2.3.5.99). The purpose of the subroutines and functions respectively can be found in headings preceding the groups of subroutines and functions respectively.

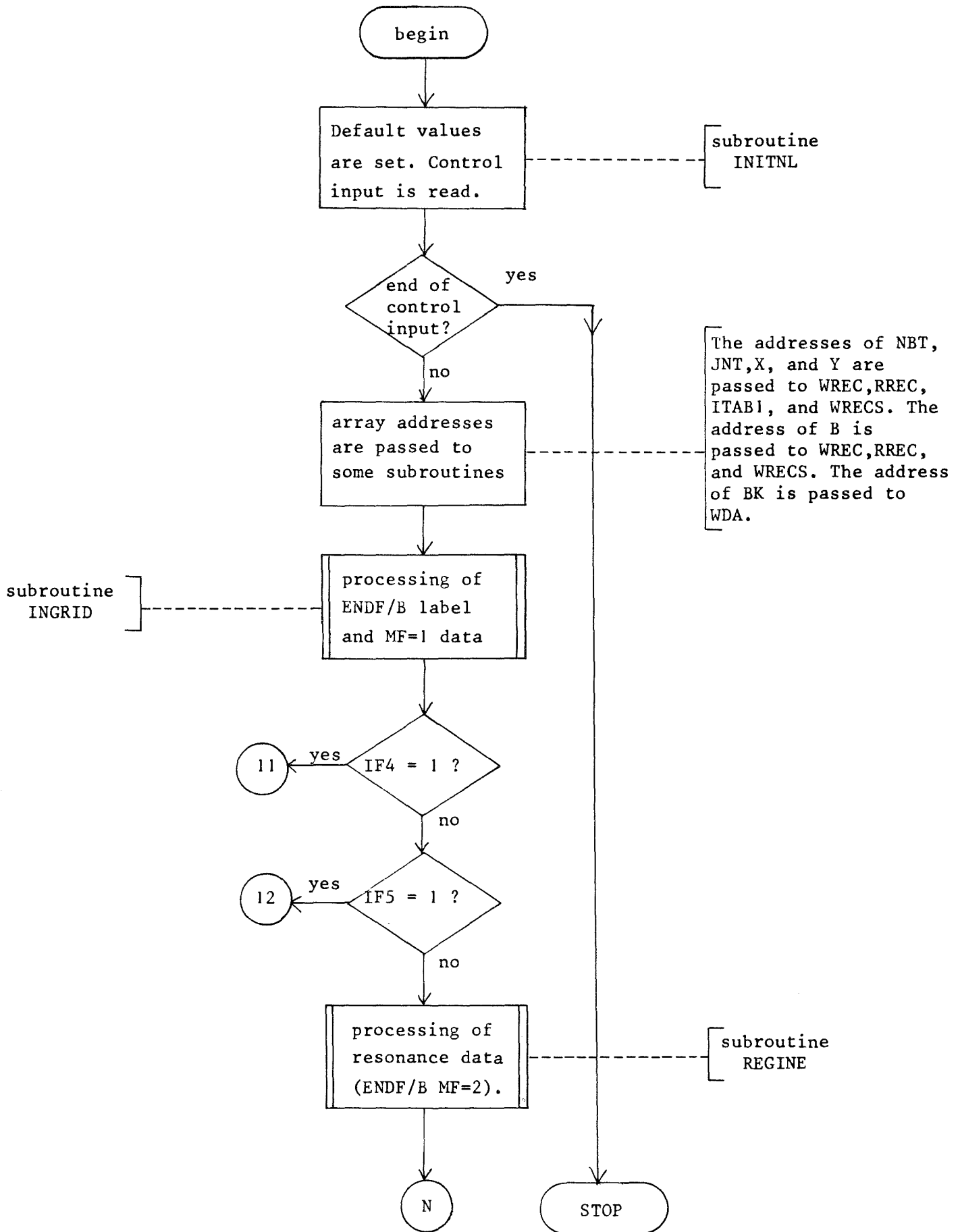
### 2.3.1 General program flow

To get a survey about the program, a block diagram of the program and data flow will be given.

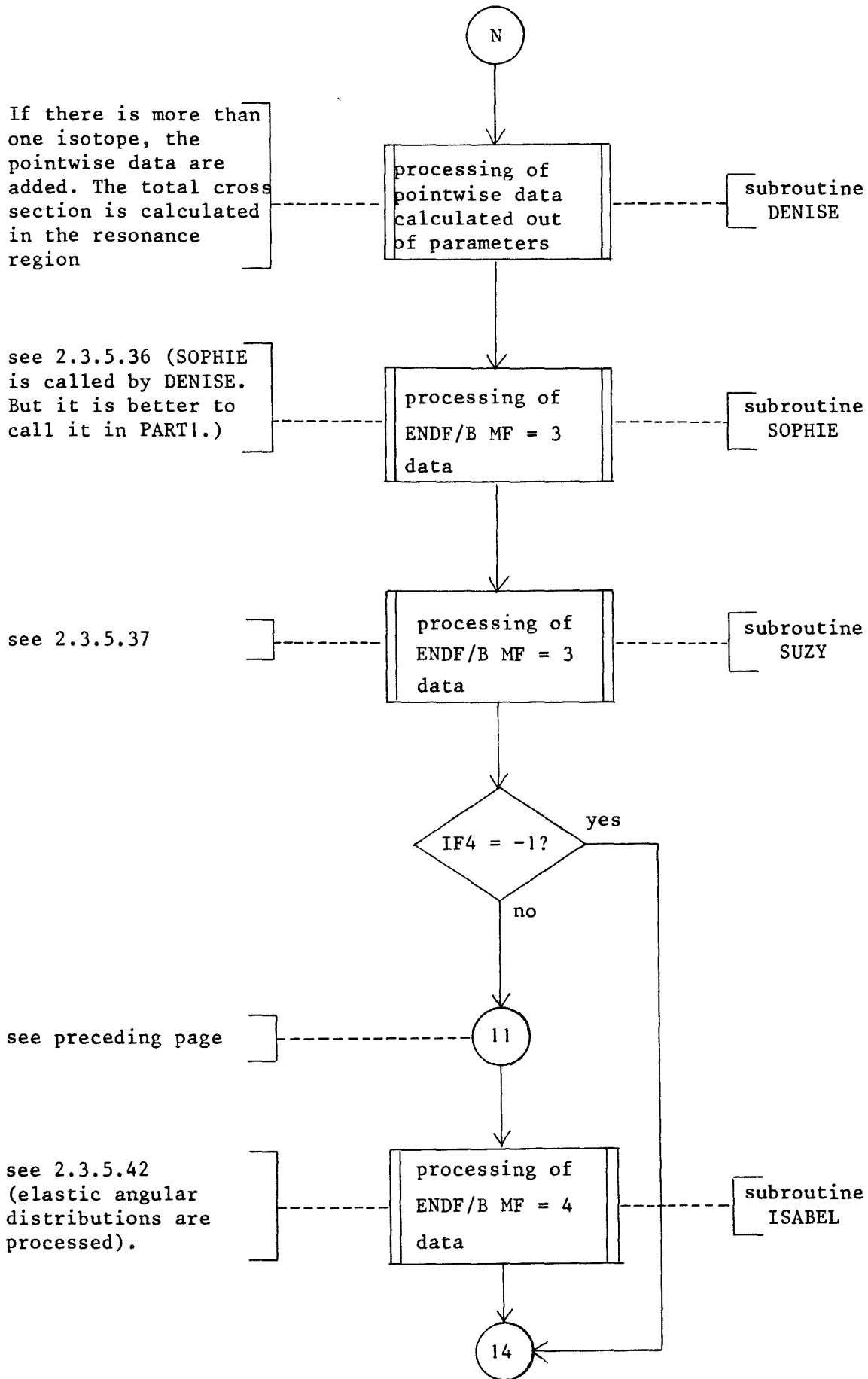
There are eight tables describing in a rough manner the program flow followed by two tables describing the data flow. The symbols used in the tables are those of DIN 66001 (see /12/). The first table "MAIN program" describes the MAIN-program. Two programs called by the MAIN-program will be described in the following tables. These are the subroutine PART1 and the subroutine SICKY. The subroutine PART1 resp. subroutines called by it does all the translation work, whereas the subroutine SICKY resp. subroutines called by it will insert the translated data into the KEDAK-dataset.

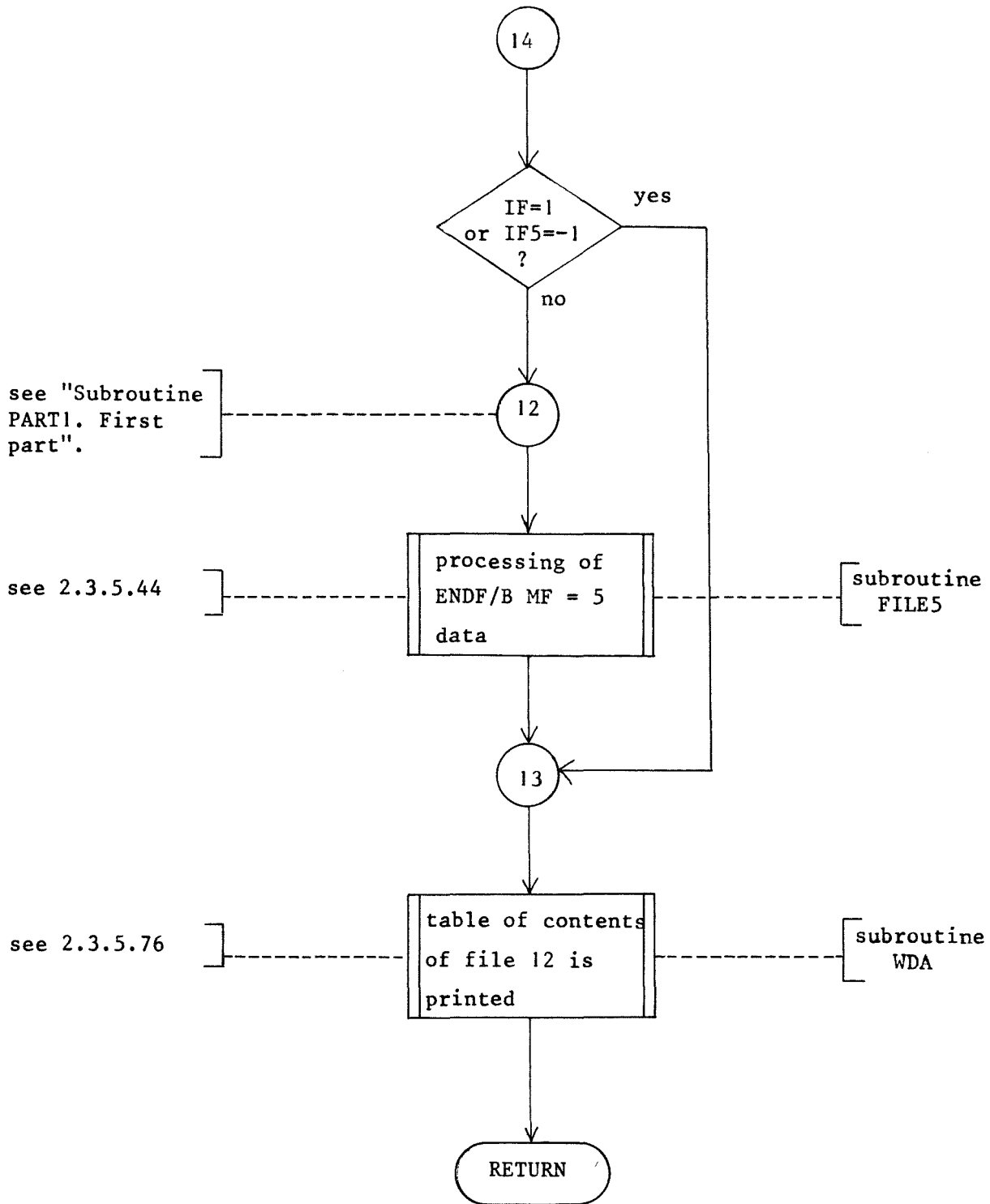
As one can see in the program flow tables, the ENDF/B-data are processed in the sequence of their file (MF)-numbers. This fact should be considered, when looking at the two data flow tables. Remember that the KEDAK-library on file 1 must have been initialized by BRIGITTE itself (see 2.2.4).



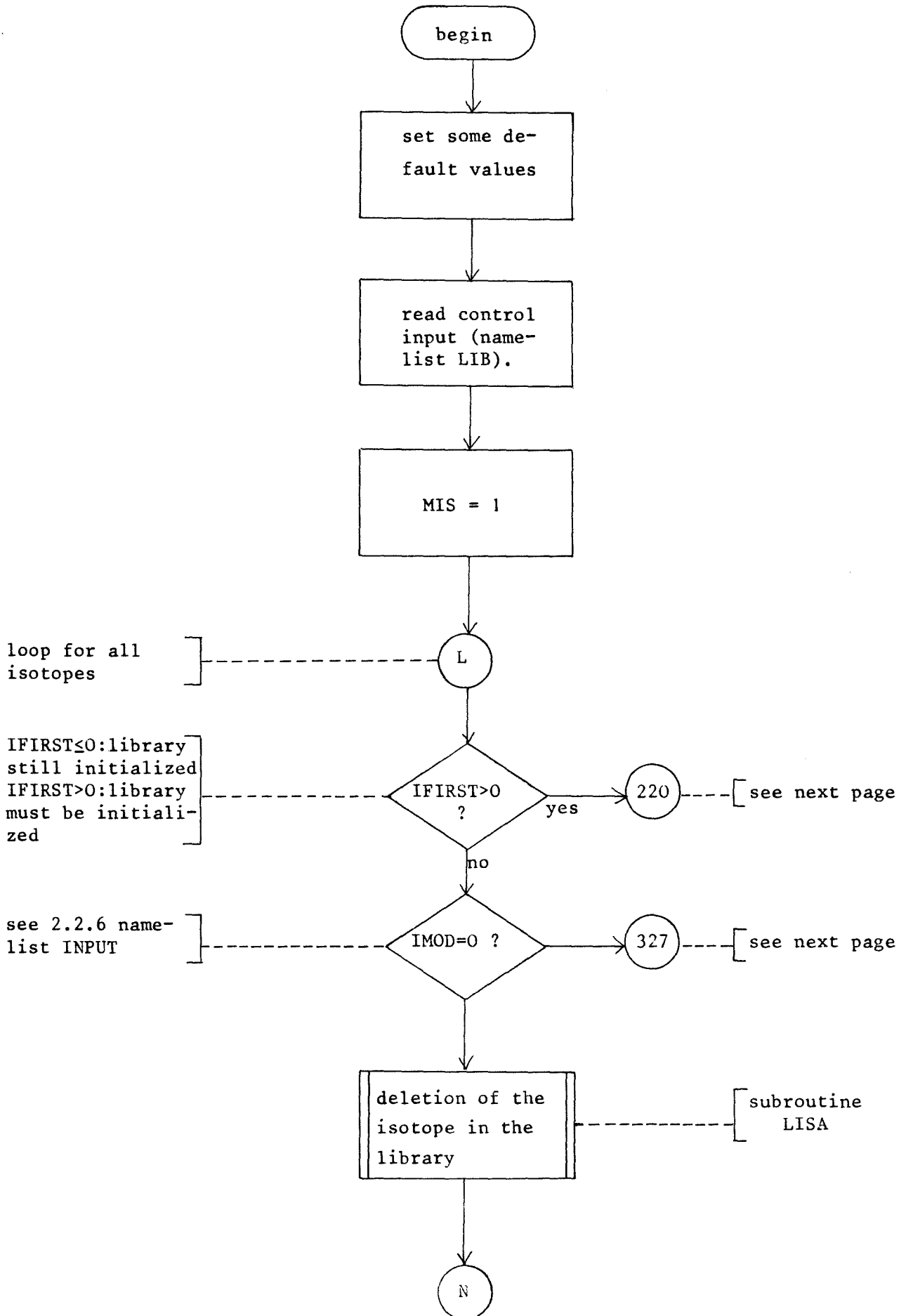


Subroutine PART1. First part (continued on next page).

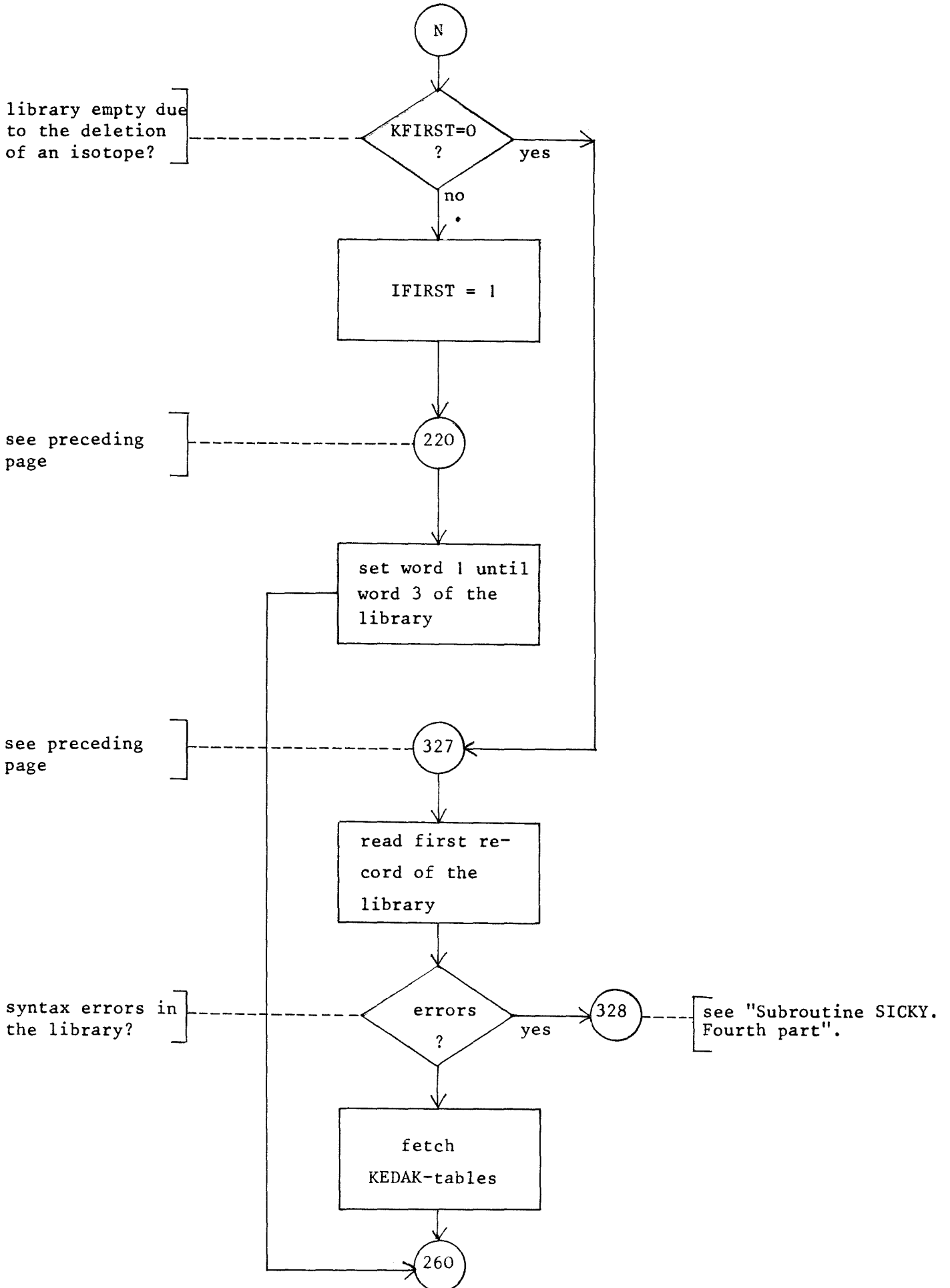


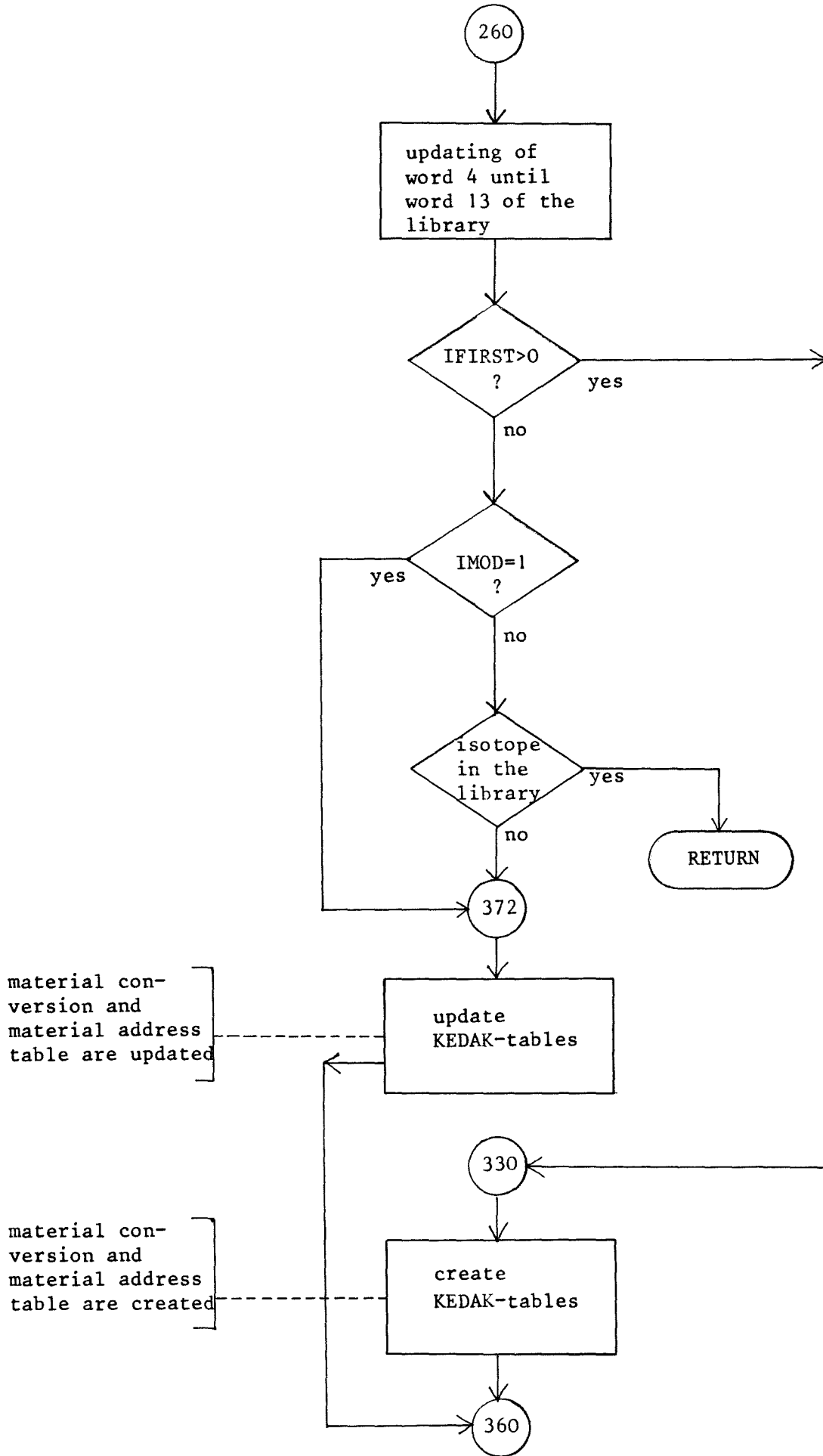


Subroutine PART1. Third part (continuation of preceding page).

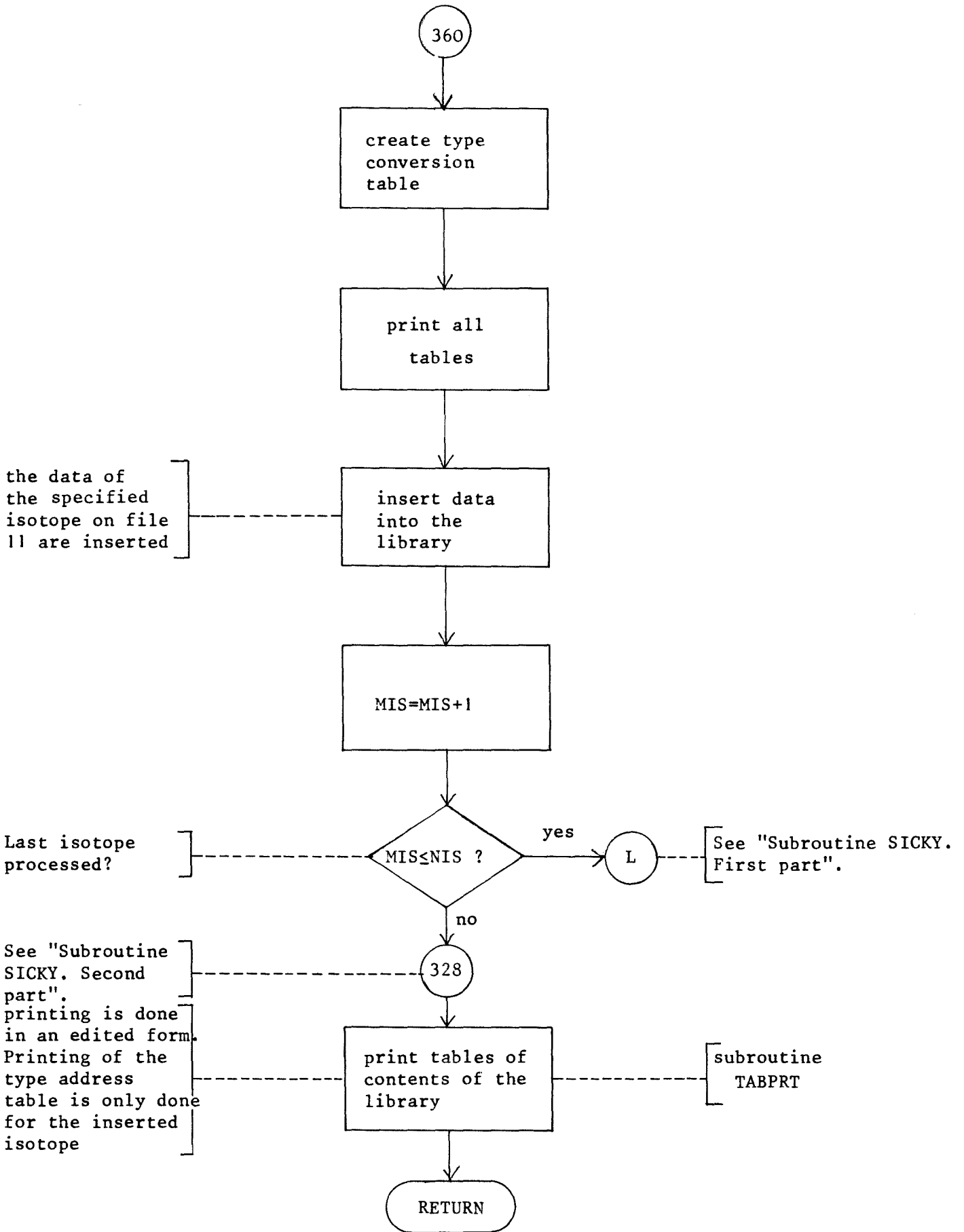


Subroutine SICKY. First part (continued on next page).

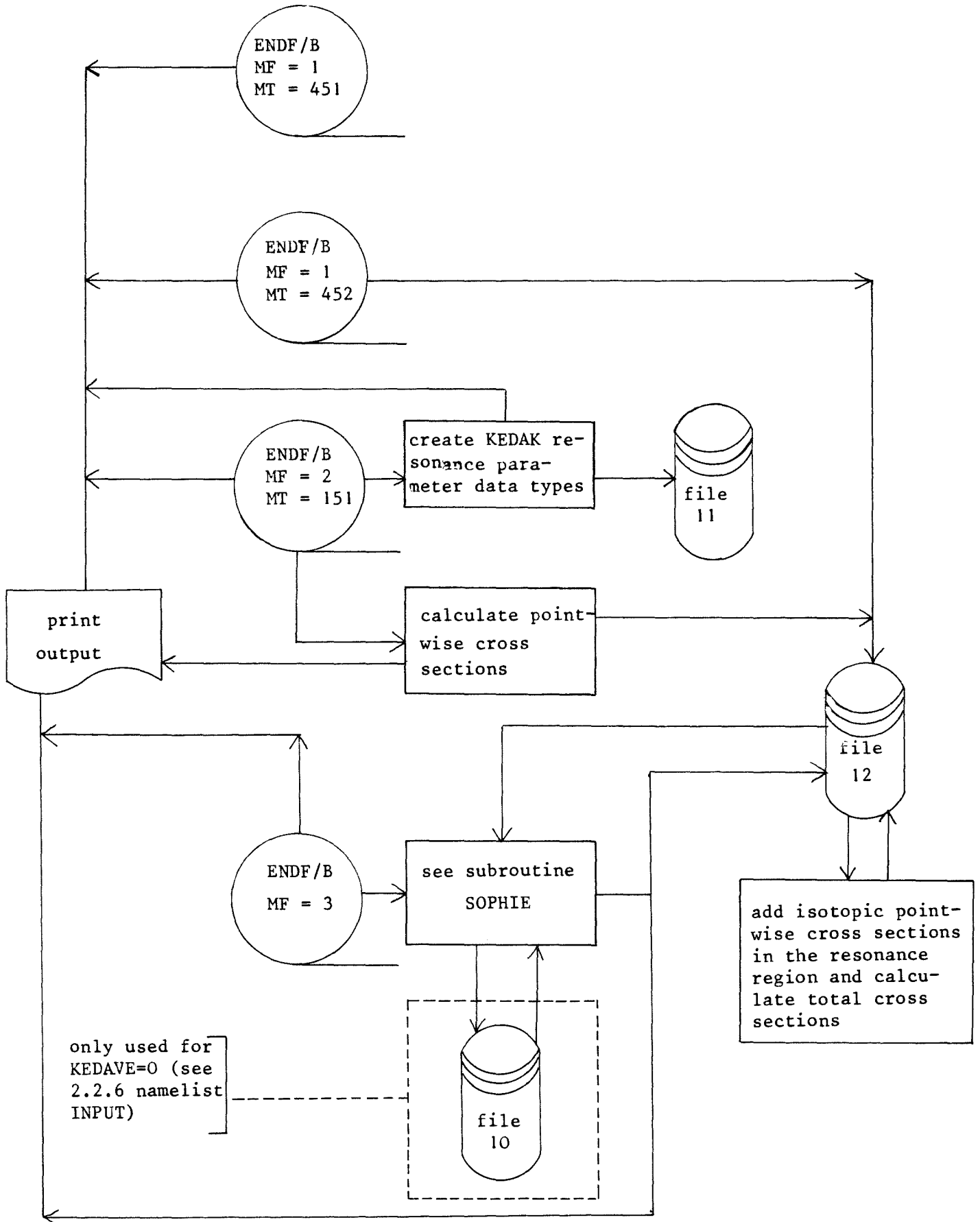




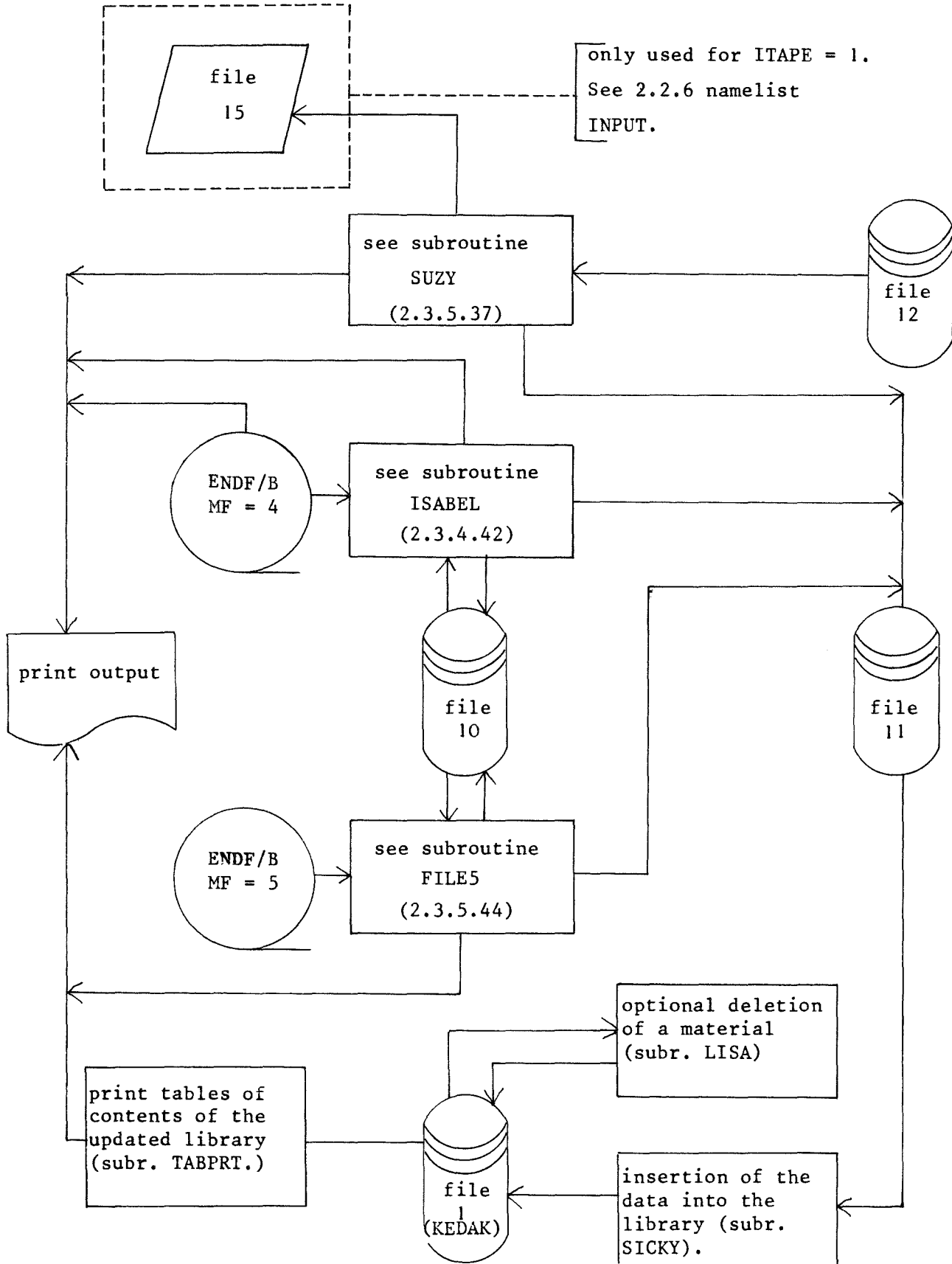




Subroutine SICKY. Fourth part (continuation of the preceding page).



Data flow in BRIGITTE. First part (continued on the next page). Note that file 10 is not obliged to reside on a disk storage.



Data flow in BRIGITTE. Second part (continuation of the preceding page). See note on the preceding page.

### 2.3.2 Overlay structure

BRIGITTE uses a two region overlay structure (see IBM System/370 literature). That means, the second region starts at the end of the largest segment of the first region. This two region structure enables all subroutines of the first region, to call any subroutine of the second region.

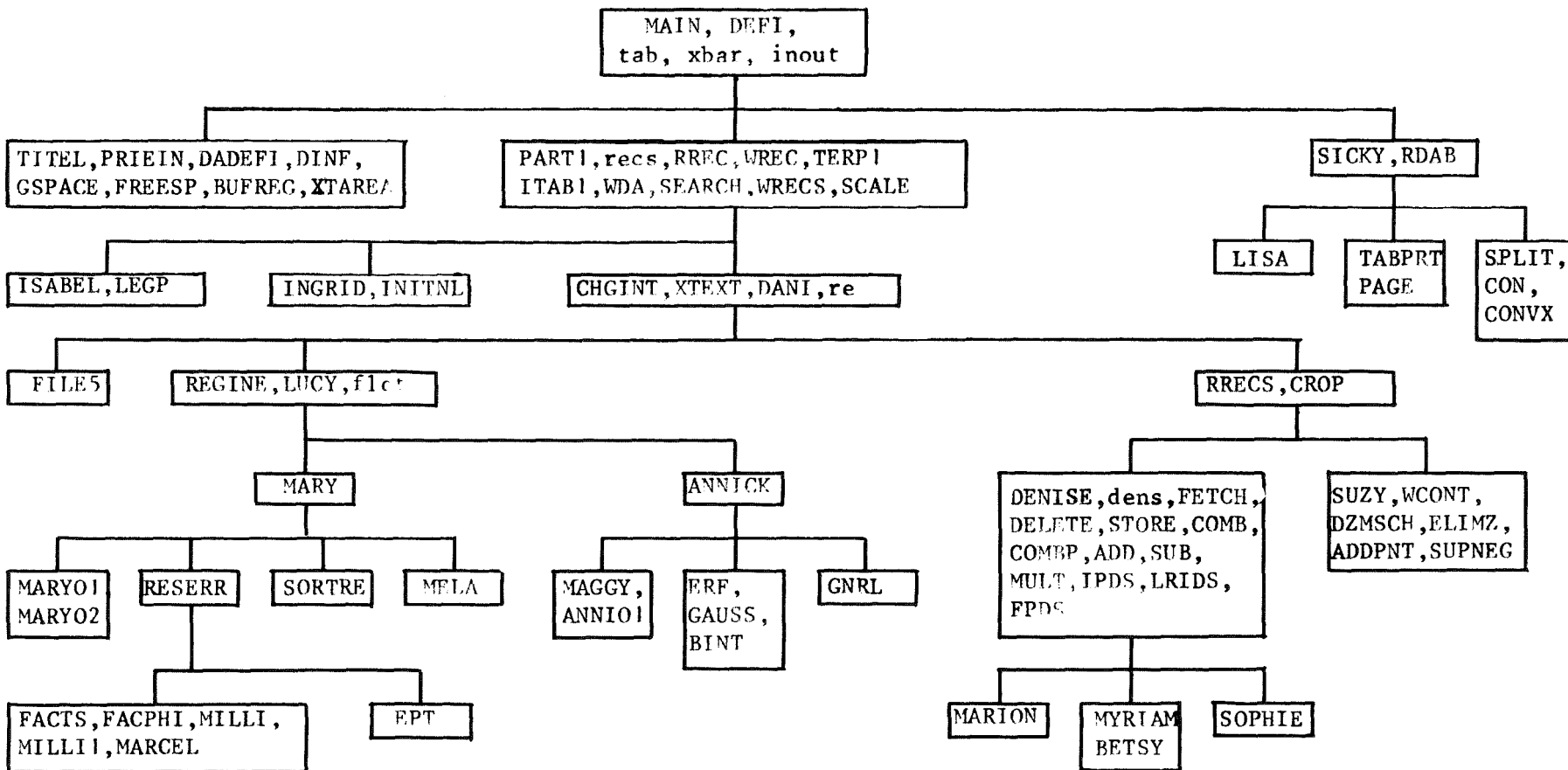
First a list of the used overlay control cards for the IBM linkage editor will be given. Second a graph will illustrate the overlay structure. To differ between subroutines and common blocks in this graph, capital letters have been used for subroutine names and small letters have been used for common block names.

Overlay control cards:

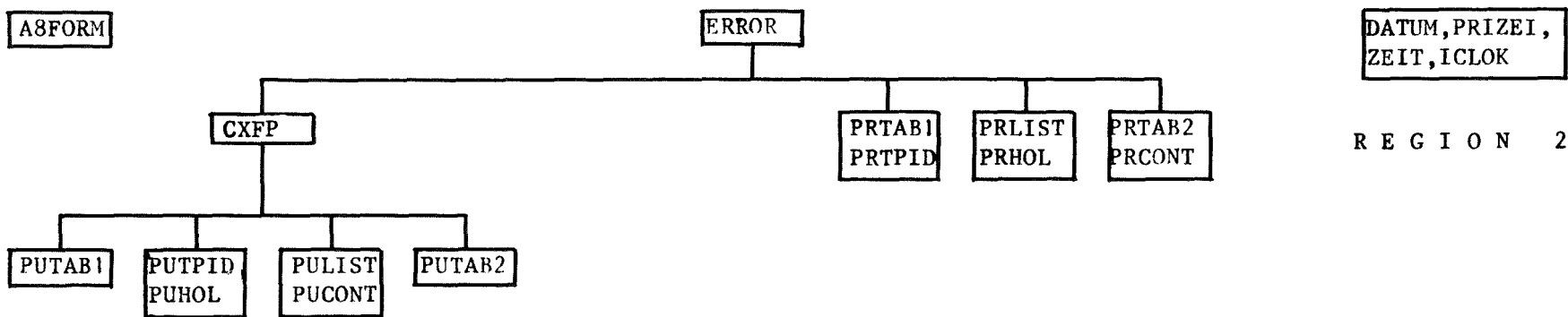
```
OVERLAY  A
INSERT  PRIEIN,DADEFI,DINF,GSPACE,FREESP
INSERT  BUFREG,XTAREA,TITEL
OVERLAY  A
INSERT  SICKY,RDAB
OVERLAY  C
INSERT  SPLIT,CON,CONVX
OVERLAY  C
INSERT  TABPRT,PAGE
OVERLAY  C
INSERT  LISA
OVERLAY  A
INSERT  PART1,RECS,RREC,WREC,TERP1,ITAB1
INSERT  WDA,SEARCH,WRECS,SCALE
OVERLAY  D
INSERT  ISABEL,LEGP
OVERLAY  D
INSERT  INGRID,INITNL
OVERLAY  D
INSERT  CHGINT,XTEXT,DANI,RE
OVERLAY  F
INSERT  FILE5
```

OVERLAY F  
INSERT REGINE,LUCY,FLCT  
OVERLAY G  
INSERT MARY  
OVERLAY G1  
INSERT MARYO1,MARYO2  
OVERLAY G1  
INSERT SORTRE  
OVERLAY G1  
INSERT RESERR  
OVERLAY G2  
INSERT EPT  
OVERLAY G2  
INSERT FACTS,FACPHI,MILLI,MILLI1,MARCEL  
OVERLAY G1  
INSERT MELA  
OVERLAY G  
INSERT ANNICK  
OVERLAY H  
INSERT MAGGY,ANNIO1  
OVERLAY H  
INSERT ERF,GAUSS,BINT  
OVERLAY H  
INSERT GNRL  
OVERLAY F  
INSERT RRECS,CROP  
OVERLAY I  
INSERT SUZY,WCONT,DZMSCH,ELIMZ,ADDPNT,SUPNEG  
OVERLAY I  
INSERT DENISE,DENS,FETCH,DELETE,STORE,COMB  
INSERT COMBP,ADD,SUB,MULT,IPDS,LRIDS,FPDS  
OVERLAY K  
INSERT MARION  
OVERLAY K  
INSERT MYRIAM,BETSY  
OVERLAY K  
INSERT SOPHIE  
OVERLAY R(REGION)  
INSERT ERROR

OVERLAY R1  
INSERT CXFP  
OVERLAY R2  
INSERT PUTAB1  
OVERLAY R2  
INSERT PUTPID,PUHOL  
OVERLAY R2  
INSERT PULIST,PUCONT  
OVERLAY R2  
INSERT PUTAB2  
OVERLAY R1  
INSERT PRTAB1,PRTPID  
OVERLAY R1  
INSERT PRLIST,PRHOL  
OVERLAY R1  
INSERT PRTAB2,PRCONT  
OVERLAY R  
INSERT A8FORM  
OVERLAY R  
INSERT DATUM,PRIZEI,ZEIT,ICLOK



Overlay structure of BRIGITTE



### 2.3.3 Organisation of data storage

Nearly all arrays used in the program are variably dimensioned. Because there are no aids in FORTRAN to make a region dependant dimensioning of arrays, ASSEMBLER subroutines must be used.

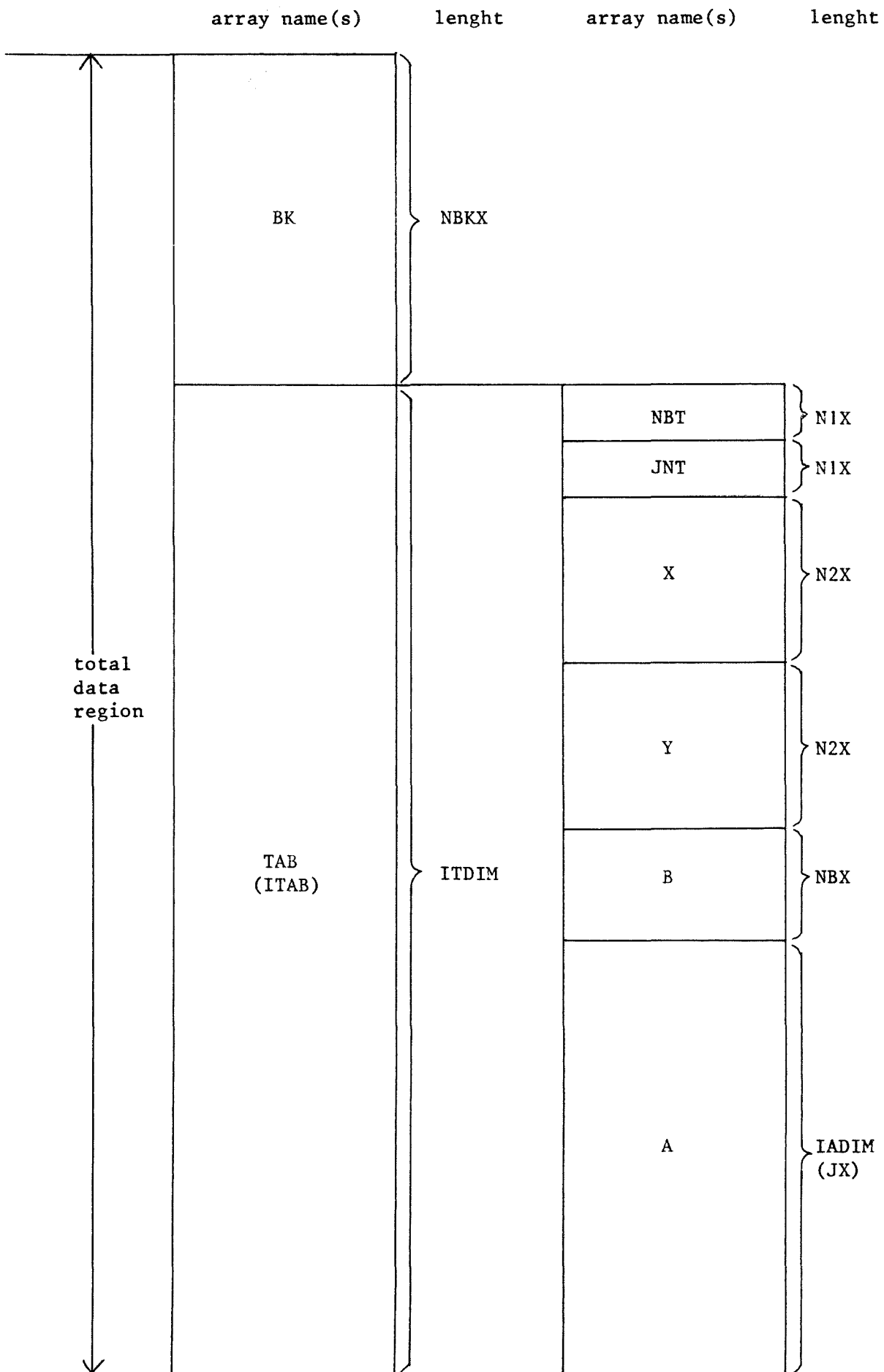
The FORTRAN-subroutine GSPACE (see 2.3.5.5) called by the main program, first evaluates the free region by using the ASSEMBLER subroutine FREESP (see 2.3.5.95). The program only continues, if this region is greater or equal 10 k-byte. Now the number of bytes needed for buffers (see subroutine BUFREG 2.3.5.6) and 6 k-byte for system use are subtracted from the total free region, which is returned by the subroutine FREESP. The remaining region is used for data arrays. The subdivision of this region in the individual arrays is shown in the following graph. To get access to this region the ASSEMBLER subroutine XTAREA (see 2.3.5.93) is used. This routine returns the absolute address of the beginning of the data storage, after having assigned ownership of this storage to the active job, by using the ASSEMBLER makro GETMAIN. Because XTAREA also returns the absolute address of the first word of a given array, which address has been defined in the calling program and passed to the subroutine, the beginning of the data region can be expressed as an index value in this array. This is done by calculating a displacement IVER, so that  $F(1+IVER)$  is the first word of the data region, if F is the given array. This displacement is returned to the calling program, which passes to the working subroutines the address  $F(1+IVER)$  for the array BK, resp.  $F(1+IVER+IVAL)$  with corresponding IVAL for the other arrays.

The given values for the array lenght - read or default values (see 2.2.6 namelist SPACE) - are checked by GSPACE. JX is set to IADIM, NBKX is made even, and ITDIM is set to maximum  $(ITDIM, 2*N1X+2*N2X+NBX+IADIM)$ .

Now NBX is set to  $ITDIM-IADIM-2*(N1X+N2X)$ , and afterwards ITDIM is rounded to a multiple of 880. By using the default values listed in 2.2.6, we get the following dimensions:

NBKX	=	12 000 words		
ITDIM	=	33 440 words	N1X	= 500 words
			N2X	= 5000 words
			NBX	= 3000 words
			IADIM	= JX = 20000 words





Overlay of the data region.

Summing up the length of all arrays, a total of 46 000 words is the result.

If the extent of the data region is not large enough to contain these 46 000 words ( $\approx 180$  k-byte), all dimensions will be shortened by multiplication with a factor  $KW/ISUM$ .  $KW$  means the extent of the existing data region,  $ISUM$  the result of summing up the lengths of all arrays, both measured in 4-byte words. The new dimensions are once more checked and eventually changed (e.g.  $NBKX$  is made even). After the calculation of the beginning of all arrays relative to  $BK(1)$ , `RETURN` is made.

If  $1.1*ISUM$  is less or equal  $KW$  all dimensions will be increased by multiplication with the factor  $KW/ISUM$ . Afterwards the same is done as in the case of shortening the dimensions. All quantities calculated by `GSPACE` are listed in the program printout.

In the table "overlay of the data region", one can see that parts of the data region are used twice. This is naturally not done simultaneously in one subroutine but in different ones. Also some arrays are subdivided in some subroutines.

To get a survey about the usage of the arrays in the program, a table is made for each array. This tables contain in the heading the names and the length of the arrays. If the name of an array in a subroutine is the same, as in the heading, only the names of the subroutines, in which this array is used, are listed. Otherwise the names of the subroutines together with the names of the arrays, or the names of the subarrays and the maximal dimensions, are tabulated. If a subroutine may use - dependant on the argument list in the call statement - different arrays, it is not contained in the following tabulations.

Some of the arrays are only used for defined data.  $NBT, JNT, X,$  and  $Y$  contain - with a few exceptions - only `TAB1`-records (see /3/ chapter R-2.3.3).  $NBT$  and  $JNT$  contain only interpolation tables. The array  $B$  is used to store `LIST`-records (see /3/ chapter R-2.3.2) read from `ENDF/B`-tapes. In some cases it is also used as an auxiliary array.

name of the array: BK		length: NBKX
subroutine name	name of the array in the subroutine, or name and dimensions of subarrays	
PART1	BK	
REGINE	BK	
MARY	BK	
MARYO1	BK	
LUCY	BK	
MILLI	NBTN(NIX),JNTN(NIX),NBTG(NIX),JNTG(NIX),NBTF(NIX), JNTF(NIX)	
MILLI1	one of the arrays NBTN,JNTN, or NBTG,JNTG, or NBTF, JNTF	
MELA	IBK	
ANNICK	BK	
ANNIO1	BK	
DENISE	BK	
SOPHIE	XIN(NBKX/2),YIN(NBKX/2)	
SUZY	BK,A,TAB,ITAB	
ISABEL	ITAB,TAB,CLEG	
FILE5	BK,IBK	
WDA	BK	
RDAB	BK	
SICKY	BK	

name of the array: NBT	length: N1X
used in the subroutines (with the name NBT): PART1,REGINE,MARY,MELA,ANNICK,ANNIO1,DENISE,SOPHIE,SUZY,ELIMZ,ADDPNT, DZMSCH,SUPNEG,ISABEL,FILE5,RREC,WREC,PRTAB1,PRTAB2,PUTAB1,PUTAB2,COMB, COMBP,ITAB1,CHGINT,CROP,WRECS,RRECS,STORE,FETCH	
name of the array: JNT	length: N1X
used in the subroutines (with the name JNT): PART1,REGINE,MARY,MELY,ANNICK,ANNIO1,DENISE,SOPHIE,SUZY,ELIMZ,ADDPNT, DZMSCH,SUPNEG,ISABEL,FILE5,RREC,WREC,PRTAB1,PRTAB2,PUTAB1,PUTAB2,COMB, COMBP,ITAB1,CHGINT,CROP,WRECS,RRECS,STORE,FETCH	
name of the array: X	length: N2X
used in the subroutines (with the name X): PART1,REGINE,MARY <sup>1)</sup> ,EPT <sup>1)</sup> ,MARYO2 <sup>1)</sup> ,MILLI <sup>1)</sup> ,ANNICK <sup>2)</sup> ,ANNIO1 <sup>2)</sup> ,MAGGY, DENISE,SOPHIE,SUZY,ELIMZ,ADDPNT,DZMSCH,SUPNEG,ISABEL,FILE5,RREC,WREC, PRTAB1,PUTAB1,COMB,COMBP,ITAB1,CHGINT,CROP,WRECS,RRECS,STORE,FETCH	
<sup>1)</sup> name of the array: ENS <sup>2)</sup> name of the array: BK2	
name of the array: Y	length: N2X
used in the subroutines (with the name Y): PART1,REGINE,MARY <sup>1)</sup> ,MELA <sup>1)</sup> ,ANNICK,ANNIO1,MAGGY,DENISE,MYRIAM,BETSY, SOPHIE,SUZY,ELIMZ,ADDPNT,DZMSCH,SUPNEG,ISABEL,FILE5,RREC,WREC,PRTAB1, PUTAB1,COMB,COMBP,ITAB1,CHGINT,CROP,WRECS,RRECS,STORE,FETCH	
<sup>1)</sup> name of the array: SIGS	
name of the array: B	length: NBX
used in the subroutines (with the name B): PART1,REGINE,MARY,ANNICK,DENISE,SOPHIE <sup>1)</sup> ,SUZY,ISABEL,FILE5 <sup>1)</sup> ,RREC,WREC, PRLIST,PRHOL,PRTPID,PULIST,PUHOL,PUTPID,WRECS,RRECS,STORE,FETCH	
<sup>1)</sup> name of the array: B and TEMP	

name of the array: A		length: IADIM
subroutine name	name of the array in the subroutine, or name and dimensions of subarrays	
PART1	A	
REGINE	AH	
MARY	ER(NDIM),AJ(NDIM),GT(NDIM),GN(NDIM),GG(NDIM),GF(NDIM), AL(NDIM),G(NDIM),SER(NDIM),PER(NDIM),AH(JXRED),NDIM = JXRES = IADIM/200; JXRED = $\frac{19}{20} * IADIM$	
SORTE	ER(NDIM),AJ(NDIM),GT(NDIM),GN(NDIM),GG(NDIM),GF(NDIM), AL(NDIM),G(NDIM),SER(NDIM),PER(NDIM) See MARY	
MARYO1	ER(NDIM),AJ(NDIM),GT(NDIM),GN(NDIM),GG(NDIM),GF(NDIM), AL(NDIM),G(NDIM) See MARY	
EPT	ER(NDIM),GTR(NDIM) See MARY(GTR is GT)	
MARCEL	ER(NDIM),GN(NDIM),GG(NDIM),GF(NDIM),AL(NDIM),G(NDIM), SER(NDIM),PER(NDIM) See MARY	
MILLI	AH	
MELA	AH	
ANNICK	SIGS(NPEMAX),E(NPEMAX),SIGG(NPEMAX),SIGF(NPEMAX), SIGT(NPEMAX),ESC(NPEMAX),XT(NPEMAX),YT(NPEMAX), TEMP(NPEMAX),D(6,NPEMAX),GX(6,NPEMAX),GNO(6,NPEMAX), GG(6,NPEMAX),GF(6,NPEMAX),AJ(6,NPEMAX),NPEMAX=IADIM/45	
ANNIO1	SIGS(NPEMAX),E(NPEMAX),SIGG(NPEMAX),SIGF(NPEMAX), SIGT(NPEMAX),ESC(NPEMAX) See ANNICK	
MAGGY	XD(NPEMAX),YD(NPEMAX) See ANNIO1. XD is E, YD is SIGS,SIGG, or SIGF	
DENISE	A	
ISABEL	A, ITAB2, TAB2	

name of the array: A (continuation)                      lenght: IADIM	
subroutine name	name of the array in the subroutine, or name and dimensions of subarrays
STORE	A,LA
FETCH	A,LA
DELETE	A
FPDS	A,LA
IPDS	A,LA

- name of the array: TAB                      lenght: ITDIM	
used in the subroutine (with the name TAB and ITAB): SICKY	

#### 2.3.4 COMMONS used in the program

There are seven named COMMONS used by the program. The COMMON names are listed in the following in alphabetic order:

DENS  
FLCT  
INOUT  
  
RE  
RECS  
TAB  
  
XBAR

For each of this COMMON blocks a separate table (which can be distributed on several pages) is given. In this table you will find for each element of the COMMON the (mainly used) name of this element, the address (measured in 4\*byte words, starting with 1), the dimension, the type, the number of the program(s), where the contents of this element are used, and the meaning of the element.

The abbreviations used,are:

in the column headed "dimension":

scalar            for scalar quantities  
(n)                for an array with n-elements  
(n,m)             for an array with n\*m elements

in the column headed "type":

I4                for INTEGER\*4 quantities  
R4                for REAL\*4 quantities  
R8                for REAL\*8 quantities  
L4                for LOGICAL\*4 quantities

in the column headed "used in program number":

n                 fetched in program 2.3.5.n  
n                 stored (and fetched) in program 2.3.5.n

- (n) in program 2.3.5.n the address of the quantity is passed to a subprogram. In this subprogram, or programs called by this subprogram, the value of the quantity is fetched.
- (n) in program 2.3.5.n the address of the quantity is passed to a subprogram. In this subprogram, or programs called by this subprogram, the value of the quantity is stored (and fetched).

Only one of these four cases is listed.

If n and (n and/or (n) and/or (n)) happens, n is listed.

If (n) and (n and/or (n)) happens, (n) is listed.

If n and (n) happens, n is listed.

E.g. 15 and (15) and 15 will force the list entry 15.



Name of the COMMON block: DENS					
word number	name	dimension	type	used in programs number	meaning
1	JMT	(100)	I4	<u>80</u> , 81, <u>82</u> , 85.	see 2.3.5.80 "problem solved"
101	JAT	(100)	I4	<u>80</u> , 81, <u>82</u> , 85.	" "
201	JTT	(100)	I4	<u>80</u> , 81, <u>82</u> .	" "
301	JLT	(100)	I4	<u>80</u> , <u>82</u> .	" "
401	JNS	scalar	I4	<u>80</u> , <u>82</u> .	" "
402	MNS	scalar	I4	<u>80</u> , 81, <u>82</u> , 85.	" "
403	MX	scalar	I4	<u>32</u> , 80, 82.	length of the arrays JMT, JAT, JTT, and JLT (see 2.3.5.80).
Name of the COMMON block: FLCT					
1	ZYN	scalar	R4	<u>25</u> , 27, 29.	number of degrees of freedom in the neutron width distribution (see 2.3.5.29)
2	ZYF	scalar	R4	<u>25</u> , 29.	number of degrees of freedom in the fission width distribution (see 2.3.5.29).
3	U	scalar	R4	<u>25</u> , 29.	used in the calculation of the Dresner factors in BINT (2.3.5.29).
4	V	scalar	R4	<u>25</u> , 29	used in the calculation of the Dresner factors in BINT (2.3.5.29).

Name of the COMMON block: FLCT (continuation)					
word number	name	dimension	type	used in programs number	meaning
5	RHON	scalar	R4	<u>25</u> , 27, 29.	used in the calculation of the Dresner factors in BINT (2.3.5.29)
6	RHOF	scalar	R4	<u>25</u> , 29.	used in the calculation of the Dresner factors in BINT (2.3.5.29).
7	NDIM6	scalar	I4	<u>11</u> , 25.	used for dimensioning. Is set to 6.
8	NPEMAX	scalar	I4	<u>11</u> , 25.	maximal number of energy values in the unresolved resonance part.
Name of the COMMON block: INOUT					
1	NO	scalar	I4	2, <u>4</u> , 5, 6, <u>9</u> , 10, 11, 12, 15, 18, 22, <u>24</u> , 25, <u>30</u> , 32, 35, 36, 37, 39, 41, 42, 44, 45, 46, 76, 78.	print output unit number
2	NI	scalar	I4	<u>4</u> , 5, <u>9</u> .	control input unit number
3	NT	scalar	I4	8, <u>9</u> , 10, 11, (12), 25, 36, 42, 44	tape unit containing ENDF/B-tape
4	MODE	scalar	I4	<u>9</u> , 10, 11, (12), (25), (36), (37), (42), (44),	arrangement of the ENDF/B-data (see /3/ chapter R-2.2).
5	LEDIT1	scalar	I4	<u>9</u> , 10	control of printing of - descriptive data and dictionary (MF = 1, MT = 451)

Name of the COMMON block: INOUT (continuation)

word number	name	dimension	type	used in programs number	meaning
6	LEDIT2	scalar	I4	<u>9</u> , 12, 15, 25, 30	control of printing of -resolved resonance parameters (in ENDF representation) -resolved resonance data in KEDAK representation -unresolved resonance parameters (in ENDF representation) -KEDAK datatypes ST and STGF
7	LEDITS	scalar	I4	<u>9</u> , 24, 30, 32, 36.	Control of printing of -pointwise cross sections calculated from resolved resonance parameters -pointwise cross sections calculated from unresolved resonance parameters -pointwise total cross sections calculated from resolved resonance parameter -pointwise total cross sections calculated from statistical resonance parameter -ENDF/B (pointwise) reaction types of MF=3
8	KEDAVE	scalar	I4	<u>9</u> , 10, 36, 37	type of KEDAK-conventions used (see 2.2.6 namelist INPUT)
9	ABIS	(20)	R4	<u>11</u> , 34	abundance (weight fraction) of the isotopes
29	NOTTRL	(25)	R8	<u>9</u> , 36	names of datatypes not to be translated (see 2.2.6 namelist INPUT).
79	BIB12	scalar	I4	<u>9</u> , 78, 79	size of table of contents of dataset on FT12FO01 (see 2.2.6 namelist INPUT).

Name of the COMMON block: INOUT (continuation)

word number	name	dimension	type	used in programs number	meaning
80	SALINE	scalar	I4	<u>9</u> , 36, 37	control of searching in MF=3 greater than MT=51 "starting" numbers (see 2.2.6 namelist INPUT).
81	NAMKDK	scalar	R8	<u>9</u> , 10, 11, <u>86</u>	alphamerical KEDAK material name (see 2.2.6 namelist INPUT).
83	MATKDK	scalar	I4	<u>9</u> , 10, 11, 12, 30, (37), 42, (44), <u>86</u>	numerical KEDAK material name (see 2.2.6 namelist INPUT).
84	NTYP	scalar	I4	<u>77</u> , <u>86</u>	number of datatypes for a specified material
85	NAMISO	(10)	R8	<u>9</u> , 12, 15, 25, 30, 86	alphamerical KEDAK isotope names (see 2.2.6 namelist ISO).
105	MATISO	(10)	I4	<u>9</u> , 11, 15, 30, 86	numerical KEDAK isotope names (see 2.2.6 namelist ISO).
115	EBISO	(10)	R4	<u>9</u> , 11	binding energies of the last neutron of the isotopes (see 2.2.6 namelist ISO).
125	ICOMB	scalar	I4	<u>9</u> , 67	automatic choice of interpolation rule or linear-linear interpolation in the subroutine COMBP (see 2.2.6 namelist INPUT)
126	NBX	scalar	I4	<u>4</u> , <u>5</u> , 49, 79, 81	maximal dimension of array B (see 2.3.3)
127	NIX	scalar	I4	<u>4</u> , <u>5</u> , 12, 23, 24, (25), (36), (37), ( <u>44</u> ), 49, 67, 75, 79, 81	maximal dimension of the arrays NBT and JNT (see 2.3.3)

## Name of the COMMON block: INOUT (continuation)

word number	name	dimension	type	used in programs number	meaning
128	N2X	scalar	I4	<u>4</u> , <u>5</u> , (12), 22, 24, 25, (36), (37), 39, (44), 49, 67, 75, 79, 81	maximal dimension of the arrays X and Y (see 2.3.3)
129	N	scalar	I4	<u>11</u> , <u>15</u> , <u>25</u> , <u>30</u> , <u>37</u> , <u>42</u> , <u>44</u> , 76, <u>77</u> , <u>86</u>	number of data to be written in subroutine WDA (see 2.3.5.76) or read in subroutine RDAB (see 2.3.5.77)
130	NBKX	scalar	I4	<u>4</u> , <u>5</u> , 11, 12, 15, 25, 30, 32, 37, <u>42</u> , <u>44</u> , 76	maximal dimension of the array BK (see 2.3.3)
131	JX	scalar	I4	<u>4</u> , <u>5</u> , 11, 80, 82	maximal dimension of the array A (see 2.3.3) JX = IADIM
132	IADIM	scalar	I4	<u>4</u> , <u>5</u> , 42	maximal dimension of the array A (see 2.3.3)
133	ITDIM	scalar	I4	<u>4</u> , <u>5</u> , 86	maximal dimension of the arrays TAB resp. ITAB (see 2.3.3)
134	IVER	scalar	I4	1, <u>5</u>	address of begin of the data region relative to F(1) (measured in 4-byte words) (see 2.3.3) (begin of array BK)
135	IVGENR	scalar	I4	1, <u>5</u>	address of the begin of the arrays TAB resp. ITAB and NBT relative to BK(1) (measured in 4-byte words) (see 2.3.3)
136	IVJNT	scalar	I4	1, <u>5</u>	address of the begin of the array JNT relative to BK(1) (measured in 4-byte words) (see 2.3.3)
137	IVX	scalar	I4	1, <u>5</u>	address of the begin of the array X relative to BK(1) (measured in 4-byte words) (see 2.3.3)

Name of the COMMON block: INOUT (continuation)

word number	name	dimension	type	used in programs number	meaning
138	IVY	scalar	I4	1, <u>5</u>	address of the begin of the array Y relative to BK(1) (measured in 4*byte words) (see 2.3.3)
139	IVB	scalar	I4	1, <u>5</u>	address of the begin of the array B relative to BK(1) (measured in 4*byte words) (see 2.3.3)
140	IVDENS	scalar	I4	1, <u>5</u>	address of the begin of the array A relative to BK(1) (measured in 4*byte words) (see 2.3.3)
141	EPSRES	scalar	R4	<u>9</u> , 22, (25), 32, 37, (44)	accuracy used in several cases (see 2.2.6 namelist INPUT)
142	SCREPS	scalar	R4	<u>9</u> , 37, (66)	accuracy used in deleting points, which can be predicted by linear interpolation (see 2.2.6 namelist INPUT)
143	LABEL	scalar	I4	(8), <u>9</u>	control of reading the TPID-record and checking the tape identification number (see 2.2.6 namelist INPUT)
144	LNER	scalar	I4	(8), <u>9</u>	number of points per resonance in EPT (see 2.2.6 namelist INPUT)
145	ITAPE	scalar	I4	(8), <u>9</u>	control of writing the converted data of ENDF/B MF=3 on unit 15 (see 2.2.6 namelist INPUT)
146	MATP	scalar	I4	(8), <u>9</u>	ENDF/B-material number (see 2.2.6 namelist INPUT)
147	EB	scalar	R4	( <u>8</u> ), <u>9</u>	binding energy of the last neutron (in eV) (see 2.2.6 namelist INPUT)
148	NISO	scalar	I4	(8), <u>9</u>	number of isotopes for which data are given in the control input (see 2.2.6 namelists INPUT and ISO)

Name of the COMMON block: INOUT (continuation)					
word number	name	dimension	type	used in programs number	meaning
149	IF4	scalar	I4	(8), <u>9</u>	control of calling the subroutine ISABEL (see 2.2.6 namelist INPUT)
150	IF5	scalar	I4	(8), <u>9</u>	control of calling the subroutine FILE5 (see 2.2.6 namelist INPUT)
151	FAICOD	scalar	R4	<u>9</u> , 67	used in combining two pointwise cross sections in the subroutine COMBP (see 2.3.5.67 and 2.2.6 namelist INPUT)
152	MLSLSW	scalar	I4	<u>9</u> , 12	switch to enforce single level Breit Wigner instead of multi level Breit Wigner (see 2.2.6 namelist INPUT)
153	ANFZEI	scalar	R4	1, 45	value produced at the first ZEIT (2.3.5.97) call
154	ELINEG	scalar	I4	<u>9</u> , 10, 41	deletion of negative SGT and SGN cross section values (see 2.2.6 namelist INPUT)
Name of the COMMON block: RE					
1	XR	(300)	R4	73, <u>74</u>	interchange of data between DANI (2.3.5.74) and CHGINT (2.3.5.73)
301	YR	(300)	R4	73, <u>74</u>	see array XR

## Name of the COMMON block: RECS

word number	name	dimension	type	used in programs number	meaning
1	MAT	scalar	I4	<u>10</u> , (37), 46, <u>48</u> , <u>49</u> , 51, 52, 53, 54, 55, 56, 57, 58, 59, 60, 61, 62, 63, 80, <u>81</u>	ENDF material number (see /3/ chapter R-2)
2	MF	scalar	I4	<u>10</u> , 36, 37, 46, 48, <u>49</u> , 51, 52, 53, 54, <u>55</u> , <u>56</u> , 58, <u>59</u> , <u>60</u> , 61, 62, 80, <u>81</u>	ENDF file number (see /3/ chapter R-2)
3	MT	scalar	I4	<u>10</u> , 36, 37, 46, 48, <u>49</u> , 51, 52, 53, 54, <u>55</u> , <u>56</u> , 58, <u>59</u> , <u>60</u> , 61, 62, 80, <u>81</u>	ENDF section (reaction type) number (see /3/ chapter R-2)
4	C1	scalar	R4	<u>10</u> , 11, 12, 24, 25, <u>37</u> , 44, <u>49</u> , 51, 52, 53, 54, 55, 56, 58, 59, <u>60</u> , 61, 62, 80, <u>81</u>	ENDF constant. (In most cases the temperature (see /3/ chapter R-2)
5	C2	scalar	R4	<u>10</u> , 11, 12, 25, 36, <u>37</u> , 42, 44, <u>49</u> , 51, 52, 53, 54, 55, <u>56</u> , 58, 59, <u>60</u> , 61, 62, 80, <u>81</u>	ENDF constant (see /3/ chapter R-2)
6	L1	scalar	I4	<u>10</u> , 11, 12, 25, <u>37</u> , 42, <u>49</u> , 51, 52, 53, 54, 55, 56, 58, 59, <u>60</u> , 61, 62, 80, <u>81</u>	ENDF integer constant. (In most cases used to indicate temperature dependence) (see /3/ chapter R-2)
7	L2	scalar	I4	<u>10</u> , 11, 25, <u>37</u> , 42, 44, <u>49</u> , 51, 52, 53, 54, 55, <u>56</u> , 58, 59, <u>60</u> , 61, 62, 80, <u>81</u>	ENDF integer constant (see /3/ chapter R-2)
8	N1	scalar	I4	<u>10</u> , 11, 12, <u>24</u> , <u>25</u> , <u>30</u> , <u>36</u> , ( <u>37</u> ), <u>39</u> , 40, 42, ( <u>44</u> ), <u>49</u> , 51, 52, 53, 54, 55, <u>56</u> , 58, 59, 60, 61, 62, <u>66</u> , <u>67</u> , 71, <u>75</u> , 78, <u>79</u> , 80, <u>81</u>	ENDF count of items in a list to follow (see /3/ chapter R-2)



Name of the COMMON block: RECS (continuation)					
word number	name	dimension	type	used in programs number	meaning
9	N2	scalar	I4	<u>10</u> , 12, <u>24</u> , <u>25</u> , <u>30</u> , 34, <u>36</u> , ( <u>37</u> ), <u>39</u> , <u>40</u> , <u>42</u> , ( <u>44</u> ), <u>49</u> , 51, 52, <u>53</u> , <u>54</u> , <u>55</u> , <u>56</u> , 58, 59, 60, 61, 62, <u>66</u> , <u>67</u> , 71, <u>75</u> , 78, <u>79</u> , 80, <u>81</u>	ENDF count of items in a second list to follow (see /3/ chapter R-2)
10	NS	scalar	I4	<u>10</u> , <u>17</u> , <u>37</u> , <u>52</u> , <u>53</u> , <u>54</u> , <u>55</u> , <u>56</u> , <u>57</u> , <u>58</u> , <u>59</u> , <u>60</u> , <u>61</u> , <u>62</u> , <u>63</u>	ENDF sequence number used in MODE=3 writing (see /3/ chapter R-2)
Name of the COMMON block: TAB					
1	ENDTAB	(25)	I4	<u>4</u> , 36, 37	ENDF MF=3 section numbers (MT) of data types to be converted
26	KEDTAB	(50)	I4	<u>4</u> , 36, 37	alphanumerical KEDAK material names associated to the ENDF-section numbers stored in the array ENDTAB
76	KEDR	(240)	I4	<u>4</u> , 42, 44, 77, 86	list of alphanumerical and associated numerical KEDAK data type names
316	NCNAME	(39)	I4	<u>9</u> , <u>42</u> , <u>44</u> , 86	number of incident energy points at which angular distributions and energy distributions of secondary neutrons are given
355	NRIF3	scalar	I4	<u>9</u> , <u>36</u> , 37, 86	number of levels of KEDAK datatype SGIZ
356	NRIF4	scalar	I4	<u>9</u> , <u>42</u> , <u>44</u> , 86	number of datatypes from ENDF MF =4 and MF=5

Name of the COMMON block: XBAR					
word number	name	dimension	type	used in programs number	meaning
1	XX	(4,10)	R4	<u>4</u> , 26	used in the calculation of the fluctuation integrals for capture, fission and elastic scattering

### 2.3.5 Description of the Main Program and the Subroutines

In the following the main program and all subroutines of BRIGITTE will be described. All the descriptions will have the following scheme:

- problem solved by the routine
- (- method of solution)
- definition with argument list (and entry names if existant)
- explanation of the quantities in the list(s)
- used common blocks and used quantities of these common blocks
- called subprograms (except those, which are FORTRAN-supplied procedures or subprograms like IBCOM)

In sections 2.3.5.92 until 2.3.5.99 a description of all auxiliary programs<sup>+</sup>, which are used by the program, will be given.

In the explanation lists of the arguments and in the used common blocks lists the following abbreviations will be used:

in the mode  
column:

- |   |  |
|---|--|
| I | is used for variables which values must be defined before calling the subroutine or must have been set in the common by another program.                         |
| O | is used for variables which are set in the subroutine (or in programs called by the subroutine) and are used outside (and within) the subroutine.                |
| L | is used for arrays (or variables) which are only used for placing some storage for the disposal of the subroutine and have only the meaning of local quantities. |

---

<sup>+</sup>) all these programs are members of local libraries.

P means the same as L with the difference that the arrays or variables are not used in the subroutine itself but passed to other subroutines or functions. The values produced in this subroutines or functions are not used outside this subroutines or functions.

(L and P are only used in the list "Explanation of the argument", but not in the list "Used common blocks".)

IO is used for a combination of I and O.

PL is used for a combination of P and L. That means e.g. the array is used in some parts of the subroutine as local quantity and in others the address is passed to another subroutine.

- is used for quantities, which are not referenced in the subprogram, or for quantities, for which none of the given classifications is true.

in the type  
column:

R4 for REAL\*4 quantities

R8 for REAL\*8 quantities

I4 for INTEGER\*4 quantities

L4 for LOGICAL\*4 quantities

B4 for 4 byte quantities of any type. The actual type will be determined by additional information.

F4 for functions supplying REAL\*4 values

In the dimension column the maximal dimension of arrays will be given. For scalar quantities the word "scalar" is used.

In the "used common blocks" list a column headed with "word" is used. The number in this column means the starting address (measured in 4-byte words and beginning with word 1) of the associate variable or array.

In the "Explanation of the argument" lists often the word energy is used for the argument of a TAB1-record (see /3/ chapter R-2.3.3). This is done, because in nearly all cases this is an energy. But also other quantities are valid as argument, as for example cosine values.

2.3.5.1 Main program

Problem solved:

Control of program flow. DEFINE FILE for direct access data sets.

Used common block:

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NR1	(133)	I4	-	1	IVY	scalar	I4	I	138
IVER	scalar	I4	I	134	IVB	scalar	I4	I	139
IVGENR	scalar	I4	I	135	IVDENS	scalar	I4	I	140
IVJNT	scalar	I4	I	136	NR2	(12)	I4	-	141
IVX	scalar	I4	I	137	ANFZEI	scalar	I4	0	153

Called subprograms:

ZEIT  
TITEL  
PRIEIN  
GSPACE  
DADEFI  
DEFI  
PART1  
SICKY

2.3.5.2 Subroutine TITEL

Problem solved:  
Printing of title page.

Definition  
Subroutine TITEL

Used common block:

name of the common block: INOUT				
name	dimension	type	mode	word
NO	scalar	I4	I	1

Called subprograms:

A8FORM  
RAMANF } (Entries of A8FORM)  
RAMEND }

2.3.5.3 Subroutine PRIEIN

Problem solved:  
Printout of the input.

Definition:  
SUBROUTINE PRIEIN(IEIN,IOUT,ZEILMA)

Explanation of the arguments:

name	mode	type	dimension	remarks
IEIN	I	I4	scalar	input unit number
IOUT	I	I4	scalar	print output unit number
ZEILMA	I	I4	scalar	maximal number of lines per printoutput page

Used common blocks:  
None

Called subprograms:  
None

2.3.5.4 BLOCK DATA

Problem solved:

Initialisation of some quantities in the common blocks.

Definition:

BLOCK DATA

Used common blocks:

name of the common block: TAB									
name	dimension	type	mode	word	name	dimension	type	mode	word
ENDTAB	(25)	I4	0	1	NCNAME	(39)	I4	-	316
KEDTAB	(50)	I4	0	26	NRIF3	scalar	I4	-	355
KEDR	(240)	I4	0	76	NRIF4	scalar	I4	-	356
name of the common block: XBAR									
XX	(4,10)	R4	0	1					
name of the common block: INOUT									
NO	scalar	I4	0	1	N	scalar	I4	-	129
NI	scalar	I4	0	2	NBKX	scalar	I4	0	130
NR1	(123)	I4	-	3	JX	scalar	I4	0	131
NBX	scalar	I4	0	126	IADIM	scalar	I4	0	132
N1X	scalar	I4	0	127	ITDIM	scalar	I4	0	133
N2X	scalar	I4	0	128	NR2	(21)	I4	-	134

Called subprograms:

None

2.3.5.5 Subroutine GSPACE

Problem solved:

Allocation of all free space in the main storage region as data working space.

Method of solution:

First the size of the region, which is not used by the program or for buffers, is determined. From this value all the region, which may be used for buffers later on, and the region for other system use is subtracted. The remaining value is used as data storage. The parcellation of this storage for the different arrays is also controlled by this routine.

Definition:

SUBROUTINE GSPACE(F)

Explanation of the arguments:

name	mode	type	dimension	remarks
F	L	R4	(1)	the subroutine uses this address as reference point

Used common block:

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NO	scalar	I4	I	1	IADIM	scalar	I4	IO	132
NI	scalar	I4	I	2	ITDIM	scalar	I4	IO	133
NR1	(123)	I4	-	3	IVER	scalar	I4	0	134
NBX	scalar	I4	IO	126	IVGENR	scalar	I4	0	135
N1X	scalar	I4	IO	127	IVJNT	scalar	I4	0	136
N2X	scalar	I4	IO	128	IVX	scalar	I4	0	137
N	scalar	I4	-	129	IVY	scalar	I4	0	138
NBKX	scalar	I4	IO	130	IVB	scalar	I4	0	139
JX	scalar	I4	IO	131	IVDENS	scalar	I4	0	140

Called subprograms:

BUFREG  
FREESP  
XTAREA



2.3.5.6 Subroutine BUFREG

Problem solved:

Determination of the region needed for buffers.

Method of solution:

All DD(data definition)-names in the TIOT (task input output table) of the form FTXXFOO1 with the attached BLKSIZE (block size) are fetched (for explanation of the terms see /10/ and other IBM system literature), the number of buffers is set to two, and with this values the place for buffers is calculated. If no values are found in the TIOT, they are fetched from the label.

Exceptions:

The file for print output (the file number is expected as first Integer\*4 byte word in the named common block INOUT) is not considered.

If there are the DD-names SYSUDUMP or SYSABEND, two k-byte will be added to the buffer region for each one.

Definition:

BUFREG START

The program is written in IBM/370 Assembler. It is called from FORTRAN-programs by the statement CALL BUFREG (INTREG,&LABEL)

Explanation of the arguments:

name	mode	type	dimension	remarks
INTREG	0	I4	scalar	region-measured in k-byte-needed for buffers. If INTREG is less than zero, the program has found an error. In this case RETURN 1 is made
&LABEL	-	-	-	label for RETURN 1

Used common block:

name of the common block: INOUT				
name	dimension	type	mode	word
NO	scalar	I4	I	1

Called subprograms:

None

### 2.3.5.7 Subroutine DADEFI

Problem solved:

The subparameters dblk1 (data block length) and nblk (number of blocks) out of SPACE = (dblk1,(nblk)) are fetched from the DD-cards FT01FO01, FT11FO01, and FT12FO01. They will be used as input in the dynamic DEFINE FILE statement program DEFI, which is called in the MAIN program.

Definition:

SUBROUTINE DADEFI(IF)

Explanation of the argument:

name	mode	type	dimension	remarks
IF	0	I4	(6)	dblk1 and nblk (see problem solved) are stored in IF  IF(1) and IF(4): dblk1 and nblk of FT01FO01.  IF(2) and IF(5): dblk1 and nblk of FT11FO01.  IF(3) and IF(6): dblk1 and nblk of FT12FO01

Used common blocks:

None.

Called subprograms:

DINF

### 2.3.5.8 Subroutine PART1

Problem solved:

Control of program flow.

Definition:

SUBROUTINE PART1(NIS,MISO,KIF,KIF5,BK,NBT,JNT,X,Y,B,A)

Explanation of the arguments:

name	mode	type	dimension	remarks
NIS	O	I4	scalar	number of isotopes for the specified ENDF/B-material
MISO	O	I4	scalar	number of isotopes for which data are specified in the control input
KIF	O	I4	scalar	value of IF4 (default value or given in the control input) (see input description)
KIF5	O	I4	scalar	value of IF5 (default value or given in the control input) see input description 2.2.6)
BK	P	R4	(NBKX)	array and dimension see 2.3.3
NBT	P	I4	(N1X)	" " " "
JNT	P	I4	(N1X)	" " " "
X	P	R4	(N2X)	" " " "
Y	P	R4	(N2X)	" " " "
B	P	R4	(NBX)	" " " "
A	P	R4	(IADIM)	" " " "

Used common block:

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NR1	(2)	I4	-	1	MATP	scalar	I4	0	146
NT	scalar	I4	I	3	EB	scalar	R4	0	147
NR2	(139)	I4	-	4	NISO	scalar	I4	0	148
LABEL	scalar	I4	0	143	IF4	scalar	I4	0	149
LNER	scalar	I4	0	144	IF5	scalar	I4	0	150
ITAPE	scalar	I4	0	145					

Called subprograms:

```

DENISE
EITAB1      (entry of ITAB1)
FILE5
INGRID
INITNL
IRREC      (entry of RREC)
ISABEL
IWREC      (entry of WREC)
REGINE
SUZY
WDA1      }
WDA3      }      (entries of WDA)
WRECS1    (entry of WRECS)
    
```

2.3.5.9 Subroutine INITNL

Problem solved:

Default values are set for most of the variables in the namelist input. The control input (namelist(s) INPUT and eventually ISO) is read.

Definition:

SUBROUTINE INITNL(MISO,\*)

Explanation of the arguments:

name	mode	type	dimension	remarks
MISO	0	I4	scalar	number of isotopes for which data are given in the control input (MISO<1 has the same meaning as MISO=1)
*	-	-	-	RETURN1 is made, if no (or nomore) namelist INPUT can be found in the control input

Used common blocks:

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NO	scalar	I4	0	1	EBISO	(10)	R4	0	115
NI	scalar	I4	0	2	ICOMB	scalar	I4	0	125
NT	scalar	I4	0	3	NR1	(15)	I4	-	126
MODE	scalar	I4	0	4	EPSRES	scalar	R4	0	141
LEDIT1	scalar	I4	0	5	SCREPS	scalar	R4	0	142
LEDIT2	scalar	I4	0	6	LABEL	scalar	I4	0	143
LEDITS	scalar	I4	0	7	LNER	scalar	I4	0	144
KEDAWE	scalar	I4	0	8	ITAPE	scalar	I4	0	145
ABIS	(20)	R4	-	9	MATP	scalar	I4	0	146
NOTTRL	(25)	R8	0	29	EB	scalar	R4	0	147
BIB12	scalar	I4	0	79	NISO	scalar	I4	0	148
SALINE	scalar	L4	0	80	IF4	scalar	I4	0	149
NAMKDK	scalar	R8	0	81	IF5	scalar	I4	0	150
MATKDK	scalar	I4	0	83	FAICOD	scalar	R4	0	151
NTYP	scalar	I4	-	84	MLSLSW	scalar	I4	0	152
NAMISO	(10)	R8	0	85	NR3	(1)	I4	-	153
MATISO	(10)	I4	0	105	ELINEG	scalar	I4	0	154

name of the common block: TAB									
NR2	(315)	I4	-	1	NRIF3	scalar	I4	0	355
NCNAME	(39)	I4	0	316	NRIF4	scalar	I4	0	356

Called subprograms:

None.

2.3.5.10 Subroutine INGRID

Problem solved:

Reads and checks the ENDF-label. Prints some text. Reads MF=1 MT=451 (Descriptive Data and Dictionary) and prints this section if desired. If existant MF=1 MT=452 (number of neutrons per fission,  $\bar{\nu}$ ) is read, printed out and stored for later processing.

Definition:

SUBROUTINE INGRID(MATP,LABEL,LRP,LFI,LNU,IF4,IF5)

Explanation of the arguments:

name	mode	type	dimension	remarks
MATP	I	I4	scalar	number of the ENDF/B-material to be converted
LABEL	I	I4	scalar	see 2.2.6 (input description)
LRP	O	I4	scalar	flag to indicate the existance of resolved and/or unresolved resonance data (see /2/)
LFI	O	I4	scalar	LFI = 0 material is <u>not</u> fissionable LFI = 1 material is fissionable
LNU	O	I4	scalar	LNU = 1 polynomial representation of $\bar{\nu}$ . LNU = 2 tabulated representation of $\bar{\nu}$ .
IF4	I	I4	scalar	see 2.2.6 (input description)
IF5	I	I4	scalar	" " "

Used common blocks:

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NO	scalar	I4	I	1	KEDAVE	scalar	I4	I	8
NI	scalar	I4	-	2	NR2	(72)	I4	-	9
NT	scalar	I4	I	3	NAMKDK	scalar	R8	I	81
MODE	scalar	I4	I	4	MATKDK	scalar	I4	I	83
LEDIT1	scalar	I4	I	5	NR3	(70)	I4	-	84
NR1	(2)	I4	-	6	ELINEG	scalar	I4	I	154

name of the common block: RECS									
MAT	scalar	I4	I	1	L1	scalar	I4	I	6
MF	scalar	I4	I	2	L2	scalar	I4	I	7
MT	scalar	I4	-	3	N1	scalar	I4	I	8
C1	scalar	R4	-	4	N2	scalar	I4	I	9
C2	scalar	R4	-	5	NS	scalar	I4	0	10

(The first nine words in the Common RECS are set to zero at the beginning of the subroutine.)

Called subprograms:

A8FORM		WREC	
ERROR		WRECS2	} (entries of WRECS)
PRIZEI		WRECS4	
RAMANF	} (entries of A8FORM)		
RAMEND			
RREC			
SEARCH			

2.3.5.11 Subroutine REGINE

Problem solved:

Processing of ENDF MF=2 (resonance data). Generation of ISOT1 and ISOT2 (and eventually ISOT3) for KEDAK.

Definition:

SUBROUTINE REGINE(MATP,LNER,LISRES,LISUNR,NIS,LFWRES,LFW,EB,LRP,NISO,BK,  
NBT,JNT,X,Y,B,AH)

Explanation of the arguments:

name	mode	type	dimension	remarks
MATP	I	I4	scalar	number of the ENDF/B-material, which is to be converted
LNER	I	I4	scalar	see 2.2.6 (input description)
LISRES	0	I4	scalar	= 0 cross sections in the resolved resonance region are to be calculated from resonance parameters plus smooth cross sections from file 3. = 1 only smooth cross sections in file 3 are used.
LISUNR	0	I4	scalar	= 0 cross sections in the unresolved resonance region are to be calculated from resonance parameters plus smooth cross sections from file 3. = 1 only smooth cross sections in file 3 are used.
NIS	0	I4	scalar	number of isotopes in ENDF MF=2 (resonance data)
LFWRES	0	I4	scalar	= 0 zero fission widths = 1 non zero fission widths in the resolved resonance region
LFW	0	I4	scalar	= 1 fission widths are given = 0 fission widths are <u>not</u> given } in the unresolved resonance region
EB	I	R4	scalar	see input description 2.2.6. In the case of more than one isotope EB is set to EBISO(MIS) (see input description 2.2.6) after storing the original value of EB in the ISOT2 data-block of the entire material



name	mode	type	dimension	remarks
LRP	I	I4	scalar	see 2.3.5.10 (explanation of the arguments)
NISO	I	I4	scalar	number of isotopes for which data are given in the input (see 2.2.6)
BK	PL	R4	(NBKX)	see 2.3.3
NBT	P	I4	(NIX)	"
JNT	P	I4	(NIX)	"
X	P	R4	(N2X)	"
Y	P	R4	(N2X)	"
B	P	R4	(NBX)	"
AH	P	R4	(JX)	"

Used common blocks:

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NO	scalar	I4	I	1	MATKDK	scalar	I4	I	83
NI	scalar	I4	-	2	NR3	(21)	I4	-	84
NT	scalar	I4	I	3	MATISO	(10)	I4	I	105
MODE	scalar	I4	I	4	EBISO	(10)	R4	I	115
NR1	(4)	I4	-	5	NR4	(4)	I4	-	125
ABIS	(20)	R4	0	9	N	scalar	I4	0	129
NR2	(52)	I4	-	29	NBKX	scalar	I4	I	130
NAMKDK	scalar	R8	I	81	JX	scalar	I4	I	131

name of the common block: RECS									
name	dimension	type	mode	word	name	dimension	type	mode	word
MAT	scalar	I4	-	1	C2	scalar	R4	I	5
MF	scalar	I4	-	2	L1	scalar	I4	I	6
MT	scalar	I4	-	3	L2	scalar	I4	I	7
C1	scalar	R4	I	4	N1	scalar	I4	I	8

name of the common block: FLCT									
NR5	(6)	I4	-	1	NPEMAX	scalar	I4	0	8
NDIM6	scalar	I4	0	7					

Called subprograms:

ANNICK  
A8FORM  
ERROR  
MARY  
PRIZEI  
RREC  
SEARCH  
WDA2 (entry of WDA)  
XTEXT

#### 2.3.5.12 Subroutine MARY

Problem solved:

Control of calculation of pointwise cross sections out of resolved resonance parameters.

Definition:

SUBROUTINE MARY(EL,EH, LRF,ZAI,ABN, LFWRES,NIS,LNER,MIS,LIS,SPI,AWRI,AP,BK,  
NBT,JNT,ENS,SIGS,B,AH,JXRED,ER,AJ,GT,GN,GG,GF,AL,G,SER,  
PER,NDIM)

Explanation of the arguments:

name	mode	type	dimension	remarks
EL	I	R4	scalar	lower energy (eV) limit for the resolved resonance region
EH	I	R4	scalar	upper energy limit (eV) for the resolved resonance region
LRF	I	I4	scalar	= 1 single level Breit-Wigner = 2 multi level Breit-Wigner
ZAI	I	R4	scalar	(Z,A) designation (see /2/)
ABN	I	R4	scalar	fractional abundance (see /2/)
LFWRES	O	I4	scalar	= 1 non zero fission widths given (as resolved resonance parameters)
NIS	I	I4	scalar	number of isotopes in ENDF MF=2 (resonance data)
LNER	I	I4	scalar	see input description (2.2.6)
MIS	I	I4	scalar	numeration of the isotope worked with
LIS	O	I4	scalar	see LISRES in 2.3.5.11 (explanation of the arguments)
SPI	O	R4	scalar	nuclear spin of the target nucleus (see /2/)
AWRI	O	R4	scalar	ratio of the mass of a particular isotope to that of a neutron (see /2/)
AP	O	R4	scalar	effective scattering radius (see /2/)
BK	P	R4	(NBKX)	see 2.3.3
NBT	P	I4	(N1X)	"
JNT	P	I4	(N1X)	"
ENS	L	R4	(N2X)	energy grid (see 2.3.3 array X)
SIGS	P	R4	(N2X)	see 2.3.3 array Y

name	mode	type	dimension	remarks
B	L	R4	(NBX)	resonance parameters (see 2.3.3)
AH	L	R4	(JXRED)	used for storing SGG,SGN and eventually SGF pointwise cross sections
JXRED	I	I4	scalar	JXRED = JX - JX/20
ER	L	R4	(NDIM)	resonance energy (see /2/)
AJ	L	R4	(NDIM)	spin of the resonance (see /2/)
GT	L	R4	(NDIM)	resonance total width $\Gamma$ evaluated at the resonance energy ER (see /2/)
GN	L	R4	(NDIM)	neutron width $\Gamma_n$ (see /2/)
GG	L	R4	(NDIM)	radiation width $\Gamma_\gamma$ (see /2/)
GF	L	R4	(NDIM)	fission width $\Gamma_f$ (see /2/)
AL	L	R4	(NDIM)	neutron angular momentum quantum number l (see /2/)
G	L	R4	(NDIM)	$(2*AJ(k)+1)/(2*(2*SPI+1))$ (see /2/)
SER	L	R4	(NDIM)	shift factors (see /2/)
PER	L	R4	(NDIM)	penetration factors (see /2/)
NDIM	I	I4	scalar	NDIM = JX/200 (JX see 2.3.3)

Used common blocks:

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NO	scalar	I4	I	1	NREF2	(22)	I4	-	105
NI	scalar	I4	-	2	N1X	scalar	I4	I	127
NT	scalar	I4	I	3	N2X	scalar	I4	I	128
MODE	scalar	I4	I	4	N	scalar	I4	-	129
LEDIT1	scalar	I4	-	5	NBKX	scalar	I4	I	130
LEDIT2	scalar	I4	I	6	NREF4	(21)	I4	-	131
NREF1	(78)	I4	-	7	MLSLSW	scalar	I4	I	152
NAMISO	(10)	R8	I	85					

name of the common block: RECS									
name	dimension	type	mode	word	name	dimension	type	mode	word
NREF3	(3)	I4	-	1	L2	scalar	I4	-	7
C1	scalar	R4	I	4	N1	scalar	I4	I	8
C2	scalar	R4	I	5	N2	scalar	I4	I	9
L1	scalar	I4	I	6					

Called subprograms:

EPT  
ERROR  
FACTS  
MARCEL  
MARGOT (entry of MARCEL)  
MARYO1  
MARYO2  
MELA  
MILLI  
PRIZEI  
RREC  
SORTRE

2.3.5.13 Subroutine FACTS

Problem solved:

Calculates penetration and shift factors liberated from Gregson et al.

Definition:

SUBROUTINE FACTS(L,RHO,SE,PE)

Explanation of the arguments:

name	mode	type	dimension	remarks
L	I	I4	scalar	l-value (see AL in 2.3.5.12)
RHO	I	R4	scalar	see $\rho$ in /2/ page D-3
SE	O	R4	scalar	shift-factor (see /2/ page D-3)
PE	O	R4	scalar	penetration-factor (see /2/ page D-4)

Used common blocks:

None

Called subprograms:

None

#### 2.3.5.14 Subroutine SORTRE

Problem solved:

Sorts (resolved) resonance data on increasing energy.

Definition:

SUBROUTINE SORTRE(NRES,ER,AJ,GT,GN,GG,GF,AL,G,SER,PER)

Explanation of the arguments:

name	mode	type	dimension	remarks
NRES	I	I4	scalar	number of (resolved) resonances
ER	IO	R4	(NDIM)	see 2.3.5.12 (argument list)
AJ	IO	R4	(NDIM)	" "
GT	IO	R4	(NDIM)	" "
GN	IO	R4	(NDIM)	" "
GG	IO	R4	(NDIM)	" "
GF	IO	R4	(NDIM)	" "

name	mode	type	dimension	remarks
AL	IO	R4	(NDIM)	see 2.3.5.12 (argument list)
G	IO	R4	(NDIM)	" "
SER	IO	R4	(NDIM)	" "
PER	IO	R4	(NDIM)	" "

Used common blocks:

None

Called subprograms:

None

### 2.3.5.15 Subroutine MARYO1

Problem solved:

Stores (resolved) resonance data in the array BK. Changes  $\Gamma_n$  for negative resonance energies. Optional printout of resolved resonance data in KEDAK representation. Printout of KEDAK type RANGRES. Stores resolved resonance data on file 11 with the name RES. Determines if fission widths are zero.

Definition:

SUBROUTINE MARYO1 (EL, EH, NRES, BK, ER, AL, AJ, G, GT, GN, GG, GF, NIS, MIS, LIS, LFWRES, \*)

Explanation of the arguments:

name	mode	type	dimension	remarks
EL	I	R4	scalar	see 2.3.5.12 argument list
EH	I	R4	scalar	" "
NRES	I	I4	scalar	number of (resolved) resonances
BK	L	R4	(NBKX)	local storage for (resolved) resonance data and other quantities
ER	I	R4	(NDIM)	see 2.3.5.12 (argument list)

name	mode	type	dimension	remarks
AL	I	R4	(NDIM)	see 2.3.5.12 (argument list)
AJ	I	R4	(NDIM)	" "
G	I	R4	(NDIM)	" "
GT	I	R4	(NDIM)	" "
GN	I	R4	(NDIM)	" "
GG	I	R4	(NDIM)	" "
GF	I	R4	(NDIM)	" "
NIS	I	I4	scalar	number of isotopes in ENDF MF=2 (resonance data)
MIS	I	I4	scalar	numeration of the isotope actually worked with
LIS	I	I4	scalar	see LISRES in 2.3.5.11 (argument list)
LFWRES	0	I4	scalar	= 1 nonzero fission widths given (as resolved resonance parameters) = 0 all fission widths are zero
*	-	-	-	RETURN 1 is made for LIS not equal zero. But LIS is always zero after the program MARYO1 has been called. Therefore RETURN 1 is never made.

Used common block:

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NO	scalar	I4	I	1	NAMISO	(10)	R8	I	85
NR1	(4)	I4	-	2	MATISO	(10)	I4	I	105
LEDIT2	scalar	I4	I	6	NR4	(14)	I4	-	115
NR2	(76)	I4	-	7	N	scalar	I4	0	129
MATKDK	scalar	I4	I	83	NBXX	scalar	I4	I	130
NR3	(1)	I4	-	84					



Called subprograms:

A8FORM

ERROR

LUCY

WDA2 (entry of WDA)

XTEXT

2.3.5.16 Subroutine LUCY

Problem solved:

Prints resolved or statistical resonance data, which are stored in the array BK, with a defined format.

Definition:

SUBROUTINE LUCY(I,BK)

Explanation of the arguments:

name	mode	type	dimension	remarks
I	I	I4	scalar	= 2 resolved resonance data format = 3 statistical data format
BK	I	R4	(NBKX)	data to be printed

Used common block:

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NO	scalar	I4	I	1	N	scalar	I4	I	129
NR1	(127)	I4	-	2					

Called subprograms:

ERROR

2.3.5.17 Subroutine EPT

Problem solved:

Calculates energy mesh based on NPR points per resonance.

Method of solution:

See comment statements in the source program listing within the subroutine EPT.

Definition:

SUBROUTINE EPT(EMIN,EMAX,NPR,NTP,NR,ER,GTR,ENS,N2X)

Explanation of the arguments:

name	mode	type	dimension	remarks
EMIN	I	R4	scalar	lower limit for the energy mesh
EMAX	I	R4	scalar	upper limit for the energy mesh
NPR	I	I4	scalar	(maximum) number of points used per resonance
NTP	0	I4	scalar	number of points to describe NR resonances from EMIN to EMAX
NR	I	I4	scalar	number of resonances
ER	I	R4	(NR)	resonance energies
GTR	I	R4	(NR)	total widths of resonances
ENS	0	R4	(N2X)	calculated energy mesh
N2X	I	I4	scalar	maximum dimension of array ENS

Used common block:

name of the common block: RECS									
name	dimension	type	mode	word	name	dimension	type	mode	word
NR1	(9)	I4	-	1	NS	scalar	I4	L	10

Called subprograms:

RESERR

2.3.5.18 Subroutine RESERR

Problem solved:

Prints error messages for the resolved resonance calculating program section.

Definition:

SUBROUTINE RESERR(N,K,NSE,IERR)

Explanation of the arguments:

name	mode	type	dimension	remarks
N	I	I4	scalar	selection of different error messages
K	I	I4	scalar	number to be printed in the subroutine
NSE	I	I4	scalar	number to be printed in the subroutine
IERR	I	I4	scalar	selection of different error messages

Used common block:

name of the common block: INOUT				
name	dimension	type	mode	word
IO	scalar	I4	I	1

Called subprograms:

None

2.3.5.19 Subroutine MARYO2

Problem solved:

Deletes all but one energy point with the same value.

Definition:

SUBROUTINE MARYO2(ENS,NSE)

Explanation of the arguments:

name	mode	type	dimension	remarks
ENS	IO	R4	(N2X)	energy points
NSE	IO	I4	scalar	number of energy points

Used common blocks:

None

Called subprograms:

None

#### 2.3.5.20 Subroutine MARCEL

Problem solved:

Calculates for one energy point E with a single level (or multi level) Breit-Wigner formulae elastic, radiative capture and depending on LFWRES fission cross sections. For LRF = 1 the single level and for LRF = 2 the multi level formalism is used.

Definition:

For initialisation:

```
SUBROUTINE MARCEL(LFWRES, LRF, C, CON3, CON4, PIM4, AP, NRES, NLS, LLSTR, ER, GN, GG,  
                  GF, G, AL, SER, PER)
```

For calculation:

```
ENTRY MARGOT(E, SGN, SGG, SGF)
```

Explanation of the arguments:

name	mode	type	dimension	remarks
LFWRES	I	I4	scalar	only for LFWRES=1 fission cross section are calculated
LRF	I	I4	scalar	for LRF=2 a multi level Breit Wigner formulae is used. Else a single level Breit Wigner formulae is used
C	I	R4	scalar	$= 2.196771E-3 \cdot AP \cdot \frac{AWRI}{AWRI+1.0}$ (AP and AWRI see 2.3.5.12 argument list)
CON3	I	R4	scalar	$= 2.196771E-3 \cdot \frac{AWRI}{AWRI+1.0} \cdot 0.1 \cdot (0.8+1.23 \cdot AWRI^{1/3})$ (AWRI see 2.3.5.12 argument list)
CON4	I	R4	scalar	$= 4\pi / (2.196771E-3)^2 \cdot \left(\frac{AWRI+1.0}{AWRI}\right)^2$ (AWRI see 2.3.5.12 argument list)
PIM4	I	R4	scalar	$= 4\pi$
AP	I	R4	scalar	see 2.3.5.12 argument list
NRES	I	I4	scalar	number of (resolved) resonances
NLS	I	I4	scalar	number of l-states (see /2/ page 7.5)
LLSTR	I	I4	(5)	l-values for the different l-states (see /2/ page 7.5)
ER	I	R4	(NDIM)	resonance energies (see /2/ page 7.5) (NDIM see 2.3.5.12 argument list)
GN	I	R4	(NDIM)	neutron width $\Gamma_n$
GG	I	R4	(NDIM)	radiation width $\Gamma_\gamma$
GF	I	R4	(NDIM)	fission width $\Gamma_f$
G	I	R4	(NDIM)	see 2.3.5.12 argument list
AL	I	R4	(NDIM)	" "
SER	I	R4	(NDIM)	shift factors (see /2/ page D-3)
PER	I	R4	(NDIM)	penetration factors (see /2/ page D-3)
E	I	R4	scalar	energy for which the pointwise cross sections are to be calculated
SGN	O	R4	scalar	$\sigma_e(E)$
SGG	O	R4	scalar	$\sigma_\gamma(E)$
SGF	O	R4	scalar	$\sigma_f(E)$ (only for LFWRES=1)

Used common blocks:

None

Called subprograms:

ERROR

FACPHI

FACTS

### 2.3.5.21 Subroutine FACPHI

Problem solved:

Calculates phase shift for the Breit Wigner formulae (see /2/ page D-4).

Definition:

SUBROUTINE FACPHI(L,RHO,PHI)

Explanation of the arguments:

name	mode	type	dimension	remarks
L	I	I4	scalar	l-value (see /2/ page 7.5)
RHO	I	R4	scalar	= $k \cdot AP$ (see /2/ page D-3 and D-4)
PHI	O	R4	scalar	calculated phase shift

Used common blocks:

None

Called subprograms:

None

### 2.3.5.22 Subroutine MILLI

Problem solved:

Checks if there are enough points describing the point resonance cross sections. If not the number of points will be enlarged. Also the best interpolation scheme is determined.

Definition:

SUBROUTINE MILLI(ENS,NSE,AH,NBTN,JNTN,NBTG,JNTG,NBTF,JNTF,LFWRES,NISGN,  
NISGG,NISGF,JXRED)

Explanation of the arguments:

name	mode	type	dimension	remarks
ENS	IO	R4	(N2X)	energy grid
NSE	IO	I4	scalar	number of energy points in ENS
AH	IO	R4	(JXRED)	1 ... NSE : $\sigma_e$ NSE+1 ... 2*NSE: $\sigma_\gamma$ For LFWRES = 1 2*NSE+1 ... 3*NSE: $\sigma_f$
NBTN	0	I4	(N1X)	interpolation ranges for $\sigma_e$
JNTN	0	I4	(N1X)	interpolation scheme identification number for $\sigma_e$ (in the resolved resonance region)
NBTG	0	I4	(N1X)	interpolation ranges for $\sigma_\gamma$
JNTG	0	I4	(N1X)	interpolation scheme identification number for $\sigma_\gamma$
NBTF	0	I4	(N1X)	interpolation ranges for $\sigma_f$
JNTF	0	I4	(N1X)	interpolation scheme identification number for $\sigma_f$
LFWRES	I	I4	scalar	Only for LFWRES=1 fission cross section are processed
NISGN	0	I4	scalar	number of interpolation ranges for $\sigma_e$
NISGG	0	I4	scalar	number of interpolation ranges for $\sigma_\gamma$
NISGF	0	I4	scalar	number of interpolation ranges for $\sigma_f$
JXRED	I	I4	scalar	maximal usable extent of array AH (measured in 4 byte words)

Used common block:

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
IOUT	scalar	I4	I	1	NR2	(12)	I4	-	129
NR1	(126)	I4	-	2	EPS	scalar	R4	I	141
N2X	scalar	I4	I	128					

Called subprograms:

ERROR  
MARGOT  
MILLI1  
RESERR  
TERP1

### 2.3.5.23 Subroutine MILLI1

Problem solved:

Inserts interpolation code IC and interpolation range K in JNT resp. NBT.

Definition:

SUBROUTINE MILLI1(NBT,JNT,N1,IC,K,NSE)

Explanation of the arguments:

name	mode	type	dimension	remarks
NBT	0	I4	(N1X)	interpolation ranges
JNT	IO	I4	(N1X)	interpolation codes
N1	IO	I4	scalar	number of interpolation ranges resp. interpolation codes
IC	I	I4	scalar	interpolation code to be inserted in JNT
K	I	I4	scalar	interpolation range to be inserted in NBT
NSE	I	I4	scalar	number of energy points processed until now



Used common block:

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NR1	(126)	I4	-	1	N1X	scalar	I4	I	127

Called subprograms:

RESERR

### 2.3.5.24 Subroutine MELA

Problem solved:

Prints and stores (on unit 12) the calculated pointwise cross sections  $\sigma_e$ ,  $\sigma_\gamma$  and  $\sigma_f$  (for LFWRES=1) in the resolved resonance region.

Definition:

SUBROUTINE MELA(LFWRES,IBK,AH,NBT,JNT,SIGS,N1SGN,N1SGG,N1SGF,NSE,MIS,EL,EH)

Explanation of the arguments:

name	mode	type	dimension	remarks
LFWRES	I	I4	scalar	fission cross sections are only processed for LFWRES=1
IBK	I	I4	(NBKX)	1 ... N1X: NBT for $\sigma_e$ N1X+1 ... 2*N1X: JNT for $\sigma_e$ 2*N1X+1 ... 3*N1X: NBT for $\sigma_\gamma$ 3*N1X+1 ... 4*N1X: JNT for $\sigma_\gamma$ only for LFWRES=1: 4*N1X+1 ... 5*N1X: NBT for $\sigma_f$ 5*N1X+1 ... 6*N1X: JNT for $\sigma_f$ (6*N1X<NBKX)
AH	I	R4	(JXRED)	$\sigma_e, \sigma_\gamma$ and for LFWRES=1 $\sigma_f$ (see 2.3.5.22 argument list)
NBT	0	I4	(N1X)	interpolation ranges
JNT	0	I4	(N1X)	interpolation codes

name	mode	type	dimension	remarks
SIGS	0	R4	(N2X)	pointwise cross sections
NISGN	I	I4	scalar	see 2.3.5.22 argument list
NISGG	I	I4	scalar	" " " "
NISGF	I	I4	scalar	" " " "
NSE	I	I4	scalar	number of energy values in the array X
MIS	I	I4	scalar	actual isotope numeration
EL	I	R4	scalar	lower limit for the resolved resonance region
EH	I	R4	scalar	upper limit for the resolved resonance region

Used common blocks:

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NO	scalar	I4	I	1	NR2	(119)	I4	-	8
NR1	(5)	I4	-	2	N1X	scalar	I4	I	127
LEDITS	scalar	I4	I	7	N2X	scalar	I4	I	128

name of the common block: RECS									
name	dimension	type	mode	word	name	dimension	type	mode	word
NR3	(3)	I4	-	1	N1	scalar	I4	0	8
C1	scalar	R4	0	4	N2	scalar	I4	0	9
NR4	(3)	I4	-	5					

Called subprograms:

ERROR

WREC

WRECS2 (entry of WRECS)

2.3.5.25 Subroutine ANNICK

Problem solved:

Calculation of pointwise cross sections in the unresolved resonance region out of average resonance parameters (see /2/ pages D-9 to D-17).

Definition:

SUBROUTINE ANNICK(EL,EH,LRF,ZAI,ABN,LFW,NIS,MIS,LIS,SPI,AWRI,AP,BK,NBT,JNT,  
BK2,Y,B,SIGS,E,SIGG,SIGF,SIGT,ESC,XT,YT,TEMP,D,GX,GNO,GG,  
GF,AJ)

Explanation of the arguments:

name	mode	type	dimension	remarks
EL	I	R4	scalar	lower limit for the unresolved resonance range
EH	I	R4	scalar	upper limit for the unresolved resonance range
LRF	I	I4	scalar	= 1 only average fission widths are energy-dependant = 2 average level spacing, competitive reaction widths, reduced neutron widths, radiation widths, and fission widths are energy dependant
ZAI	-	R4	scalar	not used in ANNICK
ABN	-	R4	scalar	not used in ANNICK
LFW	I	I4	scalar	= 0 average fission widths are <u>not</u> given = 1 average fission widths are given in the unresolved resonance region
NIS	I	I4	scalar	number of isotopes
MIS	I	I4	scalar	pointer to the actual isotope, if NIS is greater than 1
LIS	0	I4	scalar	is set to zero
SPI	0	R4	scalar	nuclear spin of target nucleus
AWRI	0	R4	scalar	ratio of the mass of the particular isotope to that of the neutron
AP	-	R4	scalar	not used in ANNICK
BK	L	R4	(NBKX)	unresolved resonance parameters

name	mode	type	dimension	remarks
NBT	L	I4	(N1X)	interpolation ranges
JNT	L	I4	(N1X)	interpolation codes
BK2	L	R4	(N2X)	(see array X in 2.3.3) data for KEDAK-type STGF (see /4/ p. 44)
Y	P	R4	(N2X)	-
B	L	R4	(NBX)	is filled in RREC with the unresolved resonance data
SIGS	L	R4	(NPEMAX)	(part of the array AH from word 1 until word NPEMAX) elastic (pointwise) scattering cross sections. (NPEMAX see common FLCT)
E	L	R4	(NPEMAX)	(AH(NPEMAX+1)...AH(2*NPEMAX)) energy grid in the unresolved resonance region with any interpolation code
SIGG	L	R4	(NPEMAX)	(AH(2*NPEMAX+1)...AH(3*NPEMAX)) radiative capture cross sections
SIGF	L	R4	(NPEMAX)	(AH(3*NPEMAX+1)...AH(4*NPEMAX)) fission cross sections
SIGT	L	R4	(NPEMAX)	(AH(4*NPEMAX+1)...AH(5*NPEMAX)) total cross sections
ESC	L	R4	(NPEMAX)	(AH(5*NPEMAX+1)...AH(6*NPEMAX)) energy grid in the unresolved resonance region with linear interpolation
XT	L	R4	(NPEMAX)	(AH(6*NPEMAX+1)...AH(7*NPEMAX)) local storage of energy grids
YT	L	R4	(NPEMAX)	(AH(7*NPEMAX+1)...AH(8*NPEMAX)) local storage of $\overline{\Gamma}_n^\sigma$ or $\overline{\Gamma}_f$
TEMP	P	R4	(NPEMAX)	(AH(8*NPEMAX+1)...AH(9*NPEMAX)) auxiliary array
D	L	R4	(NDIM6,NPEMAX)	(AH(9*NPEMAX+1)...AH(15*NPEMAX)) (NDIM6 has the value 6) mean level spacing
GX	L	R4	(NDIM6,NPEMAX)	(AH(15*NPEMAX+1)...AH(21*NPEMAX)) average competitive reaction width
GNO	L	R4	(NDIM6,NPEMAX)	(AH(21*NPEMAX+1)...AH(27*NPEMAX)) average reduced neutron width

name	mode	type	dimension	remarks
GG	L	R4	(NDIM6,NPEMAX)	(AH(27*NPEMAX+1)...AH(33*NPEMAX)) average radiation width
GF	L	R4	(NDIM6,NPEMAX)	(AH(33*NPEMAX+1)...AH(39*NPEMAX)) average fission width
AJ	L	R4	(NDIM6,NPEMAX)	(AH(39*NPEMAX+1)...AH(45*NPEMAX)) floating point value of the j-state

Used common blocks:

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NO	scalar	I4	I	1	MATISO	(10)	I4	-	105
NI	scalar	I4	-	2	NR2	(11)	I4	-	115
NT	scalar	I4	I	3	NBX	scalar	I4	-	126
MODE	scalar	I4	I	4	NIX	scalar	I4	I	127
LEDIT1	scalar	I4	-	5	N2X	scalar	I4	I	128
LEDIT2	scalar	I4	I	6	N	scalar	I4	0	129
NR1	(76)	I4	-	7	NBKX	scalar	I4	I	130
MATKDK	scalar	I4	-	83	NR3	(10)	I4	-	131
NTYP	scalar	I4	-	84	EPS	scalar	R4	I	141
NAMISO	(10)	R8	I	85					

name of the common block: FLCT									
name	dimension	type	mode	word	name	dimension	type	mode	word
ZYN	scalar	R4	0	1	RHON	scalar	R4	0	5
ZYF	scalar	R4	0	2	RHOV	scalar	R4	0	6
U	scalar	R4	0	3	NDIM6	scalar	I4	I	7
V	scalar	R4	0	4	NPEMAX	scalar	I4	I	8

name of the common block: RECS									
name	dimension	type	mode	word	name	dimension	type	mode	word
MAT	scalar	I4	-	1	L1	scalar	I4	I	6
MF	scalar	I4	-	2	L2	scalar	I4	I	7
MT	scalar	I4	-	3	N1	scalar	I4	IO	8
C1	scalar	R4	I	4	N2	scalar	I4	IO	9
C2	scalar	R4	I	5					

Called subprograms:

ANNIO1

CHGINT

ERF

ERROR

GENER (entry of SCALE)

GNRL

INITSC (entry of SCALE)

PRIZEI

RREC

### 2.3.5.26 Subroutine GNRL

Problem solved:

Calculation of the fluctuation integrals for capture, fission, and elastic scattering. The subroutine is called by ANNICK.

Definition:

SUBROUTINE GNRL(GALPHA,GBETA,GAMMA,MU,NU,LAMBDA,S,DF,ID)

Explanation of the arguments:

name	mode	type	dimension	remarks
GALPHA	I	R4	scalar	$\bar{\Gamma}_n$
GBETA	I	R4	scalar	$\bar{\Gamma}_f$
GAMMA	I	R4	scalar	$\bar{\Gamma}_\gamma$
MU	I	I4	scalar	number of degrees of freedom used in the neutron width distribution

name	mode	type	dimension	remarks
NU	I	I4	scalar	number of degrees of freedom used in the fission width distribution
LAMBDA	I	I4	scalar	number of degrees of freedom used in the competitive width distribution
S	0	R4	scalar	fluctuation integral value for elastic scattering (ID=1) capture (ID=2) fission (ID=3)
DF	I	R4	scalar	$\bar{\Gamma}_x$
ID	I	I4	scalar	see description of S

Used common block:

name of the common block: XBAR				
name	dimension	type	mode	word
XX	(4,10)	R4	I	1

Called subprograms:

None

### 2.3.5.27 Subroutine ERF

Problem solved:

Part of the calculation of the Dresner factors  $S_f$ ,  $S_\gamma$ ,  $R_f$ ,  $R_\gamma$  (see /4/, /8/ and part II.). Subdivides the integration range in parts of 0.05 and stops the integration process if a integral over such an interval is less than  $10^{-6}$ .

Definition:

SUBROUTINE ERF(EER)

Explanation of the argument:

name	mode	type	dimension	remarks
ERR	0	R4	scalar	result of the integration

Used common block:

name of the common block: FLCT									
name	dimension	type	mode	word	name	dimension	type	mode	word
ZYN	scalar	R4	I	1	V	scalar	R4	-	4
ZYF	scalar	R4	-	2	RHON	scalar	R4	I	5
U	scalar	R4	-	3	RHOF	scalar	R4	-	6

Called subprogram:

GAUSS

2.3.5.28 Subroutine GAUSS

Problem solved:

Numerical integration with a Gauss method with five points (see /7/).

Definition:

SUBROUTINE GAUSS(HU,HO,B)

Explanation of the arguments:

name	mode	type	dimension	remarks
HU	I	R4	scalar	lower integration limit
HO	I	R4	scalar	upper integration limit
B	O	R4	scalar	result of the integration

Used common blocks:

None

Called subprogram:

BINT



2.3.5.29 Function BINT

Problem solved:

Used for the calculation of the Dresner factors  $S_f$ ,  $S_\gamma$ ,  $R_f$ ,  $R_\gamma$  (see /4/, /8/ and II.). Calculation of the function to be integrated.

Definition

FUNCTION BINT(X)

Explanation of the argument:

name	mode	type	dimension	remarks
X	I	R4	scalar	value for which the functional value is to be calculated

Used common block:

name of the common block: FLCT									
name	dimension	type	mode	word	name	dimension	type	mode	word
ZYN	scalar	R4	I	1	V	scalar	R4	I	4
ZYF	scalar	R4	I	2	RHON	scalar	R4	I	5
U	scalar	R4	I	3	RHOF	scalar	R4	I	6

Called subprograms:

None

2.3.5.30 Subroutine ANNI01

Problem solved:

Optionally printing of the unresolved resonance parameters in KEDAK representation and of the pointwise cross sections. Storing of the unresolved resonance parameter blocks on logical unit 11 and of the pointwise cross sections on logical unit 12.

Definition:

SUBROUTINE ANNIO1 (NE, SIGT, SIGG, SIGS, SIGF, NIS, MIS, E, NB, BK, ISTGF, NB2,  
BK2, LFCHG, NCU, ESC, IDIMBK, INTS, NBT, JNT, Y, EL, EH, LFW)

Explanation of the arguments:

name	mode	type	dimension	remarks
NE	I	I4	scalar	number of the energy points of the pointwise cross sections in the unresolved resonance region
SIGT	L	R4	(NPEMAX)	see 2.3.5.25 argument list
SIGG	I	R4	(NPEMAX)	" "
SIGS	I	R4	(NPEMAX)	" "
SIGF	I	R4	(NPEMAX)	" "
NIS	I	I4	scalar	number of isotopes
MIS	I	I4	scalar	pointer to the actual isotope, if NIS is greater than 1
E	I	R4	(NPEMAX)	see 2.3.5.25 argument list
NB	I	I4	scalar	number of values in the array BK
BK	I	R4	(NBKX)	unresolved resonance parameters
ISTGF	I	I4	scalar	= 0 no fission widths given = 1 fission widths given
NB2	I	I4	scalar	number of values in the array BK2
BK2	I	R4	(N2X)	data for KEDAK-type STGF(I) and pointwise cross sections (0). (BK2 and X have the same starting address.)
LFCHG	I	I4	scalar	= 0 all parameters energy independent = 1 only fission-widths are energy dependent = 2 all parameters are energy dependent
NCU	I	I4	scalar	number of values in the array ESC
ESC	I	R4	(NPEMAX)	see 2.3.5.25 argument list
IDIMBK	I	I4	scalar	IDIMBK = NBKX
INTS	I	I4	(18)	interpolation codes for the unresolved parameters

name	mode	type	dimension	remarks
NBT	L	I4	(N1X)	interpolation ranges
JNT	L	I4	(N1X)	interpolation codes
Y	L	R4	(N2X)	pointwise cross section values
EL	I	R4	scalar	see 2.3.5.25 argument list
EH	I	R4	scalar	" "
LFW	I	I4	scalar	" "

Used common blocks:

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NO	scalar	I4	I	1	NTYP	scalar	I4	-	84
NI	scalar	I4	-	2	NAMISO	(10)	R8	I	85
NT	scalar	I4	-	3	MATISO	(10)	I4	I	105
MODE	scalar	I4	-	4	NR2	(11)	I4	-	115
LEDIT1	scalar	I4	-	5	NBX	scalar	I4	-	126
LEDIT2	scalar	I4	I	6	N1X	scalar	I4	-	127
LEDITS	scalar	I4	I	7	N2X	scalar	I4	-	128
NR1	(75)	I4	-	8	N	scalar	I4	0	129
MATKDK	scalar	I4	I	83	NBKX	scalar	I4	I	130

name of the common block: RECS									
name	dimension	type	mode	word	name	dimension	type	mode	word
MAT	scalar	I4	-	1	L1	scalar	I4	-	6
MF	scalar	I4	-	2	L2	scalar	I4	-	7
MT	scalar	I4	-	3	N1	scalar	I4	0	8
C1	scalar	R4	-	4	N2	scalar	I4	0	9
C2	scalar	R4	-	5					

Called subprograms:

A8FORM  
ERROR  
LUCY  
MAGGY  
TERP1  
WDA2 (entry of WDA)  
WRECS2 (entry of WRECS)  
XTEXT

2.3.5.31 Subroutine MAGGY

Problem solved:

The contents of the arrays XD and YD are stored in the arrays X and Y.

Definition:

SUBROUTINE MAGGY(XD,YD,N,X,Y)

Explanation of the arguments:

name	mode	type	dimension	remarks
XD	I	R4	(N)	input array
YD	I	R4	(N)	" "
N	I	I4	scalar	number of data in the arrays
X	O	R4	(N)	output array
Y	O	R4	(N)	" "

Used common blocks:

None

Called subprograms:

None

2.3.5.32 Subroutine DENISE

Problem solved:

The pointwise cross section data calculated out of resolved and unresolved resonance parameters are fetched from logical unit 12. The data for different isotopes - if existant - are added and stored on logical unit 12. The total cross section in the resolved and unresolved resonance region is calculated and stored on logical unit 12. Optional printout of the pointwise cross section is done.

Definition:

SUBROUTINE DENISE(MATP,LRP,LISRES,LISUNR,NIS,LFW,LNU,BK,NBT,JNT,X,Y,B,A)

Explanation of the arguments:

name	mode	type	dimension	remarks
MATP	I	I4	scalar	ENDF material number
LRP	-	I4	scalar	not used in DENISE
LISRES	I	I4	scalar	= 1 no resolved pointwise cross sections - calculated out of parameters - available † 1 resolved pointwise cross sections - calculated out of parameters - are available
LISUNR	I	I4	scalar	= 1 no unresolved pointwise cross sections - calculated out of parameters - available † 1 unresolved pointwise cross sections - calculated out of parameters - are available
NIS	I	I4	scalar	number of isotopes
LFW	I	I4	scalar	= 0 no pointwise fission cross sections - calculated out of parameters - available = 1 pointwise fission cross sections - calculated out of parameters - are available in the unresolved resonance region
LNU	I	I4	scalar	see 2.3.5.10 argument list
BK	P	R4	(NBKX)	see 2.3.5.36 argument list

name	mode	type	dimension	remarks
NBT	L	I4	(N1X)	interpolation ranges
JNT	L	I4	(N1X)	interpolation codes
X	L	R4	(N2X)	(mainly energy values for cross sections)
Y	L	R4	(N2X)	(mainly cross section values)
B	L	R4	(NBX)	see subroutines called by DENISE, in which B is used
A	L	R4	(IADIM)	see subroutines called by DENISE, in which A is used

Used common blocks:

name of the common block: DENS									
name	dimension	type	mode	word	name	dimension	type	mode	word
NR3	(402)	I4	-	1	MX	scalar	I4	0	403

name of the common block: INOUT									
NO	scalar	I4	I	1	NBKX	scalar	I4	I	130
NR1	(5)	I4	-	2	NR4	(10)	I4	-	131
LEDITS	scalar	I4	I	7	EPS	scalar	R4	I	141
NR2	(122)	I4	-	8					

Called subprograms:

DELETE  
 IFETCH (entry of FETCH)  
 IFPDS (entry of FPDS)  
 IICOMB (entry of COMB)  
 IIPDS (entry of IPDS)  
 IRRECS (entry of RRECS)  
 ISTORE (entry of STORE)  
 MARION  
 MYRIAM

SOPHIE  
 WREC  
 WRECS2 (entry of WRECS)  
 XTEXT

2.3.5.33 Subroutine MYRIAM

Problem solved:

The pointwise cross sections calculated out of resolved resp. unresolved resonance parameters are added up for the different isotopes with respect to the abundance. The result is stored in the arrays X and Y.

Definition:

SUBROUTINE MYRIAM(NIS,LISRES,LISUNR,NAME,EL1,EH1,EPS,ITEST,Y)

Explanation of the arguments:

name	mode	type	dimension	remarks
NIS	I	I4	scalar	number of isotopes
LISRES	I	I4	scalar	only for LISRES=0 adding up is done in the resolved resonance region
LISUNR	I	I4	scalar	only for LISUNR=0 adding up is done in the unresolved resonance region
NAME	I	I4	scalar	type of the cross section (SCAT,CAPT,FISS)
EL1	0	R4	scalar	lower energy limit for the resolved resp. the unresolved resonance region
EH1	0	R4	scalar	upper energy limit for the resolved resp. the unresolved resonance region
EPS	I	R4	scalar	see 2.3.5.66 argument list
ITEST	I	I4	scalar	† 1 resolved cross sections are processed = 1 unresolved cross sections are processed
Y	L	R4	(N2X)	cross section values for the different isotopes

Used common blocks:

None

Called subprograms:

ADD                    FETCH  
 BETSY                 RRECS  
 COMB                  STORE  
 ERROR

2.3.5.34 Subroutine BETSY

Problem solved:

The pointwise cross sections values are multiplied with the abundance associated to the processed isotope.

Definition:

SUBROUTINE BETSY(MIS,Y)

Explanation of the arguments:

name	mode	type	dimension	remarks
MIS	I	I4	scalar	pointer to the isotope, which is to be processed
Y	IO	R4	(N2X)	cross section values

Used common blocks:

name of the common block: RECS									
name	dimension	type	mode	word	name	dimension	type	mode	word
NR2	(8)	I4	-	1	N2	scalar	I4	I	9
name of the common block: INOUT									
NR1	(8)	I4	-	1	ABIS	(20)	R4	I	9

Called subprograms:

None



2.3.5.35 Subroutine MARION

Problem solved:

The total cross section in the resolved (NAME2='RES') or the unresolved (NAME2='STAT') resonance region is calculated by adding  $\sigma_e$ ,  $\sigma_\gamma$  and  $\sigma_f$  if existant.

Definition:

SUBROUTINE MARION(LFW,EL,EH,EPS,NAME2)

Explanation of the arguments:

name	mode	type	dimension	remarks
LFW	I	I4	scalar	$\sigma_f$ is only added if LFW=1
EL	I	R4	scalar	lower energy limit for the resolved resp. the unresolved resonance region
EH	I	R4	scalar	upper energy limit for the resolved resp. the unresolved resonance region
EPS	IO	R4	scalar	see 2.3.5.66 argument list Accuracy used for calculating $\sigma_t$
NAME2	I	I4	scalar	= 'RES' cross sections of the resolved resonance region are processed = 'STAT' cross sections of the unresolved resonance region are processed

Used common blocks:

name of the common block: INOUT				
name	dimension	type	mode	word
NO	scalar	I4	I	1

Called subprograms:

ADD

COMB

ERROR

FETCH

RRECS

STORE

### 2.3.5.36 Subroutine SOPHIE

Problem solved:

Processing of ENDF/B MF=3 data. First following ENDF/B data types from MF=3 with the MT-numbers listed below are searched, and read if found (the names in brackets are the corresponding KEDAK names):

1(SGT),2(SGN),3(SGX),4(SGI),16(SG2N),17(SG3N),18(SGF),22(SGIA),23(SGI3A),  
24(SG2NA),25(SG3NA),27(SGA),28(SGIP),29(SGI2A),51...90(SGIZ),91(SGIZC),  
102(SGG),103(SGP),104(SGD),105(SGH3),106(SGHE3),107(SGALP),108(SG2HE),  
201(SGTR),206(ETA),207(ALPHA),251(MUEL).

In the case of SGT,SGN,SGF and SGG the resolved and unresolved resonance part - if existant - is added to the pointwise cross sections. The pointwise cross sections are stored on logical unit number 12.

Optionally (only used for an old KEDAK version) for KEDAVE=0 a common energy grid for SGI, all SGIZ, and SGIZC is generated after changing the interpolation code of all this cross sections to linear-linear. All cross section values for the types SGI, all SGIZ, and SGIZC are calculated and stored at this energy values.

If no absorption cross section has been found on ENDF/B it is generated by adding all the following existing cross section types:

SGG,SGF,SGP,SGD,SGH3,SGHE3,SGALP,SG2HE.

If possible SGTR is calculated using the formulae  $SGTR=SGT-MUEL*SGN$ . If the  $\bar{v}$ -data, read in the subroutine INGRID, had been given in polynomial representation, they will be calculated pointwise and stored on logical unit number 12.

Definition:

SUBROUTINE SOPHIE(MATP,LISRES,LISUNR,LFW,LNU,EPS,BK,XIN,YIN,NCMAX,  
NBT,JNT,X,Y,B,TEMP)

Explanation of the arguments:

name	mode	type	dimension	remarks
MATP	I	I4	scalar	ENDF/B material number
LISRES	I	I4	scalar	flag for the existence of pointwise cross sections calculated out of resolved resonance parameters = 1 no data available
LISUNR	I	I4	scalar	flag for the existence of pointwise cross sections calculated out of un-resolved resonance parameters = 1 no data available
LFW	I	I4	scalar	flag for the existence of pointwise fission cross sections calculated out of resonance parameters = 1 data available
LNU	I	I4	scalar	for LNU=1 pointwise $\bar{v}$ -data are calculated out of the polynomial representation
EPS	I	R4	scalar	accuracy used in the subroutines COMB (see 2.3.5.66) and CHGINT (see 2.3.5.73)
BK	-	R4	(NBKX)	not used in SOPHIE
XIN	L	R4	(NCMAX)	used for the common energy grid for the total inelastic and the inelastic excitation cross sections (BK(1)...BK(NBKX/2))
YIN	L	R4	(NCMAX)	used for the cross section values of SGI,SGIZ and SGIZC calculated at the energy points, which are stored in XIN (BK(NBKX/2+1)...BK(NBKX))
NCMAX	I	I4	scalar	NCMAX = NBKX/2
NBT	L	I4	(NIX)	interpolation ranges
JNT	L	I4	(NIX)	interpolation codes
X	L	R4	(N2X)	energy values
Y	L	R4	(N2X)	cross section values
B	L	R4	(NBX)	polynomial coefficients for $\bar{v}$
TEMP	L	R4	(NBX)	(array B) auxiliary array used in GENER (see 2.3.5.65)

Used common blocks:

name of the common block: RECS									
name	dimension	type	mode	word	name	dimension	type	mode	word
MAT	scalar	I4	-	1	L1	scalar	I4	-	6
MF	scalar	I4	0	2	L2	scalar	I4	-	7
MT	scalar	I4	0	3	N1	scalar	I4	0	8
C1	scalar	R4	-	4	N2	scalar	I4	IO	9
C2	scalar	R4	I	5					

name of the common block: INOUT									
NO	scalar	I4	I	1	NOTTRL	(50)	I4	I	29
NI	scalar	I4	-	2	BIB12	scalar	I4	-	79
NT	scalar	I4	I	3	SALINE	scalar	L4	I	80
MODE	scalar	I4	I	4	NR2	(2)	I4	-	81
LEDIT1	scalar	I4	-	5	MATKDK	scalar	I4	-	83
LEDIT2	scalar	I4	-	6	NR3	(42)	I4	-	84
LEDITS	scalar	I4	I	7	NBX	scalar	I4	-	126
KEDAVE	scalar	I4	I	8	N1X	scalar	I4	I	127
NR1	(20)	I4	-	9	N2X	scalar	I4	I	128

name of the common block: TAB									
ENDTAB	(25)	I4	I	1	NCNAME	(39)	I4	-	316
KEDTAB	(50)	I4	I	26	NRIF3	scalar	I4	IO	355
KEDR	(240)	I4	-	76	NRIF4	scalar	I4	-	356

Called subprograms:

ADD  
CHGINT  
COMB  
ERROR  
GENER } (entries of SCALE)  
INITSC }  
ITAB1  
MULT  
PRIZEI  
RREC  
RRECS  
RRECS4 (entry of RRECS)  
SEARCH  
STORE  
SUB  
WREC  
WRECS2 (entry of WRECS)  
XTEXT

2.3.5.37 Subroutine SUZY

Problem solved:

The cross sections, which have been written by the subroutine SOPHIE (see 2.3.5.36) on logical unit number 12, are read. The interpolation is changed to linear-linear, points with zero cross section values are suppressed, except those points, which are preceding the threshold of a cross section. At the threshold also zeros may be generated. Points, which can be predicted by linear interpolation, are deleted. Cross section values with energy values, which fulfill the condition  $(E_2 - E_1) / E_2 < 5.0 * 10^{-6}$ , are replaced by mean values. Points with negative cross section values for the types  $\sigma_t$  or  $\sigma_e$  are deleted. The result is printed out and written together with KEDAK type names on logical unit number 11. Optionally it can also be written in ENDF/B-format on logical unit number 15.

Definition:

SUBROUTINE SUZY(LNU,ITAPE,BK,A,TAB,ITAB,NBT,JNT,X,Y,B)

Explanation of the arguments:

name	mode	type	dimension	remarks
LNU	-	I4	scalar	not used in SUZY
ITAPE	I	I4	scalar	control of writing on unit 10 = 1 the data produced in the subroutine are written on the external unit 10 ≠ 1 no writing on unit 10 is done
BK	L	R4	(NBKX)	energy grid and cross sections which will be written in WDA2 on an external unit
A	L	R4	(NBKX)	(same starting address as BK) same contents as BK
TAB	L	R4	(NBKX)	(same starting address as BK) used for building up the KEDAK further name address table for the excited in- elastic cross section
ITAB	L	I4	(NBKX)	(same starting address as BK) same contents as TAB
NBT	L	I4	(N1X)	interpolation ranges
JNT	L	I4	(N1X)	interpolation codes
X	L	R4	(N2X)	energy grids
Y	L	R4	(N2X)	cross section values
B	P	R4	(NBX)	address is passed to the subroutine RREC

Used common blocks:

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NO	scalar	I4	I	1	NR4	(42)	I4	-	84
NI	scalar	I4	-	2	NBX	scalar	I4	-	126
NT	scalar	I4	-	3	N1X	scalar	I4	I	127
MODE	scalar	I4	I	4	N2X	scalar	I4	I	128
NR1	(3)	I4	-	5	N	scalar	I4	0	129
KEDAVE	scalar	I4	I	8	NBKX	scalar	I4	I	130
NR2	(71)	I4	-	9	NR5	(10)	I4	-	131
SALINE	scalar	L4	I	80	EPS	scalar	R4	I	141
NR3	(2)	I4	-	81	SCREPS	scalar	R4	I	142
MATKDK	scalar	I4	I	83					

name of the common block: RECS									
MAT	scalar	I4	I	1	L1	scalar	I4	0	6
MF	scalar	I4	0	2	L2	scalar	I4	0	7
MT	scalar	I4	0	3	N1	scalar	I4	0	8
C1	scalar	R4	0	4	N2	scalar	I4	0	9
C2	scalar	R4	0	5	NS	scalar	I4	0	10

name of the common block: TAB									
ENDTAB	(25)	I4	I	1	NCNAME	(39)	I4	-	316
KEDTAB	(50)	I4	I	26	NRIF3	scalar	I4	I	355
KEDR	(240)	I4	-	76	NRIF4	scalar	I4	-	356

Called subprograms:

ADDPNT  
A8FORM  
CHGINT  
CROP  
DZMSCH  
ELIMZ  
ERROR  
IRRECS (entry of RRECS)  
PRIZEI  
RRECS  
SUPNEG  
WCONT  
WDA2 (entry of WDA)  
WREC  
XTEXT

2.3.5.38 Subroutine ELIMZ

Problem solved:

Points with  $y = 0.0$  are suppressed, except the point preceding the threshold of the cross section. (The interpolation scheme must be linear-linear.)

Definition:

SUBROUTINE ELIMZ(N1,N2,NBT,JNT,X,Y)

Explanation of the arguments:

name	mode	type	dimension	remarks
N1	I	I4	scalar	must have the value 1
N2	IO	I4	scalar	number of points in the arrays X and Y
NBT	IO	I4	(N1X)	interpolation ranges. NBT(1) must be equal N2
JNT	I	I4	(N1X)	interpolation codes. JNT(1) must be equal 2 or equal 1
X	IO	R4	(N2X)	energy grid
Y	IO	R4	(N2X)	cross section values



Used common blocks:

None

Called subprograms:

ERROR

### 2.3.5.39 Subroutine ADDPNT

Problem solved:

For threshold reactions two zero cross section values at the beginning of the energy range must exist. A threshold reaction is assumed, if the first energy point is greater than  $1.0 \cdot 10^{-5}$  eV and if the first cross section value is less or equal  $1.0 \cdot 10^{-20}$ . The subroutine generates such cross section values, if they don't exist. If the first cross section value is less or equal  $1.0 \cdot 10^{-20}$  it is set to zero. (Interpolation must be linear-linear.)

Definition:

SUBROUTINE ADDPNT(X,Y,NBT,JNT,MT)

Explanation of the arguments:

name	mode	type	dimension	remarks
X	IO	R4	(N2X)	energy grid
Y	IO	R4	(N2X)	cross section values
NBT	IO	I4	(N1X)	interpolation ranges. NBT(1) must be equal N2
JNT	I	I4	(N1X)	interpolation codes. JNT(1) must be less or equal 2
MT	I	I4	scalar	ENDF/B section number

Used common blocks:

name of the common block: RECS									
name	dimension	type	mode	word	name	dimension	type	mode	word
NR1	(7)	I4	-	1	N2	scalar	I4	I0	9
N1	scalar	I4	I	8					

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NO	scalar	I4	I	1	N2X	scalar	I4	I	128
NR2	(126)	I4	-	2					

Called subprograms:

ERROR

#### 2.3.5.40 Subroutine DZMSCH

Problem solved:

Cross section values at energies with 6 similar digits will be replaced by mean values:

$$X_{NEW} = 1/N*(X_1+X_2+\dots+X_N)$$

$$Y_{NEW} = 1/N*(Y_1+Y_2+\dots+Y_N)$$

The accuracy with which the replaced points could have been interpolated is determined and ordered in classes.

Definition:

SUBROUTINE DZMSCH(X,Y,NBT,JNT,NDZ,EPS,SCRCLA,NUMSCR)

Explanation of the arguments:

name	mode	type	dimension	remarks
X	IO	R4	(N2X)	energy grid
Y	IO	R4	(N2X)	cross section values
NBT	IO	I4	(N1X)	interpolation ranges. NBT(1) must be equal N2
JNT	I	I4	(N1X)	interpolation codes. JNT(1) must be less or equal 2
NDZ	0	I4	scalar	total number of scratched values
EPS	I	R4	scalar	used for the construction of the scratch classes
SCRCLA	0	R4	(10)	scratch classes. Accuracies, with which the replaced points could have been interpolated (see 2.2.7).
NUMSCR	0	I4	(10)	number of points in a specified scratch class

Used common block:

name of the common block: RECS									
name	dimension	type	mode	word	name	dimension	type	mode	word
NR1	(7)	I4	-	1	N2	scalar	I4	IO	9
N1	scalar	I4	I	8					

Called subprograms:

ERROR

2.3.5.41 Subroutine SUPNEG

Problem solved:

For the cross section types  $\sigma_t$  and  $\sigma_e$  negative cross section values (mainly in the resolved resonance region) are diagnosed and optionally eliminated.

Definition:

SUBROUTINE SUPNEG(MT,N1,N2,NBT,JNT,X,Y,)

Explanation of the arguments:

name	mode	type	dimension	remarks
MT	I	I4	scalar	ENDF/B section number. Must be equal 1 or 2
N1	I	I4	scalar	number of interpolation ranges. Must be equal 1, if points must be eliminated
N2	IO	I4	scalar	number of energy points resp. cross section values
NBT	IO	I4	(N1X)	interpolation ranges. NBT(1) must be equal N2, if points must be eliminated
JNT	I	I4	(N1X)	interpolation codes. JNT(1) must be less or equal 2, if points must be eliminated
X	IO	R4	(N2X)	energy grid
Y	IO	R4	(N2X)	cross section values

Used common block:

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NO	scalar	I4	I	1	ELINEG	scalar	I4	I	154
NR1	(152)	I4	-	2					

Called subprogram:

ERROR

2.3.5.42 Subroutine ISABEL

Problem solved:

Processing of ENDF/B MF=4 data (angular distributions of secondary neutrons).

Only elastic angular distributions are processed. If the distributions are given in tabulated form, only these data (SGNC or SGNL see /4/, /5/ or II) are written on logical unit number 11 for later insertion into KEDAK. If the data are given as Legendre polynomial coefficients, these coefficients are written as LEGNC or LEGNL (see /4/, /5/ or II) on logical unit number 11 together with the angular pointwise data calculated out of the coefficients (81 cosine values are used for each energy point).

Definition:

SUBROUTINE ISABEL(MATP,IF4,ITAB,TAB,CLEG,NBT,JNT,X,Y,B,A,ITAB2,TAB2)

Explanation of the arguments:

name	mode	type	dimension	remarks
MATP	I	I4	scalar	ENDF/B-number of the material to be converted
IF4	I	I4	scalar	see 2.2.6. If IF4=1 only ENDF/B-data of MF=4 are converted. Then ISABEL is the first subroutine writing on file 11. Then this file is initialized in this subroutine
ITAB	L	I4	(NBKX)	(array BK) used for storing Legendre polynomial coefficients in KEDAK representation
TAB	L	R4	(NBKX)	(array BK) see ITAB
CLEG	L	R4	(NBKX)	(array BK) see ITAB
NBT	L	I4	(NIX)	interpolation ranges
JNT	L	I4	(NIX)	interpolation codes
X	L	R4	(N2X)	cosine values
Y	L	R4	(N2X)	distribution probabilities
B	L	R4	(NBX)	Legendre coefficients
A	L	R4	(IADIM)	used for storing tabulated distribution probabilities in KEDAK representation

name	mode	type	dimension	remarks
ITAB2	L	I4	(IADIM)	(array A) see A
TAB2	L	R4	(IADIM)	(array A) see A

Used common blocks:

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NO	scalar	I4	I	1	NR2	(45)	I4	-	84
NI	scalar	I4	-	2	N	scalar	I4	L	129
NT	scalar	I4	I	3	NBKX	scalar	I4	I	130
MODE	scalar	I4	I	4	JX	scalar	I4	-	131
NR1	(78)	I4	-	5	IADIM	scalar	I4	I	132
MATKDK	scalar	I4	I	83					

name of the common block: RECS									
name	dimension	type	mode	word	name	dimension	type	mode	word
MAT	scalar	I4	-	1	L1	scalar	I4	I	6
MF	scalar	I4	-	2	L2	scalar	I4	I	7
MT	scalar	I4	-	3	N1	scalar	I4	IO	8
C1	scalar	R4	-	4	N2	scalar	I4	IO	9
C2	scalar	R4	I	5					

name of the common block: TAB									
name	dimension	type	mode	word	name	dimension	type	mode	word
ENDTAB	(25)	I4	-	1	NCNAME	(39)	I4	0	316
KEDTAB	(50)	I4	-	26	NRIF3	scalar	I4	-	355
KEDR	(240)	I4	I	76	NRIF4	scalar	I4	0	356

Called subprograms:

A8FORM  
ERROR  
ITAB1  
LEGP  
PRIZEI  
RREC  
SEARCH  
WREC

2.3.5.43 Function LEGP

Problem solved:

Calculation of Legendre polynomials.

Definition:

REAL FUNCTION LEGP(X,NL,CON)

Explanation of the arguments:

name	mode	type	dimension	remarks
X	I	R4	scalar	cosine value
NL	I	I4	scalar	order of the Legendre polynom.
CON	I	R4	(NL)	Legendre polynomial coefficients

Used common blocks:

None

Called subprograms:

None

2.3.5.44 Subroutine FILE5

Problem solved:

Processing of ENDF/B MF=5 data (energy distributions of secondary neutrons).

Following cross section types can be produced by the subroutine

KEDAK-names	ENDF section numbers
SED2N	MT = 16
SED3N	MT = 17
SEDF	MT = 18
SEDIC	MT = 91

These types can be produced, if the ENDF-representation is LF = 3 (K = 4 excitation of discrete levels), LF = 7 (K = 2 Maxwellian spectrum), LF = 9 (K = 1 evaporation spectrum) or LF = 10 (K = 3 Watt spectrum). (The values K refer to the KEDAK data types.) All interpolation is changed to linear-linear. The result of the subroutine is written on logical unit number 11 in a representation, which is used to built up the KEDAK-file.

Definition:

SUBROUTINE FILE5(MATP,IF5,BK,IBK,NBT,JNT,X,Y,B,TEMP)

Explanation of the arguments:

name	mode	type	dimension	remarks
MATP	I	I4	scalar	ENDF/B-number of the material, which is to be converted
IF5	-	I4	scalar	not used in FILE5
BK	L	R4	(NBKX)	used for building up the energy distributions in a KEDAK similar representation
IBK	L	I4	(NBKX)	(array BK) see BK
NBT	L	I4	(N1X)	interpolation ranges
JNT	L	I4	(N1X)	interpolation codes
X	L	R4	(N2X)	energy grids



name	mode	type	dimension	remarks
Y	L	R4	(N2X)	distributions or probabilities
B	L	R4	(NBX)	used for LF=10 (Watt spectrum) for a and b (see /2/ pages 10.3 and 10.8)
TEMP	P	R4	(NBX)	(array B) auxiliary array

Used common blocks:

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NO	scalar	I4	I	1	NBX	scalar	I4	-	126
NI	scalar	I4	-	2	N1X	scalar	I4	I	127
NT	scalar	I4	I	3	N2X	scalar	I4	I	128
MODE	scalar	I4	I	4	N	scalar	I4	0	129
NR1	(78)	I4	-	5	NBKX	scalar	I4	I	130
MATKDK	scalar	I4	I	83	NR3	(10)	I4	-	131
NR2	(42)	I4	-	84	EPS	scalar	R4	I	141

name of the common block: RECS									
MAT	scalar	I4	-	1	L1	scalar	I4	-	6
MF	scalar	I4	-	2	L2	scalar	I4	I	7
MT	scalar	I4	-	3	N1	scalar	I4	I	8
C1	scalar	R4	I	4	N2	scalar	I4	I	9
C2	scalar	R4	I	5					

name of the common block: TAB									
ENDTAB	(25)	I4	-	1	NCNAME	(39)	I4	0	316
KEDTAB	(50)	I4	-	26	NRIF3	scalar	I4	-	355
KEDR	(240)	I4	I	76	NRIF4	scalar	I4	I0	356

Called subprograms:

A8FORM

CHGINT

ERROR

GENER } (entries of SCALE)

INITSC }

ITAB1

PRIZEI

RREC

SEARCH

WDA2 (entry of WDA)

WREC

2.3.5.45 Subroutine PRIZEI

Problem solved:

Prints date, elapsed time and CPU-time.

Definition:

SUBROUTINE PRIZEI

Used common block:

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NO	scalar	I4	I	1	ANFZEI	scalar	R4	I	153
NR1	(151)	I4	-	2					

Called subprograms:

DATUM

ZEIT

2.3.5.46 Subroutine ERROR

Problem solved:

Prints the error number N, the quantities MAT, MF, and MT of the common RECS and the contents of NBT,JNT,X, and Y. Afterwards STOP1 is made.

Definition:

SUBROUTINE ERROR(N)

Explanation of the argument:

name	mode	type	dimension	remarks
N	I	I4	scalar	error number

Used common blocks:

name of the common block: RECS									
name	dimension	type	mode	word	name	dimension	type	mode	word
MAT	scalar	I4	I	1	MT	scalar	I4	I	3
MF	scalar	I4	I	2					

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NOUT	scalar	I4	I	1					

Called subprogram:

WREC

### 2.3.5.47 Subroutine XTEXT

Problem solved:

The value of I is assigned to J. The contents of J is normally text of the form 'ABCD'.

Definition:

SUBROUTINE XTEXT(I,J)

Explanation of the arguments:

name	mode	type	dimension	remarks
I	I	I4	scalar	see problem solved
J	O	I4	scalar	see problem solved

Used common blocks:

None

Called subprograms:

None

### 2.3.5.48 Subroutine SEARCH

Problem solved:

Search ENDF/B tape NT for a particular section denoted by MATP,MFP,MTP. Search is in forward direction only. Zero values of MATP,MFP,MTP are read as (any). Thus, MATP,MFP,0 means any section in material MATP, file MFP. 0,MFP,0 means any (next) section in file MFP for any (next) material.

Definition:

SUBROUTINE SEARCH(NT,MODE,MATP,MFP,MTP,LNT)

(The subroutine is taken out of /3/ part R)

Explanation of the arguments:

name	mode	type	dimension	remarks
NT	I	I4	scalar	logical ENDF/B tape number
MODE	I	I4	scalar	tape mode. May be 1,2,3. (see /2/ chapter 4.)
MATP	I	I4	scalar	desired material number or zero
MFP	I	I4	scalar	desired file number or zero
MTP	I	I4	scalar	desired section number or zero
LNT	0	I4	scalar	= 0 section located. MAT,MF,MT stored in common RECS. = 1 section not located, because end of tape encountered, or because desired record is located prior to the current position of the tape

Used common block:

name of the common block: RECS									
name	dimension	type	mode	word	name	dimension	type	mode	word
MAT	scalar	I4	0	1	MT	scalar	I4	0	3
MF	scalar	I4	0	2					

Called subprograms:

ERROR

RREC

#### 2.3.5.49 Subroutine RREC

Problem solved:

One record is read from the ENDF/B tape. The result is stored in the  
common RECS for CONT-records

or in the

common RECS and the array B for  
LIST-records

or in the

common RECS and the arrays NBT,JNT,  
X,Y, for TAB1-records

or in the

common RECS and the arrays NBT,JNT,  
for TAB2-records

or in the

common RECS and the array B for  
HOL-records

or in the

quantities MAT,JF,MT of the common RECS  
and the first 17 words of the array B  
for TPID-records.

For explanation of the record-types see /2/ chapter 4.

Definition:

SUBROUTINE RREC(JT,NT,MODE,T)

ENTRY IRREC(NBT,JNT,X,Y,B)

(The entry is only used for passing the addresses of the arrays to the sub-  
routine. It must be called before the first RREC-call.) (The subroutine has  
been taken with some changes out of /3/ part R.)

Explanation of the arguments:

name	mode	type	dimension	remarks
JT	I	I4	scalar	record type. May be 1,2,3,4,5,6. (see /2/)
NT	I	I4	scalar	logical ENDF/B-tape number
MODE	I	I4	scalar	ENDF/B-tape mode. May be 1,2,3 (see /2/)
T	I	R4	scalar	temperature (Kelvin). For T>0 all data will be interpolated for tempera- ture T, if there is a temperature dependance given in the record. For T<0 only the first record will be read
NBT	P <sup>+</sup>	I4	(N1X)	used to store retrieved data
JNT	P <sup>+</sup>	I4	(N1X)	" " " " "
X	P <sup>+</sup>	R4	(N2X)	" " " " "
Y	P <sup>+</sup>	R4	(N2X)	" " " " "
B	P <sup>+</sup>	R4	(NBX)	" " " " "

Used common blocks:

name of the common block: RECS									
name	dimension	type	mode	word	name	dimension	type	mode	word
MAT	scalar	I4	0	1	L1	scalar	I4	0	6
MF	scalar	I4	0	2	L2	scalar	I4	0	7
MT	scalar	I4	0	3	N1	scalar	I4	0	8
C1	scalar	R4	0	4	N2	scalar	I4	0	9
C2	scalar	R4	0	5					

<sup>+</sup>The addresses are passed to the subroutine in the entry. These arrays will be filled by RREC-calls with respect to the rules given in the section "problem solved".

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NR1	(125)	I4	-	1	N1X	scalar	I4	I	127
NBX	scalar	I4	I	126	N2X	scalar	I4	I	128

Called subprograms:

ERROR

TERP1

### 2.3.5.50 Subroutine WCONT

Problem solved:

Writes ENDF/B end-record on tape unit NT (see /2/ chapter 4.).

Definition:

SUBROUTINE WCONT(NT,MODE,MAT,MF,NS)

Explanation of the arguments:

name	mode	type	dimension	remarks
NT	I	I4	scalar	logical tape unit number
MODE	I	I4	scalar	ENDF/B-mode. May be 1,2,3. (see /2/ chapter 4.)
MAT	I	I4	scalar	ENDF/B mat number or zero or -1 (see /2/ chapter 4.)
MF	I	I4	scalar	ENDF/B file number or zero (see /2/ chapter 4.)
NS	I	I4	scalar	sequence number written in card columns 76-80

Used common blocks:

None

Called subprograms:

None



2.3.5.51 Subroutine WREC

Problem solved:

Writes one record in ENDF representation. The data in the common RECS together with the contents of the arrays NBT,JNT,X,Y, and B (depending on the record type) are written in the ENDF representation according to the given mode.

The subroutine has been taken out of /3/ chapter R (see R-34 to R-35 and R-63) but some modifications have been made before inserting it in BRIGITTE.

E.g.: the arrays NBT,KNT,X,Y, and B, which have been of fixed length and placed in the common RECS, are now of variable length, and are passed to the subroutine by an entry.

Definition:

```
SUBROUTINE WREC(JT,NT,MODE)
ENTRY IWREC(NBT,JNT,X,Y,B)
```

(A call of the entry only passes the addresses of the arrays to the subroutine)

Explanation of the arguments:

name	mode	type	dimension	remarks
JT	I	I4	scalar	record type. May be 1,2,3,4,5,6. (see /3/ chapter R-2.3)
NT	I	I4	scalar	logical output tape number. For $NT \leq 0$ writing is ignored
MODE	I	I4	scalar	arrangement of the data. May be 1,2,3,4 (see /3/ chapter R-2.2)
NBT	I	I4	(N1X)	data to be written according to the record type
JNT	I	I4	(N1X)	" "
X	I	R4	(N2X)	" "
Y	I	R4	(N2X)	" "
B	I	R4	(NBX)	" "

Used common block:

name of the common block: RECS									
name	dimension	type	mode	word	name	dimension	type	mode	word
MAT	scalar	I4	I	1	L1	scalar	I4	I	6
MF	scalar	I4	I	2	L2	scalar	I4	I	7
MT	scalar	I4	I	3	N1	scalar	I4	I	8
C1	scalar	R4	I	4	N2	scalar	I4	I	9
C2	scalar	R4	I	5					

Called subprograms:

ERROR  
PRCONT  
PRHOL  
PRLIST  
PRTAB1  
PRTAB2  
PRTPID  
PUCONT  
PUHOL  
PULIST  
PUTAB1  
PUTAB2  
PUTPID

2.3.5.52 Subroutine PRCONT

Problem solved:

Called by the subroutine WREC to do the expanded interpreted print of a CONT-record. The subroutine has been taken from /3/ part R with the modifications mentioned in 2.3.5.51 (Original routine see /3/ page R-66).

Definition:

SUBROUTINE PRCONT(NT)

Explanation of the argument:

name	mode	type	dimension	remarks
NT	I	I4	scalar	print output unit number

Used common block:

name of the common block: RECS									
name	dimension	type	mode	word	name	dimension	type	mode	word
MAT	scalar	I4	I	1	L1	scalar	I4	I	6
MF	scalar	I4	I	2	L2	scalar	I4	I	7
MT	scalar	I4	I	3	N1	scalar	I4	I	8
C1	scalar	R4	I	4	N2	scalar	I4	I	9
C2	scalar	R4	I	5	NS	scalar	I4	IO	10

Called subprogramms:

None

2.3.5.53 Subroutine PRLIST

Problem solved:

Called by the subroutine WREC to do the expanded interpreted print of a LIST-record. The subroutine has been taken from /3/ part R with the modifications mentioned in 2.3.5.51 (Original routine see /3/ page R-67).

Definition:

SUBROUTINE PRLIST(NT,B)

Explanation of the arguments:

name	mode	type	dimension	remarks
NT	I	I4	scalar	print output unit number
B	I	R4	(NBX)	data to be printed

Used common blocks:

See 2.3.5.52

Called subprograms:

None

2.3.5.54 Subroutine PRTAB1

Problem solved:

Called by the subroutine WREC to do the expanded interpreted print of a TAB1-record. The subroutine has been taken from /3/ part R with the modifications mentioned in 2.3.5.51 (Original routine see /3/ page R-68).

Definition:

SUBROUTINE PRTAB1(NT,NBT,JNT,X,Y)

Explanation of the arguments:

name	mode	type	dimension	remarks
NT	I	I4	scalar	print output unit number
NBT	I	I4	(N1X)	data to be printed
JNT	I	I4	(N1X)	" " " "
X	I	R4	(N2X)	" " " "
Y	I	R4	(N2X)	" " " "

Used common blocks:

See 2.3.5.52

Called subprograms:

None

2.3.5.55 Subroutine PRTAB2

Problem solved:

Called by the subroutine WREC to do the expanded interpreted print of a TAB2-record. The subroutine has been taken from /3/ part R with the modifications mentioned in 2.3.5.51 (Original routine see /3/ page R-70).

Definition:

SUBROUTINE PRTAB2(NT,NBT,JNT)

Explanation of the arguments:

name	mode	type	dimension	remarks
NT	I	I4	scalar	print output unit number
NBT	I	I4	(N1X)	data to be printed
JNT	I	I4	(N1X)	" " " "

Used common blocks:

See 2.3.5.52

Called subprograms:

None

2.3.5.56 Subroutine PRHOL

Problem solved:

Called by the subroutine WREC to do the expanded interpreted print of a HOL-record. The subroutine has been taken from /3/ part R with the modifications mentioned in 2.3.5.51 (Original routine see /3/ page R-65).

Definition:

SUBROUTINE PRHOL(NT,B)

Explanation of the arguments:

name	mode	type	dimension	remarks
NT	I	I4	scalar	print output unit number
B	I	R4	(NBX)	data to be printed

Used common blocks:

See 2.3.5.52

Called subprograms:

None

2.3.5.57 Subroutine P RTPID

Problem solved:

Called by the subroutine WREC to do the expanded interpreted print of a TPID-record. The subroutine has been taken from /3/ part R with the modifications mentioned in 2.3.5.51 (Original routine see /3/ page R-65).

Definition:

SUBROUTINE P RTPID(NT,B)

Explanation of the arguments:

name	mode	type	dimension	remarks
NT	I	I4	scalar	print output unit number
B	I	R4	(NBX)	data to be printed

Used common block:

name of the common block: RECS									
name	dimension	type	mode	word	name	dimension	type	mode	word
MAT	scalar	I4	I	1	NS	scalar	I4	IO	10
NR1	(8)	I4	-	2					

Called subprograms:

None

2.3.5.58 Subroutine PUCONT

Problem solved:

Called by the subroutine WREC to produce in a BCD card image tape a CONT-record. The subroutine has been taken from /3/ part R with the modifications mentioned in 2.3.5.51 (Original routine see /3/ page R-67).

Definition:

SUBROUTINE PUCONT(NT)

Explanation of the argument:

name	mode	type	dimension	remarks
NT	I	I4	scalar	logical tape unit number

Used common block:

See 2.3.5.52

Called subprogram:

CXFP

2.3.5.59 Subroutine PULIST

Problem solved:

Called by the subroutine WREC to produce in a BCD card image tape a LIST-record. The subroutine has been taken from /3/ part R with the modifications mentioned in 2.3.5.51 (Original routine see /3/ page R-67).

Definition:

SUBROUTINE PULIST(NT,B)

Explanation of the arguments:

name	mode	type	dimension	remarks
NT	I	I4	scalar	logical tape unit number
B	I	R4	(NBX)	data to be written on NT



Used common block:

See 2.3.5.52

Called subprogram :

CXFP

2.3.5.60 Subroutine PUTABI

Problem solved:

Called by the subroutine WREC to produce in a BCD card image tape a TABI-record. The subroutine has been taken from /3/ part R with the modifications mentioned in 2.3.5.51 (Original routine see /3/ page R-69).

Definition:

SUBROUTINE PUTABI (NT,NBT,JNT,X,Y)

Explanation of the arguments:

name	mode	type	dimension	remarks
NT	I	I4	scalar	logical tape unit number
NBT	I	I4	(N1X)	data to be written on tape NT
JNT	I	I4	(N1X)	" "
X	I	R4	(N2X)	" "
Y	I	R4	(N2X)	" "

Used common block:

See 2.3.5.52

Called subprogram:

CXFP

### 2.3.5.61 Subroutine PUTAB2

Problem solved:

Called by the subroutine WREC to produce in a BCD card image tape a TAB2-record. The subroutine has been taken from /3/ part R with the modifications mentioned in 2.3.5.51 (Original routine see /3/ page R-71).

Definition:

SUBROUTINE PUTAB2(NT,NBT,JNT)

Explanation of the arguments:

name	mode	type	dimension	remarks
NT	I	I4	scalar	logical tape unit number
NBT	I	I4	(N1X)	data to be written on tape NT
JNT	I	I4	(N1X)	" "

Used common block:

See 2.3.5.52

Called subprogram:

CXFP

### 2.3.5.62 Subroutine PUHOL

Problem solved:

Called by the subroutine WREC to produce in a BCD card image tape a HOL-record. The subroutine has been taken from /3/ part R with the modifications mentioned in 2.3.5.51 (Original routine see /3/ page R-65).

Definition:

SUBROUTINE PUHOL(NT,B)

Explanation of the arguments:

name	mode	type	dimension	remarks
NT	I	I4	scalar	logical tape unit number
B	I	R4	(NBX)	data to be written on tape NT

Used common block:

See 2.3.5.52

Called subprogram:

CXFP

2.3.5.63 Subroutine PUTPID

Problem solved:

Called by the subroutine WREC to produce in a BCD card image tape a TPID-record. The subroutine has been taken from /3/ part R with the modifications mentioned in 2.3.5.51 (Original routine see /3/ page R-65).

Definition:

SUBROUTINE PUTPID(NT,B)

Explanation of the arguments:

name	mode	type	dimension	remarks
NT	I	I4	scalar	logical tape unit number
B	I	R4	(NBX)	data to be written on tape NT

Used common block:

name of the common block: RECS									
name	dimension	type	mode	word	name	dimension	type	mode	word
MAT	scalar	I4	I	1	NS	scalar	I4	I0	10
NR1	(8)	I4	-	2					

Called subprograms:

None

2.3.5.64 Subroutine CXFP

Problem solved:

Convert X for punching. X is represented as  $F*10.0^{**SN}$  with

F  $0.999995 \leq F \leq 9.999995$

S sign either (EBCDIC) + or - .

N exponent

(see /3/ page R-66).

Definition:

SUBROUTINE CXFP(X,F,S,N)

Explanation of the arguments:

name	mode	type	dimension	remarks
X	I	R4	scalar	number to be converted
F	O	R4	scalar	see problem solved
S	O	R4	scalar	" " "
N	O	I4	scalar	" " "

Used common blocks:

None

Called subprograms:

None

2.3.5.65 Subroutine SCALE

Problem solved:

Two or more energy scales are combined. If the combined scale is shorter than NCMAX it is kept in main storage, otherwise datasets NIN, NOUT are used (unit numbers 3 and 4). An energy point XNEW is added to the combined scale if no old points XOLD exists with

$$XOLD \in [XNEW \cdot 0.99999, XNEW \cdot 1.00001]$$

Definition:

SUBROUTINE SCALE (This entry point is never used.)

ENTRY INITSC(NCMAX) (Initialisation. Must be called before the first GENER-call and for deletion of a generated scale.)

ENTRY GENER(MAX,E,TEMP,EC,ND,NCU) (This entry creates the scale.)

Explanation of the arguments:

name	mode	type	dimension	remarks
NCMAX	I	I4	scalar	maximal dimension of array E.
MAX	I	I4	scalar	number of data in array E.
E	I	R4	(MAX)	energy points to be inserted in array EC.
TEMP	L	R4	(NCMAX)	auxiliary array.
EC	IO	R4	(NCMAX)	complete energy scale.
ND	0	I4	scalar	0 or number of the unit where the complete scale can be found.
NCU	0	I4	scalar	number of energy points in array EC.

Used common blocks:

None.

Called subprograms:

None.

### 2.3.5.66 Subroutine COMB

Problem solved:

Combine two TAB1 functions (see /3/ chapters R-2.3.3, R-9.2 and page R-87 for explanation).

Two TAB1 functions stored in dense storage (see 2.3.5.80 to 2.3.5.85) are combined and the result is stored in the arrays NBT, JNT, X and Y. Sufficient energy points are generated in order to guarantee the interpolation of the produced TAB1 record accurate to EPS. If FA(X) is the first TAB1 function and FB(X) is the second TAB1 function, then the resulting function FC(X) is defined by

$$FC(X) = FA(X) \text{ O } FB(X)$$

O means any function OPER, which address had been passed to the subroutine COMB. This function must be defined by the statement

```
REAL FUNCTION oper*4(YA,YB,CON)
```

where oper is to be replaced by a name. YA and YB mean the input values. CON is an input array of the length 20, which can be used by the function. Examples for such functions can be found in 2.3.5.68 to 2.3.5.70. The subroutine has been taken from /3/ part R with the changes mentioned in 2.3.5.51.

Definition:

```
SUBROUTINE COMB(OPER,CON,MA,MB,XL,XH,EPS,LOF,LNT)
```

```
ENTRY IICOMB(NBT,JNT,X,Y) (This entry passes the addresses of the  
arrays NBT, JNT, X and Y to the subroutine.  
It must be called before the first COMB-  
call and if the addresses are destroyed in  
the subroutine by using an overlay-structure.)
```

Explanation of the arguments:

name	mode	type	dimension	remarks
OPER	-	F4	scalar	name of a function, which combines the two TAB1 records (see problem solved)
CON	I	R4	(20)	array passed to the function OPER
MA	I	I4	scalar	identification of first TAB1-record in dense storage (see 2.3.5.80 to 2.3.5.85)
MB	I	I4	scalar	identification of second TAB1-record in dense storage (see 2.3.5.80 to 2.3.5.85)
XL	I	R4	scalar	lower limit of energy array X
XH	I	R4	scalar	upper limit of energy array X
EPS	I	R4	scalar	relative error criterion for generation of the energy grid of the combined TAB1 function
LOF	0	I4	scalar	overflow indicator. Normally 0. If 1, one or some of the arrays NBT, JNT, X or Y are too small
LNT	0	I4	scalar	normal exit = 0 = 1 record MA or MB not in dense storage
NBT	0	I4	(N1X)	produced interpolation ranges
JNT	0	I4	(N1X)	produced interpolation codes
X	0	R4	(N2X)	produced energy grid
Y	0	R4	(N2X)	produced Y (e.g. cross section) values

Used common blocks:

name of the common block: RECS									
name	dimension	type	mode	word	name	dimension	type	mode	word
NR1	(7)	I4	-	1	N2	scalar	I4	0	9
N1	scalar	I4	0	8					

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NR2	(141)	I4	-	1	SCREPS	scalar	R4	I	142

Called subprograms:

COMBP

CROP

ERROR

FPDS

IPDS

LRIDS

OPER (The function name is defined at the subroutine call statement)



2.3.5.67 Subroutine COMBP

Problem solved:

Two functions, given in the interval  $[x_1, x_2]$  by values at the beginning and end point and by interpolation codes are combined by using the function OPER.

Sufficient points in the interval  $[x_1, x_2]$  are created in order to guarantee the interpolation of the produced function accurate to EPS. The produced function is added to the function given in the arrays NBT, JNT, X and Y.

The subroutine has been taken from /3/ page R-85 with the following changes:

- (1) NBT, JNT, X and Y are now passed to the subroutine by an entry-call instead of a common. The length of these arrays now depends only from the available main storage region.
- (2) If  $|x_1 - x_2| / |x_1| \leq 10^{-6}$  no intermediate values are generated. The interpolation scheme of the combined function will be linear-linear.
- (3) In the case of (2) or a discontinuity a jump to RETURN after adding the values to NBT, JNT, X and Y has been built in.
- (4) The interpolation scheme of the produced function can be chosen by the program or can be always linear-linear.
- (5) To avoid frequent switching of the interpolation code, following method is used:  
Let EPS4 be the interpolation code of a previous interval and EPS5 the best interpolation code for the processed interval. If  $EPS4 \leq FAICOD \cdot EPS5$  the code of the previous interval EPS4 is used. Otherwise EPS5 is used. FAICOD is passed to the program by a common and can be set in the input.
- (6) The test on the elimination of points, which can be predicted by interpolation, has been replaced by a CROP-call in the

subroutine COMB (see 2.3.5.66 and 2.3.5.75). The old method had the incorrect effect to smooth faintly varying functions.

Definition:

SUBROUTINE COMBP (X1P, YA1P, YB1P, X2P, YA2P, YB2P, IAP, IBP, OPER, CON, EPSP, LOF, NBT, JNT, X, Y)

Explanation of the arguments:

name	mode	type	dimension	remarks
X1P	I	R4	scalar	lower limit of the interval
YA1P	I	R4	scalar	value of function A at X1P
YB1P	I	R4	scalar	value of function B at point X1P
X2P	I	R4	scalar	upper limit of the interval
YA2P	I	R4	scalar	value of the function A at X2P
YB2P	I	R4	scalar	value of the function B at X2P
IAP	I	I4	scalar	interpolation code for the function A
IBP	I	I4	scalar	interpolation code for the function B
OPER	-	F4	scalar	name of the function, which combines the functions A and B
CON	I	R4	(20)	array passed to the function OPER
EPSP	I	R4	scalar	relative error criterion for the generation of the energy grid of the combined function
LOF	0	I4	scalar	overflow indicator. Normally 0. If 1, one or some of the arrays NBT, JNT, X or Y have been too small
NBT	0	I4	(N1X)	interpolation ranges
JNT	0	I4	(N1X)	interpolation codes
X	0	R4	(N2X)	energy grid
Y	0	R4	(N2X)	Y (e.g. cross section) values

Used common blocks:

name of the common block: RECS									
name	dimension	type	mode	word	name	dimension	type	mode	word
MAT	scalar	I4	-	1	L1	scalar	I4	-	6
MF	"	I4	-	2	L2	"	I4	-	7
MT	"	I4	-	3	N1	"	I4	IO	8
C1	"	R4	-	4	N2	"	I4	IO	9
C2	"	R4	-	5					

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NR1	(124)	I4	-	1	N2X	scalar	I4	I	128
ICOMB	scalar	I4	I	125	NR2	(22)	I4	-	129
NBX	"	I4	-	126	FAICOD	scalar	R4	I	151
NIX	"	I4	I	127					

Called subprograms:

OPER (This name is defined at the subroutine  
call statement.)

TERP1

2.3.5.68 Function ADD

Problem solved:

The quantities YA and YB are added and the result is assigned to the function name (ADD=YA+YB).

Definition:

FUNCTION ADD(YA,YB,CON)

Explanation of the arguments:

name	mode	type	dimension	remarks
YA	I	R4	scalar	first operand
YB	I	R4	"	second operand
CON	-	R4	(20)	not used in ADD

Used common blocks:

None

Called subprograms:

None

2.3.5.69 Function SUB

Problem solved:

The quantities YA and YB are subtracted and the result is assigned to the function name (SUB=YA-YB).

Definition:

FUNCTION SUB(YA,YB,CON)

Explanation of the arguments:

name	mode	type	dimension	remarks
YA	I	R4	scalar	first operand
YB	I	R4	"	second operand
CON	-	R4	(20)	not used in SUB

Used common blocks:

None

Called subprograms:

None

2.3.5.70 Function MULT

Problem solved:

The quantities YA and YB are multiplied and the result is assigned to the function name (MULT=YA\*YB).

Definition:

```
REAL FUNCTION MULT(YA,YB,CON)
```

Explanation of the arguments:

name	mode	type	dimension	remarks
YA	I	R4	scalar	first operand
YB	I	R4	"	second operand
CON	-	R4	(20)	not used in MULT

Used common blocks:

None

Called subprograms:

None

2.3.5.71 Subroutine ITAB1

Problem solved:

Given a TAB1 function by the arrays NBT, JNT, X and Y (see /3/ chapter R-2.3.3). The subroutine interpolates the value YP at the x-value XP. If XP is outside the range of the TAB1 function, YP is set to zero.

Definition:

SUBROUTINE ITAB1(XP,YP)

ENTRY EITAB1(NBT,JNT,X,Y)

(An entry call must be made before the first ITAB1-call. By this entry call the addresses of NBT, JNT, X and Y are passed to the subroutine.)

Explanation of the arguments:

name	mode	type	dimension	remarks
XP	I	R4	scalar	x-value
YP	O	R4	scalar	desired y-value for x=XP
NBT	I	I4	(N1X)	interpolation ranges of the given TAB1-record
JNT	I	I4	(N1X)	interpolation codes of the given TAB1-record
X	I	R4	(N2X)	energy grid of the given TAB1-record
Y	I	R4	(N2X)	functional values of the given TAB1-record

Used common block:

name of the common block: RECS									
name	dimension	type	mode	word	name	dimension	type	mode	word
MAT	scalar	I4	-	1	L1	scalar	I4	-	6
MF	"	I4	-	2	L2	"	I4	-	7
MT	"	I4	-	3	N1	"	I4	-	8
C1	"	R4	-	4	N2	"	I4	I	9
C2	"	R4	-	5					

Called subprograms:

ERROR

TERP1



2.3.5.72 Subroutine TERP1

Problem solved:

Interpolate one point. An interpolated value Y at the desired X is computed by using the end points of the line (X1,Y1) and (X2,Y2) and the interpolation code I.

If X1 and/or X2 are negative or zero and interpolation code I=3 or 5 is desired, it is changed to I=2 or 4. If Y1 and/or Y2 are negative or zero and interpolation code I=4 or 5 is desired, it is changed to I=2 or 3. For negative or zero X and I=3 or 5 an error stop occurs.

The program given in /3/ pages R-30 resp. R-59 has been used after doing some changes. One of these changes is given in the following: The interpolation formulae for I=3 (y linear in log(x)) and I=5 (log(4) linear in log (x)) are

$$y_p = y_a + \log\left(\frac{x_p}{x_a}\right) \cdot \frac{y_b - y_a}{\log\left(\frac{x_b}{x_a}\right)} \quad (I=3)$$

$$y_p = y_a \cdot \text{EXP}\left(\log\left(\frac{x_p}{x_a}\right) \cdot \frac{\log\left(\frac{y_b}{y_a}\right)}{\log\left(\frac{x_b}{x_a}\right)}\right) \quad (I=5)$$

To avoid numerical difficulties,  $\log\left(\frac{x_p}{x_a}\right)$  and  $\log\left(\frac{x_b}{x_a}\right)$  is expanded in a series, if the quotients  $\frac{x_p}{x_a}$  and  $\frac{x_b}{x_a}$  have values near by one.

Definition:

SUBROUTINE TERP1(X1,Y1,X2,Y2,X,Y,I)

Explanation of the arguments:

name	mode	type	dimension	remarks
X1	I	R4	scalar	lower end of the interpolation interval
Y1	I	R4	scalar	functional value at the lower end of the interpolation interval
X2	I	R4	scalar	upper end of the interpolation interval
Y2	I	R4	scalar	functional value at the upper end of the interpolation interval
X	I	R4	scalar	point for which Y is to be calculated
Y	O	R4	scalar	interpolated value
I	I	I4	scalar	interpolation code in the interval $[X1, X2]$ . May be 1,2,3,4,5

Used common blocks:

None

Called subprogram:

ERROR

2.3.5.73 Subroutine CHGINT

Problem solved:

The interpolation code of the function given in the TAB1-record (see /3/ chapter R-2.3.3) is changed to linear-linear. Discontinuities are eliminated by small changes in the x (energy)-values.

Definition:

SUBROUTINE CHGINT(EPS,N1,N2,NBT,JNT,X,Y,N1X,N2X,\*)

Explanation of the arguments:

name	mode	type	dimension	remarks
EPS	I	R4	scalar	passed to subroutine DANI (see 2.3.5.74)
N1	IO	I4	scalar	number of interpolation ranges (array NBT) and interpolation codes (array JNT)
N2	IO	I4	scalar	number of energy points (array X) and functional values (array Y)
NBT	IO	I4	(N1X)	interpolation ranges
JNT	IO	I4	(N1X)	interpolation codes
X	IO	R4	(N2X)	energy grid
Y	IO	R4	(N2X)	functional values (e.g. cross sections)
N1X	I	I4	scalar	maximal dimension of the arrays NBT and JNT
N2X	I	I4	scalar	maximal dimension of the arrays X and Y
*	-	-	-	Return 1 is made, if the dimension of the arrays X and Y is too small

Used common block:

name of the common block: RE									
name	dimension	type	mode	word	name	dimension	type	mode	word
XR	(300)	R4	I	1	YR	(300)	R4	I	301

Called subprograms:

DANI  
ERROR

2.3.5.74 Subroutine DANI

Problem solved:

For a given interval  $[X1, X2]$  with functional values Y1 resp. Y2 and any permitted interpolation code JNTX, an energy grid from X1 to X2 is generated and the associate functional values are calculated. The grid is made by bisections of the interval  $[X1, X2]$  and the produced subintervals.

Bisectioning is stopped, if the functional value at a point  $x_h = 0.5 \cdot (x_n + x_m)$  -  $x_n$  and  $x_m$  are neighboured points - can be predicted by linear interpolation with an error less EPS. The interval from X1 to X2 can have maximally 300 points, including X1 and X2.

Definition:

SUBROUTINE DANI(X1, Y1, X2, Y2, JNTX, N, EPS)

Explanation of the arguments:

name	mode	type	dimension	remarks
X1	I	R4	scalar	lower limit of the interval
Y1	I	R4	scalar	functional value at the lower end of the interval
X2	I	R4	scalar	upper limit of the interval
Y2	I	R4	scalar	functional value at the upper end of the interval
JNTX	I	I4	scalar	predefined interpolation code
N	0	I4	scalar	number of produced energy points (and functional values) in the array XR (resp. YR). (XR and YR see common RE) ( $2 \leq N \leq 300$ )
EPS	I	R4	scalar	accuracy (see problem solved)

Used common block:

name of the common block: RE									
name	dimension	type	mode	word	name	dimension	type	mode	word
XR	(300)	R4	0	1	YR	(300)	R4	0	301

Called subprogram:

ERROR

2.3.5.75 Subroutine CROP

Problem solved:

A TAB1-record (see /3/ chapter R-2.3.3) stored in the arrays NBT, JNT, X and Y is cropped, and only the portion between XL and XH is retained. Portions of the range not defined by the original TAB1 record are set to zero. Unnecessary points are eliminated, if they can be predicted by interpolation between adjacent points to relative accuracy of EPS.

The subroutine originates from /3/ (see /3/ chapter R-6.4 and page R-75). Yet following changes have been made before using it:

- (1) EPS is equal EPSMIN in the case of LOSSTE equal .FALSE.  
In this case the variables EPSMAX, SCRCLA and NUMSCR have no meaning. If LOSSTE is equal .TRUE., EPS has also normally the value EPSMIN. It has the value EPSMAX, if two adjacent points fulfil the condition  $\left| \frac{X_2 - X_1}{X_1} \right| < 5.0 \cdot 10^{-6}$ .

In this case the number of points which have been eliminated, will be counted. To do this, first so called scratch classes are built in the following way. If EPSMIN is less than EPSMAX, five intervals are made from 0.0 up to EPSMIN, and another five from EPSMIN to EPSMAX. If EPSMIN is equal EPSMAX ten intervals are made from zero to EPSMIN. In the subroutine CROP the accuracy, by which the cropped points could have been predicted, is calculated, and the number of these points is added up with respect to the associated scratch class. The array element NUMSCR(I) contains the number of cropped points that could have been predicted with an relative error in the range

$$\text{SCRCLA}(I-1) < \text{EPS} \leq \text{SCRCLA}(I)$$

(SCRCLA(0) is defined to be zero).

- (2) Points in an interval are only eliminated, if all points between can be obtained by interpolation with an error less EPS.

Definition:

SUBROUTINE CROP (XL,XH,EPSMIN,LOSSTE,EPSSMAX,SCRCLA,NUMSCR,LOF,NBT,JNT,X,Y)

Explanation of the arguments:

name	mode	type	dimension	remarks
XL	I	R4	scalar	lower energy limit
XH	I	R4	scalar	upper energy limit
EPSMIN	I	R4	scalar	see "problem solved"
LOSSTE	I	L4	scalar	" " "
EPSSMAX	I	I4	scalar	" " "
SCRCLA	O	R4	(10)	" " "
NUMSCR	O	I4	(10)	" " "
LOF	O	I4	scalar	= 0 normal return = 1 the dimension of the arrays X and Y is too small to add zeros at the lower or upper end
NBT	IO	I4	(N1X)	interpolation ranges
JNT	IO	I4	(N1X)	interpolation codes
X	IO	R4	(N2X)	energy grid
Y	IO	R4	(N2X)	functional values

Used common blocks:

name of the common block: RECS									
name	dimension	type	mode	word	name	dimension	type	mode	word
MAT	scalar	I4	-	1	L1	scalar	I4	-	6
MF	scalar	I4	-	2	L2	scalar	I4	-	7
MT	scalar	I4	-	3	N1	scalar	I4	IO	8
C1	scalar	R4	-	4	N2	scalar	I4	IO	9
C2	scalar	R4	-	5					

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NR1	(125)	I4	-	1	N1X	scalar	I4	I	127
NBX	scalar	I4	-	126	N2X	scalar	I4	I	128

Called subprograms:

ERROR

TERP1

2.3.5.76 Subroutine WDA

Problem solved:

The contents of the array BK from I=1 to I=N (N given in common INOUT) are written on the direct access data set with the logical unit number NF=11. All written records have the length 400 4-Byte words. The data set has the following structure:

1. record (table of contents and addresses)

word number	contents	remarks
1	MATKDK	numerical material name
2	MIS	number of the isotope
3	NAME1	} name of the (KEDAK) data type in alphanumerical representation
4	NAME2	
5	IR	number of the record, where the data can be found
6	N	number of data measured in 4-Byte words
.		
.		
.		
word 1 until 6 are repeated for each entry		
399	IR	first free record
400	IW	first free word in record number 1

2. and following records  
contents of the array BK.



Definition:

SUBROUTINE WDA  
(not used entry)

ENTRY WDA1(BK)  
(passes the address of BK to the subroutine and initializes the data set on unit NF=11)

ENTRY WDA2(MATKDK,MIS,NAME1,NAME2)  
(does the writing of the contents of BK on unit NF=11)

ENTRY WDA3(IT)  
(prints the table of contents and addresses (record 1) of the dataset on unit NF=11 and returns the length of the filled portion of record 1.)

Explanation of the arguments:

name	mode	type	dimension	remarks
BK	I	R4	(NBX)	data to be written on logical unit number NF=11
MATKDK	I	I4	scalar	see problem solved
MIS	I	I4	scalar	" " "
NAME1	I	I4	scalar	" " "
NAME2	I	I4	scalar	" " "
IT	0	I4	scalar	length of the filled portion in record 1 (table of contents and addresses)

Used common block:

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NO	scalar	I4	I	1	N	scalar	I4	I	129
NR1	(127)	I4	-	2	NBKX	scalar	I4	I	130

Called subprogram :

ERROR

### 2.3.5.77 Subroutine RDAB

Problem solved:

The data written by the subroutine WDA (see 2.3.5.76) on the direct access data set with the logical unit number NF=11 are read into the array BK.

Organisation of the dataset see 2.3.5.76.

Definition:

SUBROUTINE RDAB

(This entry is never called.)

ENTRY RDA1(BK)

(This entry must be called before using all other entries. It passes the address of the array BK to the subroutine and reads the first record - table of contents and addresses - of the dataset.)

ENTRY RDA4(MATKDK,MIS,NAME1,NAME2,LNT,IR)

(The datatype "NAME1,NAME2" of the isotope number MIS of the material MATKDK is searched in the table of contents and addresses. If found the record number is stored in IR, the number of data in the quantity N of the common INOUT and RETURN is made.)

ENTRY RDA2(MATKDK,MIS,NAME1,NAME2,LNT)

(The datatype "NAME1,NAME2" of the isotope number MIS of the material MATKDK is searched in the table of contents and addresses. If found the data are read into the array BK.)

ENTRY RDA3(MATKDK)

(The number of datatypes, which are descending from ENDF /B MF=1,2, and 3, and are to be inserted in KEDAK, are counted in the variable NTYP of the common INOUT.)

Explanation of the arguments:

name	mode	type	dimension	remarks
BK	0	R4	(NBKX)	used to store data read from the direct access data set NF=11
MATKDK	I	I4	scalar	numerical KEDAK material name
MIS	I	I4	scalar	number of the isotope
NAME1	I	I4	scalar	first part of the KEDAK alphanumerical data type name
NAME2	I	I4	scalar	second part of the KEDAK alphanumerical data type name
LNT	0	I4	scalar	0 normal RETURN. Data found. 1 data not found
IR	0	I4	scalar	number of the record of the direct access data set NF=11, where the desired data starts

Used common blocks:

name of the common block: TAB									
name	dimension	type	mode	word	name	dimension	type	mode	word
ENDTAB	(25)	I4	-	1	NCNAME	(39)	I4	-	316
KEDTAB	(50)	I4	-	26	NRIF3	scalar	I4	-	355
KEDR	(240)	I4	I	76	NRIF4	scalar	I4	-	356

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NR1	(83)	I4	-	1	NR2	(44)	I4	-	85
NTYP	scalar	I4	0	84	N	scalar	I4	0	129

Called subprograms:

None

2.3.5.78 Subroutine WRECS

Problem solved:

TAB1 and LIST records (see /3/ chapter R-2.3.3 and chapter R-2.3.2) are stored on the direct access data set NF=12. One record has the length of 400 4-Byte words.

Organisation of the data set:

Record 1 until record BIB12 (table of contents and addresses)

word number	contents	remarks
1	MIS	number of the isotope for a specified material
2	NAME1	} name of a datatype in alphabetical form
3	NAME2	
4	EL for TAB1 records	lower energy limit
	0 for LIST records	
5	EH for TAB1 records	upper energy limit
	0 for LIST records	
6	IR	number of the record, where the data starts
7	N1	length of the arrays NBT and JNT for TAB1-records length of the array B for LIST-records
8	N2 for TAB1 records	length of the arrays X and Y
	0 for LIST records	

words number 1 until 8 are repeated for each entry.

Records BIB12+1, BIB12+2,... contain the data.

Definition:

SUBROUTINE WRECS

(This entry is never used)

ENTRY WRECS1(NBT,JNT,X,Y,B)

(This entry must be called first before using any other entry.

The addresses of NBT,JNT,X,Y and B are passed to the subroutine and the dataset is initialized.)

ENTRY WRECS2(MIS,NAME1,NAME2,EL,EH)

(TAB1-records are written on the dataset.)

ENTRY WRECS4(MIS,NAME1,NAME2)

(LIST-records are written on the dataset.)

Explanation of the arguments:

name	mode	type	dimension	remarks
NBT	I	I4	(N1X)	interpolation ranges
JNT	I	I4	(N1X)	interpolation codes
X	I	R4	(N2X)	energy grid
Y	I	R4	(N2X)	functional values
B	I	R4	(NBX)	"LIST-records"
MIS	I	I4	scalar	number of the isotope
NAME1	I	I4	scalar	first part of the alphanumerical data type name
NAME2	I	I4	scalar	second part of the alphanumerical data-type name
EL	I	R4	scalar	lower energy limit of the TAB1-record
EH	I	R4	scalar	upper energy limit of the TAB1-record

Used common blocks:

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NO	scalar	I4	I	1	BIB12	scalar	I4	I	79
NR1	(77)	I4	-	2					

name of the common block: RECS									
MAT	scalar	I4	-	1	L1	scalar	I4	-	6
MF	scalar	I4	-	2	L2	scalar	I4	-	7
MT	scalar	I4	-	3	N1	scalar	I4	I	8
C1	scalar	R4	-	4	N2	scalar	I4	I	9
C2	scalar	R4	-	5					

Called subprograms:

None

### 2.3.5.79 Subroutine RRECS

Problem solved:

The data written by the subroutine WRECS (see 2.3.5.78) on the direct access data set NF=12 are read into the arrays NBT, JNT, X, and Y for TAB1-records respectively B for LIST-records. Organisation of the dataset see 2.3.5.78.

Definition:

SUBROUTINE RRECS (MIS, NAME1, NAME2, EL, EH, LNT)

(Reads TAB1-records)

ENTRY RRECS4 (MIS, NAME1, NAME2, LNT)

(Reads LIST-records)

ENTRY IRRECS(NBT,JNT,X,Y,B)

(This entry must be called first. The addresses of NBT,JNT,X,Y, and B are passed to the subroutine.)

Explanation of the arguments:

name	mode	type	dimension	remarks
MIS	I	I4	scalar	number of the isotope for a specified material
NAME1	I	I4	scalar	first part of the desired datatype name in alphanumerical form
NAME2	I	I4	scalar	second part of the desired datatype name in alphanumerical form
EL	0	R4	scalar	lower energy limit for the TAB1-record
EH	0	R4	scalar	upper energy limit for the TAB1-record
LNT	0	I4	scalar	0 data found 1 data not found
NBT	0	I4	(N1X)	interpolation ranges
JNT	0	I4	(N1X)	interpolation codes
X	0	R4	(N2X)	energy grids
Y	0	R4	(N2X)	functional values
B	0	R4	(NBX)	"LIST-records"

Used common blocks:

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NR1	(78)	I4	-	1	NBX	scalar	I4	I	126
BIB12	scalar	I4	I	79	N1X	scalar	I4	I	127
NR2	(46)	I4	-	80	N2X	scalar	I4	I	128

name of the common block: RECS									
name	dimension	type	mode	word	name	dimension	type	mode	word
MAT	scalar	I4	-	1	L1	scalar	I4	-	6
MF	scalar	I4	-	2	L2	scalar	I4	-	7
MT	scalar	I4	-	3	N1	scalar	I4	0	8
C1	scalar	R4	-	4	N2	scalar	I4	0	9
C2	scalar	R4	-	5					

Called subprogram:

ERROR

#### 2.3.5.80 Subroutine STORE

Problem solved:

The subroutines described in 2.3.5.80 until 2.3.5.85 work with the so called "Dense Record Storage" (see /3/ chapter R-3.2). This is a fast core storage to store all possible ENDF/B record types in a dense manner. To work with this storage the subroutines given in /3/ pages R-37, R-39, and R-40 and pages R-71, R-72, R-74, R-77, and R-78 have been used. Some arrays, located in a common in the original routines have yet been replaced by arrays, which are now passed to the subroutines by call-statements.

The subroutine STORE stores one record, which is given in the arrays NBT, JNT,X,Y and B, in the dense storage. The organisation of the dense storage is as following.

A table of contents and addresses is hold in the common DENS (see commons used in the program 2.3.4). The meaning of these quantities is:

JMT(M)            arbitrary identification of the  $M^{\text{th}}$  record stored in array A.

JAT(M)            starting location in the array A of the  $M^{\text{th}}$  record.



JTT(M)            type of the M<sup>th</sup> record (see /3/ chapter R-2.3)

JLT(M)            number of words in array A used by the M<sup>th</sup>-record

JNS                next available location in the array A.

MNS                next available location in the arrays JMT,JAT,JTT and JLT.

MX                 length of the arrays JMT,JAT,JTT and JLT (= 100)

The dense storage itself is realized in the array A (see 2.3.3). The organization of this array can be found in /3/ chapter R-3.2.

Definition:

SUBROUTINE STORE(JT,MA,LOF)

ENTRY ISTORE(NBT,JNT,X,Y,B,A,LA)

(Passes the addresses of NBT,JNT,X,Y,B,A and LA to the subroutine.  
Must be called first.)

Explanation of the arguments:

name	mode	type	dimension	remarks
JT	I	I4	scalar	record type. May be 1,2,3,4,5,6
MA	I	I4	scalar	identification of the record to be stored. If a record with the same identification is already stored, it is deleted
LOF	O	I4	scalar	0 normal RETURN 1 dimension of array A too small
NBT	I	I4	(N1X)	interpolation ranges
JNT	I	I4	(N1X)	interpolation codes
X	I	R4	(N2X)	energy grid
Y	I	R4	(N2X)	functional values
B	I	R4	(NBX)	"LIST-records"
A	O	R4	(JX)	dense storage
LA	O	I4	(JX)	dense storage. Same starting address as array A

Used common blocks:

name of the common block: DENS									
name	dimension	type	mode	word	name	dimension	type	mode	word
JMT	(100)	I4	IO	1	JNS	scalar	I4	IO	401
JAT	(100)	I4	0	101	MNS	scalar	I4	IO	402
JTT	(100)	I4	0	201	MX	scalar	I4	I	403
JLT	(100)	I4	0	301					

name of the common block: INOUT									
NR1	(130)	I4	-	1	JX	scalar	I4	I	131

name of the common block: RECS									
MAT	scalar	I4	I	1	L1	scalar	I4	I	6
MF	scalar	I4	I	2	L2	scalar	I4	I	7
MT	scalar	I4	I	3	N1	scalar	I4	I	8
C1	scalar	R4	I	4	N2	scalar	I4	I	9
C2	scalar	R4	I	5					

Called subprograms:

DELETE

ERROR

2.3.5.81 Subroutine FETCH

Problem solved:

One record from dense storage is fetched (see 2.3.5.80 for definitions and sources).

Definition:

SUBROUTINE FETCH(MA,LNT)

ENTRY IFETCH(NBT,JNT,X,Y,B,A,LA)

(Passes the addresses of the arrays NBT,JNT,X,Y,B,A and LA to the subroutine. Must be called first.)

Explanation of the arguments:

name	mode	type	dimension	remarks
MA	I	I4	scalar	identification of the record to be fetched
LNT	0	I4	scalar	0 record has been found and moved to the common RECS and some of the arrays NBT,JNT,X,Y,B 1 record not found
NBT	0	I4	(NIX)	interpolation ranges
JNT	0	I4	(NIX)	interpolation codes
X	0	R4	(N2X)	energy grid
Y	0	R4	(N2X)	functional values
B	0	R4	(NBX)	"LIST-record"
A	I	R4	(JX)	dense storage
LA	I	I4	(JX)	dense storage (Same starting address as array A.)

Used common blocks:

name of the common block: RECS									
name	dimension	type	mode	word	name	dimension	type	mode	word
MAT	scalar	I4	0	1	L1	scalar	I4	0	6
MF	scalar	I4	0	2	L2	scalar	I4	0	7
MT	scalar	I4	0	3	N1	scalar	I4	0	8
C1	scalar	R4	0	4	N2	scalar	I4	0	9
C2	scalar	R4	0	5					

name of the common block: DENS									
JMT	(100)	I4	I	1	JNS	scalar	I4	-	401
JAT	(100)	I4	I	101	MNS	scalar	I4	I	402
JTT	(100)	I4	I	201	MX	scalar	I4	-	403
JLT	(100)	I4	-	301					

name of the common block: INOUT									
NR1	(125)	I4	-	1	N1X	scalar	I4	I	127
NBX	scalar	I4	I	126	N2X	scalar	I4	I	128

Called subprogram:

ERROR

2.3.5.82 Subroutine DELETE

Problem solved:

Deletes one record from dense storage or clears total dense storage.

(See 2.3.5.80 for definitions and sources.)

Definition:

SUBROUTINE DELETE(MA,A)

Explanation of the arguments:

name	mode	type	dimension	remarks
MA	I	I4	scalar	identification of the record to be deleted. If MA=0 all records in dense storage are deleted
A	IO	R4	(JX)	dense storage

Used common blocks:

name of the common block: DENS									
name	dimension	type	mode	word	name	dimension	type	mode	word
JMT	(100)	I4	IO	1	JNS	scalar	I4	IO	401
JAT	(100)	I4	IO	101	MNS	scalar	I4	IO	402
JTT	(100)	I4	IO	201	MX	scalar	I4	I	403
JLT	(100)	I4	IO	301					

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NR1	(130)	I4	-	1	JX	scalar	I4	I	131

Called subprograms:

None

2.3.5.83 Subroutine FPDS

Problem solved:

One point from a TAB1-record stored in dense storage is fetched. The data are stored in XP,YP and IP. The following output is produced:

NP	XP	YP	IP
$\leq 0$	$-1.0 \cdot 10^{20}$	0.0	1
$> 0$ and $< N2$	X(NP)	Y(NP)	interpolation code stored in dense storage
= N2	X(N2)	Y(N2)	2
= N2+1	X(N2)	0.0	2
$> N2+1$	$1.0 \cdot 10^{20}$	0.0	2

(TAB1-record see /3/ chapter R-2.3.3. For definition of dense storage and sources for the program see 2.3.5.80)

Definition:

SUBROUTINE FPDS(JA,NP,XP,YP,IP)

ENTRY IFPDS(A,LA)

(Must be called first before any FPDS-call. Passes the address of A resp. LA to the subroutine.)

Explanation of the arguments:

name	mode	type	dimension	remarks
JA	I	I4	scalar	starting index in dense storage (realized by the array A resp. LA) of the TAB1-record, which is constructed out of the arrays NBT,JNT,X and Y
NP	I	I4	scalar	desired point number
XP	O	R4	scalar	= X(NP),X(N2), $-1.0 \cdot 10^{20}$ , or $1.0 \cdot 10^{20}$ depending from the value of NP (see problem solved)
YP	O	R4	scalar	= Y(NP),Y(N2),or 0.0 depending from the value of NP (see problem solved)

name	mode	type	dimension	remarks
IP	0	I4	scalar	interpolation code used between X(NP) and X(NP+1), or 1 or 2 depending from the value of NP (see problem solved).
A	I	R4	(JX)	dense storage array
LA	I	I4	(JX)	dense storage array (Same starting address as array A.)

Used common blocks:

None

Called subprograms:

None

#### 2.3.5.84 Subroutine IPDS

Problem solved:

Interpolates one point in a TAB1-record in the dense storage.

If the given value of X, which is named XP, is less than X(1), the corresponding (desired) value of Y named YP is set to zero, NP (address of X and Y arrays, such that XP lies between X(NP) and X(NP+1)) is also set to zero, and the interpolation code IP to 1. If XP is greater than X(N2), YP is set to zero, NP to N2+1, and IP to 1. (TAB1-record see /3/ chapter R-2.3.3. For definition of dense storage and sources for the program see 2.3.5.80.)

Definition:

```
SUBROUTINE IPDS(JA, NP, XP, YP, IP)
```

```
ENTRY IIPDS(A, LA)
```

(Must be called first before any IPDS-call. Passes the address of A resp. LA to the subroutine.)

Explanation of the arguments:

name	mode	type	dimension	remarks
JA	I	I4	scalar	starting index in dense storage, which is realized by the array A resp. LA, of the TAB1-record
NP	IO	I4	scalar	index of X and Y arrays, such that XP lies between X(NP) and X(NP+1). If NP is greater than zero on input, it is used to start search. Otherwise search is started at NP=1
XP	I	R4	scalar	given value of X
YP	O	R4	scalar	corresponding (desired) value of Y
IP	O	I4	scalar	interpolation code used to compute YP, or 1 if XP is outside the TAB1-record range
A	I	R4	(JX)	dense storage array
LA	I	I4	(JX)	dense storage array. (Same starting address as array A.)

Used common blocks:

None

Called subprograms:

ERROR

TERP1

#### 2.3.5.85 Subroutine LRIDS

Problem solved:

Locates record MA in the dense storage. (For definition of dense storage and sources for the program see 2.3.5.80.)

Definition:

SUBROUTINE LRIDS(MA,JA,LNT)



Explanation of the arguments:

name	mode	type	dimension	remarks
MA	I	I4	scalar	identification of the record
JA	0	I4	scalar	starting index in array A for the record identified by MA
LNT	0	I4	scalar	= 0 address JA of the record identified by MA found = 1 record identified by MA not in dense storage

Used common block:

name of the common block: DENS									
name	dimension	type	mode	word	name	dimension	type	mode	word
JMT	(100)	I4	I	1	JNS	scalar	I4	-	401
JAT	(100)	I4	I	101	MNS	scalar	I4	I	402
JTT	(100)	I4	-	201	MX	scalar	I4	-	403
JLT	(100)	I4	-	301					

Called subprograms:

None

### 2.3.5.86 Subroutine SICKY

Problem solved:

Inserts the data in a KEDAK library<sup>+</sup> on the direct access data set IFILE=1. If no library exists, one is built up. If an isotope is yet present in the library and IMOD (see 2.2.6) has been set to 1, the isotope is replaced.

---

<sup>+</sup>This KEDAK library must not have been processed with or created by KEMA (see II or /4/). It must have been created by BRIGITTE itself.

If IMOD has the value zero, and an isotope is yet present in the library, the data are not inserted.

Deleting a specified isotope and compressing the library is done by the subroutine LISA (see 2.3.5.87). The data to be inserted in the KEDAK library are read from direct access dataset NF=11 by the subroutine RDAB.

Definition:

SUBROUTINE SICKY(NIS,NISO,IF4,IF5,BK,TAB,ITAB)

Explanation of the arguments:

name	mode	type	dimension	remarks
NIS	I0	I4	scalar	number of isotopes. Is set to 1 for NISO equal zero
NISO	I	I4	scalar	number of isotopes. Only used to set NIS to 1 for NISO equal zero
IF4	I	I4	scalar	= 1 only data from ENDF/B MF=4 have been converted ≠ 1 data from all files can be found in direct access data set NF=11
IF5	I	I4	scalar	= 1 only data from ENDF/B MF=5 have been converted ≠ 1 data from all files can be found in direct access data set NF=11
BK	L	R4	(NBKX)	used to store data read from direct access dataset NF=11
TAB	L	R4	(ITDIM)	used to built up the data to be written in KEDAK
ITAB	L	I4	(ITDIM)	same array as TAB

Used common blocks:

name of the common block: TAB									
name	dimension	type	mode	word	name	dimension	type	mode	word
ENDTAB	(25)	I4	-	1	NCNAME	(39)	I4	I	316
KEDTAB	(50)	I4	-	26	NRIF3	scalar	I4	I	355
KEDR	(240)	I4	I	76	NRIF4	scalar	I4	I	356

name of the common block: INOUT									
name	dimension	type	mode	word	name	dimension	type	mode	word
NRE1	(80)	I4	-	1	MATISO	(10)	I4	I	105
NAMKDK	(2)	I4	IO	81	NRE2	(14)	I4	-	115
MATKDK	scalar	I4	I	83	NKED	scalar	I4	IO	129
NTYP	scalar	I4	IO	84	NRE3	(3)	I4	-	130
NAMISO	(20)	I4	I	85	ITDIM	scalar	I4	I	133

Called subprograms:

DATUM

LISA

PRIZEI

RDA1

RDA2

RDA3

RDA4

(entries of RDAB)

SPLIT

TABPRT

### 2.3.5.87 Subroutine LISA

Problem solved:

The material "INAM1,INAM2" in the KEDAK library on the direct access data set on NT=1 is deleted. The library is compressed<sup>+</sup>.

Definition:

SUBROUTINE LISA(INAM1,INAM2,IFIRST)

---

<sup>+</sup>This subroutine can not be used for KEDAK libraries, which have been created by or processed with KEMA (see II), because the libraries generated by BRIGITTE have a special form. First in word 880 of the first record the number of the next available record can be found. Second each material starts on a new record.

Explanation of the arguments:

name	mode	type	dimension	remarks
INAM1	I	I4	scalar	first part of the alphanumerical KEDAK material name to be deleted
INAM2	I	I4	scalar	second part of the alphanumerical KEDAK material name to be deleted
IFIRST	0	I4	scalar	0 normal RETURN 1 the deleted material is the only one in the library. The library is regarded as empty and must be initialized in the calling program

Used common blocks:

None

Called subprograms:

None

#### 2.3.5.88 Subroutine TABPRT

Problem solved:

A table of contents of the KEDAK library, which contains the inserted material, is printed out. This routine can also be used for libraries, which have been created by KEMA (see II).

Definition:

SUBROUTINE TABPRT(MAT,NMAT,IFIL)

Explanation of the arguments:

name	mode	type	dimension	remarks
MAT	IO	R8	(NMAT)	MAT(1) = 'ALL':the material conversion table, the type conversion table, the material address table and the type address tables of all materials are printed out  = 'material names': like MAT(1)= 'ALL' with the exception that only the type address tables of the specified materials are printed out
NMAT	I	I4	scalar	number of material names in the array MAT
IFIL	I	I4	scalar	logical unit number of the KEDAK library, for which the table of contents is to be printed out

Used common blocks:

None

Called subprograms:

INITPG (entry of PAGE)

PAGE

PRIZEI

#### 2.3.5.89 Subroutine PAGE

Problem solved:

Called by TABPRT (see 2.3.5.88). Prints the heading of each page created by TABPRT.

Definition:

SUBROUTINE PAGE(TITLE,IL)

ENTRY INITPG(IFIL)

(Must be called first. The page counter is set to zero and the value of IFIL is passed to the subroutine.)

Explanation of the arguments:

name	mode	type	dimension	remarks
TITLE	I	R8	scalar	text printed in the heading of each page (e.g. a material name)
IL	0	I4	scalar	number of lines per page printed in the subroutine
IFIL	I	I4	scalar	logical file number of the KEDAK dataset

Used common blocks:

None

Called subprograms:

None

#### 2.3.5.90 Subroutine SPLIT

Problem solved:

The data given in the form dd.mm.jj in EBCDIC are splitted up in three integer values dd, mm, and jj.

Definition:

SUBROUTINE SPLIT(A,I1,I2,I3)

Explanation of the arguments:

name	mode	type	dimension	remarks
A	I	R8	scalar	date in the form dd.mm.jj
I1	0	I4	scalar	dd
I2	0	I4	scalar	mm
I3	0	I4	scalar	jj

Used common blocks:

None

Called subprogram:

CON

2.3.5.91 Subroutine CON

Problem solved:

A number in the range 0 until 99 represented in EBCDIC, is converted in a machine internal representation.

Definition:

SUBROUTINE CON(IA,IN)

Explanation of the arguments:

name	mode	type	dimension	remarks
IA	I	I2	scalar	number in EBCDIC
IN	O	I2	scalar	number in internal representation

Used common blocks:

None

Called subprogram:

CONVX

2.3.5.92 Subroutine A8FORM

Auxiliary program (member of a local library) used by BRIGITTE (see /13/).

Problem solved:

Prints texts with "big" letters. One "line" can contain up to 8 characters. One page can contain 3 lines.

Because the subroutine is not a part of the program BRIGITTE, only examples how to use the program together with an explanation of the arguments will be given in the following.

Program call:

CALL A8FORM(IA,T)

Effect of the call:

The text T is written on unit IA.

Explanation of the arguments:

name	mode	type	dimension	remarks
IA	I	I4	scalar	print output unit number
T	I	R8	scalar	text (8 characters) to be printed

Program call:

CALL RAMANF(IA)

CALL RAMEND(IA)

Effect of the call:

The top and the bottom of a frame of asterisk is produced. All A8FORM-calls between will produce additionally with the text the left and right parts of the frame.

Explanation of the argument:

name	mode	type	dimension	remarks
IA	I	I4	scalar	print output unit number

Examples for the use of the routine in the source program:

See subroutine TITEL (2.3.5.2).



2.3.5.93 Subroutine\_DEFI

Auxiliary program (member of a local library) used by BRIGITTE (see /14/).

Problem solved:

The static IBM-FORTRAN IV DEFINE FILE statement (see /9/) is made dynamic.

Because the subroutine is not a part of the program BRIGITTE, only an example how to use the program together with an explanation of the arguments will be given in the following.

Program call:

CALL DEFI(UNIT,NPRIQ,FORMAT,NBLK,NASS)<sup>+</sup>

Effect of the call:

The direct access dataset on unit number UNIT is defined.

Explanation of the arguments:

name	mode	type	dimension	remarks
UNIT	I	I4	scalar	logical unit number
NPRIQ	I	I4	scalar	number of records for the direct access data set (see 2.3.5.93)
FORMAT	I	R4	scalar	format of the records. May be 'L', 'E' or 'U' (see /9/)
NBLK	I	I4	scalar	number of words for FORMAT='U' or number of bytes for FORMAT='L' or 'E' of one record
NASS	P	I4	scalar	address of the associated variable (see /9/)

Example for the use of the routine in the source program:

See MAIN program (2.3.5.1)

---

<sup>+</sup>The argument list can be repeated

### 2.3.5.94 Subroutine DINF

Auxiliary program (member of a local library) used by BRIGITTE (see /14/).

#### Problem solved:

The data block length and the primary quantity, both given in the SPACE parameter of a DD-card (see /10/), are fetched and passed to the calling program.

Because the subroutine is not a part of the program BRIGITTE, only examples how to use the program together with an explanation of the arguments will be given in the following.

#### Program call:

```
CALL DINF(DDNAME,NBLK,NPRIQ)
```

#### Effect of the call:

NBLK is filled with the data block length and NPRIQ with the primary quantity of the SPACE parameter of the DD-card DDNAME.

#### Explanation of the arguments:

name	mode	type	dimension	remarks
DDNAME	I	R8	scalar	DD-name of the DD-card in interest
NBLK	0	I4	scalar	data block length (in byte) given in the SPACE parameter
NPRIQ	0	I4	scalar	primary quantity of the SPACE parameter. Number of records for the data-set

Examples for the use of the routine in the source program:

See subroutine DADEFI (2.3.5.7)

### 2.3.5.95 Subroutine FREESP

Auxiliary program (member of a local library) used by BRIGITTE (see /15/).

Problem solved:

The size of the unused part of the region (see /10/) - measured in units of 2 k-bytes - is fetched.

Because the subroutine is not a part of the program BRIGITTE, only an example how to use the program together with an explanation of the argument will be given in the following.

Program call:

CALL FREESP(I)

Effect of the call:

I contains the size of the unused region, measured in units of 2 k-bytes.

Explanation of the argument:

name	mode	type	dimension	remarks
I	0	I4	scalar	unused part of the region measured in units of 2 k-bytes

Example for the use of the routine in the source program:

See subroutine GSPACE (2.3.5.5).

### 2.3.5.96 Subroutine XTAREA

Auxiliary program (member of a local library) used by BRIGITTE (see /16/).

Problem solved:

Portions of the unused region or all of it are placed at the program disposal on execution time by using the IBM/370 macros GETMAIN and FREEMAIN (see /11/).

Because the subroutine is not a part of the program BRIGITTE, only an example how to use the program together with an explanation of the arguments will be given in the following.

Program call:

CALL XTAREA(K1,K2,K3,AREA)

Effect of the call:

The absolute address of the desired space is filled in K1. K2 is filled with zero, if the desired space was not available. K3 returns the address of the array AREA.

Explanation of the arguments:

name	mode	type	dimension	remarks
K1	0	I4	scalar	absolute address - measured in bytes - of the beginning of the desired space
K2	I0	I4	scalar	on input: number of bytes of the desired space. on output: if greater zero: normal RETURN. If equal zero: space was not available
K3	0	I4	scalar	absolute address of the beginning of the array AREA. Used to calculate the relative address of the desired space in the FORTRAN program
AREA	P	-	(1)	used as reference point (see K3)

Program call:

CALL REXTAR(K1,K2)

Effect of the call:

K2 bytes starting on the absolute address K1 are released

Explanation of the arguments:

name	mode	type	dimension	remarks
K1	I	I4	scalar	absolute address of area to be released
K2	I	I4	scalar	number of bytes to be released

Example for the use of the routine in the source program:

See subroutine GSPACE (2.3.5.5)

### 2.3.5.97 Function\_ZEIT

Auxiliary program (member of a local library) used by BRIGITTE (see /17/).

Problem solved:

The used CPU-time - relative to a fixed zero point - is measured.

Because the function is not a part of the program BRIGITTE, only examples how to use the program together with an explanation of the arguments will be given in the following.

Program call:

UHRZEI=0.0

ANFZEI=ZEIT(UHRZEI)

Effect of the call:

This call defines a reference point. All following calls will give times relative to this call.

Explanation of the argument:

name	mode	type	dimension	remarks
UHRZEI	I(0)	R4	scalar	used to define the reference point UHRZEI must have the value 0.0. It can be changed by this first ZEIT-call

Program call:

DURA=ZEIT(ANFZEI)

Effect of the call:

The CPU-time passed since the ZEIT(UHRZEI)-call is stored in DURA. It is given in seconds as REAL\*4 number.

Explanation of the argument:

name	mode	type	dimension	remarks
ANFZEI	I	R4	scalar	value produced at the first ZEIT-call

Examples for the use of the routine in the source program:

See MAIN-program (2.3.5.1) and subroutine  
PRIZEI (2.3.5.45).

#### 2.3.5.98 Subroutine DATUM

Auxiliary program (member of a local library) used by BRIGITTE (see /18/).

Problem solved:

The date and time of day is returned.

Because the subroutine is not a part of the program BRIGITTE, only an example how to use the program together with an explanation of the arguments will be given in the following.

Program:call:

CALL DATUM(DATE,UHRZEI)

Effect of the call:

The date and time of day is returned.

Explanation of the arguments:

name	mode	type	dimension	remarks
DATE	0	R8	scalar	date in the form dd.mm.yy (day, month, year)
UHRZEI	0	R8	scalar	time of day in the form hh.mm.ss (hour, minute, second)

Example for the use of the routine in the source program:

See subroutine PRIZEI (2.3.5.45)

### 2.3.5.99 Subroutine CONVX

Auxiliary program (member of a local library) used by BRIGITTE (see /19/).

Problem solved:

Conversion of machine internal numbers in alphanumerical representation (EBCDIC).

Because the subroutine is not a part of the program BRIGITTE, only examples how to use the program together with an explanation of the arguments will be given in the following.

Program call:

CALL CONVX(ARG,ALPH,FIELDD)

Effect of the call:

The machine internal number ARG is stored in alphanumerical representation by using the format FIELDD in the variable ALPH.

Explanation of the arguments:

name	mode	type	dimension	remarks
ARG	I	-	scalar	machine internal number
ALPH	0	- <sup>+</sup>	- <sup>+</sup>	alphanumerical representation
FIELDD	I	- <sup>+</sup>	- <sup>+</sup>	conversion format

<sup>+</sup>If no type or dimension is given, several possibilities are permitted.

Example for the use of the routine in the source program:

See subroutine CON (2.3.5.91).



## 2.4 Literature

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## 2.5 Source program

In the section 2.5.3 a complete listing of the source program is given. This listing includes all programs described in 2.3.5.1 until 2.3.5.91. No listing of the auxiliary library programs (2.3.5.92 - 2.3.5.99) is given.

To facilitate the use of the source program listing two tabulations will be given. In the first (2.5.1) you will find the subroutine and function names in the sequence of their appearance in the list of the source program. The subroutines and functions are numbered (first column). The names are given in the second column. The type, listed in the third column, means:

M	main program
SF	FORTRAN subroutine
B	BLOCK DATA program
RF4	FORTRAN function delivering REAL*4 results
SA	ASSEMBLER subroutine

In the fourth column the number of the section, where the description is given, can be found. Finally the fifth column contains entry names, if such one are existing. In section 2.5.2 an alphabetic listing of all subroutine-, function-, and entry names is given (excluding the auxiliary programs (see 2.3.5.92 - 2.3.5.99), which are contained in program libraries). The number defined in 2.5.1 is associated to each name. If the specified name is an entry of a subroutine, the subroutine name is listed in column 2.

2.5.1 List of the subroutine and function names in the sequence of their appearance in the source program list

number	name	type	described	entry names
1	MAIN	M	2.3.5.1	-
2	TITEL	SF	2.3.5.2	-
3	PRIZEI	SF	2.3.5.45	-
4	DADEFI	SF	2.3.5.7	-
5	GSPACE	SF	2.3.5.5	-
6	BLOCK DATA	B	2.3.5.4	-
7	PART1	SF	2.3.5.8	-
8	INITNL	SF	2.3.5.9	-
9	INGRID	SF	2.3.5.10	-
10	REGINE	SF	2.3.5.11	-
11	MARY	SF	2.3.5.12	-
12	SORTRE	SF	2.3.5.14	-
13	MARCEL	SF	2.3.5.20	MARGOT
14	MILLI	SF	2.3.5.22	-
15	MILLI1	SF	2.3.5.23	-
16	RESERR	SF	2.3.5.18	-
17	MELA	SF	2.3.5.24	-
18	MARYO1	SF	2.3.5.15	-
19	MARYO2	SF	2.3.5.19	-
20	FACTS	SF	2.3.5.13	-
21	FACPHI	SF	2.3.5.21	-
22	ANNICK	SF	2.3.5.25	-
23	ANNIO1	SF	2.3.5.30	-
24	ERF	SF	2.3.5.27	-
25	GAUSS	SF	2.3.5.28	-

number	name	type	described	entry names
26	BINT	RF4	2.3.5.29	-
27	GNRL	SF	2.3.5.26	-
28	ADD	RF4	2.3.5.68	-
29	SUB	RF4	2.3.5.69	-
30	MULT	RF4	2.3.5.70	-
31	MARION	SF	2.3.5.35	-
32	MYRIAM	SF	2.3.5.33	-
33	SOPHIE	SF	2.3.5.36	-
34	ELIMZ	SF	2.3.5.38	-
35	SUZY	SF	2.3.5.37	-
36	DZMSCH	SF	2.3.5.40	-
37	ADDPNT	SF	2.3.5.39	-
38	SUPNEG	SF	2.3.5.41	-
39	ISABEL	SF	2.3.5.42	-
40	ITAB1	SF	2.3.5.71	EITAB1
41	WCONT	SF	2.3.5.50	-
42	BETSY	SF	2.3.5.34	-
43	LEGP	RF4	2.3.5.43	-
44	RDAB	SF	2.3.5.77	RDA1
	-	-	-	RDA2
	-	-	-	RDA3
	-	-	-	RDA4
45	FILE5	SF	2.3.5.44	-
46	CHGINT	SF	2.3.5.73	-
47	DANI	SF	2.3.5.74	-
48	COMB	SF	2.3.5.66	IICOMB
49	ERROR	SF	2.3.5.46	-
50	MAGGY	SF	2.3.5.31	-

number	name	type	described	entry names
51	SEARCH	SF	2.3.5.48	-
52	TERP1	SF	2.3.5.72	-
53	WDA	SF	2.3.5.76	WDA1
	-	-	-	WDA2
	-	-	-	WDA3
54	WREC	SF	2.3.5.51	IWREC
55	WRECS	SF	2.3.5.78	WRECS1
	-	-	-	WRECS2
	-	-	-	WRECS3
56	CROP	SF	2.3.5.75	-
57	LUCY	SF	2.3.5.16	-
58	RRECS	SF	2.3.5.79	IRRECS
	-	-	-	RRECS4
59	PRTPID	SF	2.3.5.57	-
60	PUTPID	SF	2.3.5.63	-
61	RREC	SF	2.3.5.49	IRREC
62	PRHOL	SF	2.3.5.56	-
63	PUHOL	SF	2.3.5.62	-
64	PRCONT	SF	2.3.5.52	-
65	PUCONT	SF	2.3.5.58	-
66	PRLIST	SF	2.3.5.53	-
67	PULIST	SF	2.3.5.59	-
68	PRTAB1	SF	2.3.5.54	-
69	PUTAB1	SF	2.3.5.60	-
70	PRTAB2	SF	2.3.5.55	-
71	PUTAB2	SF	2.3.5.61	-
72	CXFP	SF	2.3.5.64	-
73	XTEXT	SF	2.3.5.47	-

number	name	type	described	entry names
74	EPT	SF	2.3.5.17	-
75	SCALE	SF	2.3.5.65	GENER
	-	-	-	INITSC
76	COMBP	SF	2.3.5.67	-
77	DENISE	SF	2.3.5.32	-
78	FETCH	SF	2.3.5.81	IFETCH
79	STORE	SF	2.3.5.80	ISTORE
80	DELETE	SF	2.3.5.82	-
81	FPDS	SF	2.3.5.83	IFPDS
82	IPDS	SF	2.3.5.84	IIPDS
83	LRIDS	SF	2.3.5.85	-
84	SICKY	SF	2.3.5.86	-
85	LISA	SF	2.3.5.87	-
86	TABPRT	SF	2.3.5.88	-
87	PAGE	SF	2.3.5.89	INITPG
88	SPLIT	SF	2.3.5.90	-
89	CON	SF	2.3.5.91	-
90	PRIEIN	SF	2.3.5.3	-
91	BUFREG	SA	2.3.5.6	-

The meaning of the abbreviations in the type column may be found in section 2.5.



2.5.2 Index of the subroutine, function and entry names in alphabetic order

name	entry of	number in section 2.5.1
ADD	-	28
ADDPNT	-	37
ANNICK	-	22
ANNIO1	-	23
BETSY	-	42
BINT	-	26
BLOCK DATA	-	6
CHGINT	-	46
COMB	-	48
COMBP	-	76
CON	-	89
CROP	-	56
CXFP	-	72
DADEFI	-	4
DANI	-	47
DELETE	-	80
DENISE	-	77
DZMSCH	-	36
EITAB1	ITAB1	40
ELIMZ	-	34
EPT	-	74
ERF	-	24
ERROR	-	49
FACPHI	-	21
FACTS	-	20
FETCH	-	78

name	entry of	number in section 2.5.1
FILE5	-	45
FPDS	-	81
GAUSS	-	25
GENER	SCALE	75
GNRL	-	27
GSPACE	-	5
IFETCH	FETCH	78
IFPDS	FPDS	81
IICOMB	COMB	48
IIPDS	IPDS	82
INGRID	-	9
INITNL	-	8
INITPG	PAGE	87
INITSC	SCALE	75
IPDS	-	82
IRREC	RREC	61
IRRECS	RRECS	58
ISABEL	-	39
ISTORE	STORE	79
ITAB1	-	40
IWREC	WREC	54
LEGP	-	43
LISA	-	85
LRIDS	-	83
LUCY	-	57
MAGGY	-	50
MAIN	-	1

name	entry of	number in section 2.5.1
MARCEL	-	13
MARGOT	MARCEL	13
MARION	-	31
MARY	-	11
MARYO1	-	18
MARYO2	-	19
MELA	-	17
MILLI	-	14
MILLI1	-	15
MULT	-	30
MYRIAM	-	32
PAGE	-	87
PART1	-	7
PRIEIN	-	90
PRIZEI	-	3
PRCONT	-	64
PRHOL	-	62
PRLIST	-	66
PRTAB1	-	68
PRTAB2	-	70
PRTPID	-	59
PUCONT	-	65
PUHOL	-	63
PULIST	-	67
PUTAB1	-	69
PUTAB2	-	71
PUTPID	-	60

name	entry of	number in section 2.5.1
RDAB	-	44
RDA1	RDAB	44
RDA2	RDAB	44
RDA3	RDAB	44
RDA4	RDAB	44
REGINE	-	10
RESERR	-	16
RREC	-	61
RRECS	-	58
RRECS4	RRECS	58
SCALE	-	75
SEARCH	-	51
SICKY	-	84
SOPHIE	-	33
SORTRE	-	12
SPLIT	-	88
STORE	-	79
SUB	-	29
SUPNEG	-	38
SUZY	-	35
TABPRT	-	86
TERP1	-	52
TITEL	-	2
WCONT	-	41
WDA	-	53
WDA1	WDA	53
WDA2	WDA	53

name	entry of	number in section 2.5.1
WDA3	WDA	53
WREC	-	54
WRECS	-	55
WRECS1	WRECS	55
WRECS2	WRECS	55
WRECS4	WRECS	55
XTEXT	-	73

2.5.3 Source program listing

The following pages contain a listing of the whole source program. This is a listing of all the subroutines and functions described in 2.3.5.1 until 2.3.5.91. A listing of the auxiliary programs described in 2.3.5.92 until 2.3.5.99 is not given.



RETURN	100	WRITE(NO,SPACE)	500
END	110	C JX WIRD GLEICH IADIM GESETZT.	510
		C NBKX MUSS EIN VIELFACHES VON 2 SEIN.	520
		C ITDIM WIRD GLEICH MAXJ(ITDIM,NBX+2*(N1X+N2X)+IADIM) GESETZT.	530
		C NBX=ITDIM-IADIM-2*(N1X+N2X)	540
		C ITDIM WIRD AUF VIELFACHES VON 880 ABGERUNDET.	550
		C	560
	10	IDUR=1	570
C SUBROUTINE GSPACE(F)	20	6 JX=IADIM	580
C DIE GROESSE DES VON PROGRAMMEN UND PUFFERN NICHT BENUTZTEN TEILS	30	NBKX=(NBKX/2)*2	590
C DER REGION WIRD BESTIMMT. DAVON ABGEZOGGEN WIRD DER PLATZ, DER	40	ITDIM=MAXJ(ITDIM,NBX+2*(N1X+N2X)+IADIM)	600
C SPATER FUER PUFFER VERWENDUNG FINDEN KOENNTE, SCHIE DER VCM	50	NBX=ITDIM-IADIM-2*(N1X+N2X)	610
C SYSTEM BENOETIGTE PLATZ. DER SO VERBLEIBENDE TEIL WIRD DEM	60	ITDIM=(ITDIM/880)*880	620
C PROGRAMM ALS DATENSPEICHER ZUR VERFUEGUNG GESTELLT.	70	IF(IDUR.EQ.1) WRITE(NO,112)	630
	80	112 FORMAT(' PLANNED DIMENSIONS:')	640
	90	IF(IDUR.EQ.2) WRITE(NO,113)	650
	100	113 FORMAT(' NEW DIMENSIONS:')	660
C DIMENSION F(1)	110	WRITE(NO,107) NBX,N1X,N1X,N2X,N2X,NBKX,IADIM,ITDIM	670
C COMMON/INJUT/ NO,NI,NR1(123),NBX,N1X,N2X,N,NBKX,JX,IADIM,ITDIM,	120	107 FORMAT(' B(',I5,')',NBT(',I4,')',JNT(',I4,')',X(',I6,')',Y(',	680
C * IVER,IVGENR,IVJNT,IVX,IVY,IVB,IVDENS	130	2I6,')',BK(',I6,')',A(',I6,')',TAB(',I6,')')	690
C NAMELIST/SPACE/ NBX,N1X,N2X,NBKX,IADIM,ITDIM	140	ISUM=NBKX+NBX+2*(N1X+N2X)+IADIM	700
C WRITE(NO,100)	150	IF(ISUM.LE.KW) GOTO 7	710
100 FORMAT('ODYNAMIC WORKING SPACE ALLOCATION.'/'+',33('_'))	160	WRITE(NO,108)	720
C CALL FREESP(J)	170	108 FORMAT('ODATA REGION NOT LARGE ENOUGH TO FULFILL PLANNED DIMENSION	730
C WRITE(NO,101) J	180	1S.'/ ' DIMENSIONS WILL BE SHORTENED.')	740
101 FORMAT(' FREE REGION=',I4,' K-BYTE.')	190	9 FAK=(1.0*KW)/ISUM	750
IF(J.GE.10) GOTO 1	200	NBX=NBX*FAK	760
4 WRITE(NO,102)	210	N1X=N1X*FAK	770
102 FORMAT(' ***** NOT ENOUGH REGION.')	220	N2X=N2X*FAK	780
2 WRITE(NO,103)	230	NBKX=NBKX*FAK	790
103 FORMAT(' ***** PROGRAM ENDED DUE TO ERROR.')	240	JX=JX*FAK	800
STOP	250	IADIM=IADIM*FAK	810
3 WRITE(NO,111) IBUF	260	ITDIM=ITDIM*FAK	820
111 FORMAT(' ***** ERROR NUMBER ',I2,' IN BUFREG.')	270	IDUR=2	830
GOTO 2	280	GOTO 6	840
	290	7 IF(1.1*ISUM.GT.KW.OR.IDUR.GE.2) GOTO 9	850
C PLATZ FUER PUFFER	300	WRITE(NO,114)	860
	310	114 FORMAT(' OPLANNED DIMENSIONS WILL BE INCREASED, BECAUSE THERE IS EN	870
	320	OUGH DATA REGION.')	880
	330	GOTO 9	890
	340	C	900
	350	C DISPLACEMENTS RELATIV ZU BK(1).	910
	360	C	920
	370	8 IVGENR=NBKX	930
	380	IVJNT=IVGENR+N1X	940
	390	IVX=IVJNT+N1X	950
	400	IVY=IVX+N2X	960
	410	IVB=IVY+N2X	970
	420	IVDENS=IVB+NBX	980
	430	WRITE(NO,109) NBKX,IVGENR,ITDIM,IVGENR,N1X,IVJNT,N1X,IVX,N2X,IVY,	990
	440	N2X,IVB,NBX,IVDENS,IADIM	1000
	450	109 FORMAT(' OVERLAY OF DATA REGION.')	1010
	460	1' DISPLACEMENT RELATIV TO BK(1) ARRAY NAME LENGHT'/	1020
	470	2/'+',62['_')//	1030
	480	3' = 0 BK NBKX=',	1040
	490		



```

3 I7/
4'   IVGENR=' ,I9,12X,'TAB',8X,'ITDIM=' ,I6/
5'   IVGENR=' ,I9,12X,'NBT',8X,'N1X=' ,I8/
6'   IVJNT =' ,I9,12X,'JNT',8X,'N1X=' ,I8/
7'   IVX   =' ,I9,12X,'X  ',8X,'N2X=' ,I8/
8'   IVY   =' ,I9,12X,'Y  ',8X,'N2X=' ,I8/
9'   IVB   =' ,I9,12X,'B  ',8X,'NBX=' ,I8/
A'   IVDENS=' ,I9,12X,'A  ',8X,'IADIM=' ,I6//)
WRITE(ND,110)
110 FORMAT(' END OF DYNAMIC WORKING SPACE ALLOCATIGN.'/ '1')
RETURN
END

```

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1050
1060
1070
1080
1090
1100
1110
1120
1130
1140
1150
1160

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```

BLOCK DATA
COMMON/TAB/ENDTAB(25),KEDTAB(50),KEDR(240),NCNAME(39),NRIF3,NRIF4
DIMENSION KDR1(123),KDR2(84)
EQUIVALENCE (KEDR(1),KDR1(1)),(KEDR(124),KDR2(1))
INTEGER ENDTAB
DATA ENDTAB/1,2,3,4,16,17,18,22,23,24,25,27,28,29,
* 102,103,104,105,106,107,108,201,206,207,251/
DATA KEDTAB/'SGT',,,'SGN',,,'
* 'SGX',,,'SGI',,,'
* 'SG2N',,,'SG3N',,,'
* 'SGF',,,'SGIA',,,'
* 'SGI3',,A,'SG2N',,A
* 'SG3N',,A,'SGA',,,'
* 'SGIP',,,'SGNI',,,'
* 'SGG',,,'SGP',,,'
* 'SGD',,,'SGH3',,,'
* 'SGHE',,3,'SGAL',,P
* 'SG2H',,E,'SGTR',,,'
* 'ETA',,,'ALPH',,A
* 'MUEL',,,'
COMMON/XBAR/XX(4,10)
DATA XX/0.005252,0.051755,0.112925,0.169150,0.037171,0.163095,
* 0.265600,0.349780,0.103126,0.288421,0.404385,0.480571,0.207836,
* 0.431766,0.547724,0.617825,0.359852,0.599210,0.704048,0.762381,
* 0.574283,0.803560,0.882440,0.922898,0.879334,1.053224,1.096835,
* 1.111387,1.334810,1.393010,1.374373,1.350285,2.105227,1.916230,
* 1.786357,1.697511,4.390800,3.301643,2.824583,2.546602/
DATA KDR1/'RANG',,RES',,14511,'ISOT',,1',,14580,'ISOT',,2',,
X 14590,'ISOT',,3',,14600,'PLNL',,E',,14570,
X 'CHIC',,R',,14560,'RES',,,' ,21520,
X 'ST',,,' ,21530,'STD',,,' ,21540,
X 'STGF',,,' ,21550,'SGT',,,' ,30010,
X 'SGN',,,' ,30020,'SGX',,,' ,30020,
X 'SGI',,,' ,30040,'SGIZ',,,' ,30050,'SGIZ',,C',,
X 30051,'SG2N',,,' ,30160,'SG3N',,,' ,30170,
X 'SGF',,,' ,30190,'SGIA',,,' ,30220,
X 'SGI3',,A',,30230,'SG2N',,A',,30240,
X 'SG3N',,A',,30250,'SGA',,,' ,30270,
X 'SGIP',,,' ,30280,'SGNI',,,' ,30290,

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X 'SGG',,,' ,31020,'SGP',,,' ,31030,
X 'SGD',,,' ,31040,'SGH2',,,' ,31050,
X 'SGHE',,3',,31060,'SGAL',,P',,31070,
X 'SG2H',,E',,31080,'SGTR',,,' ,32010,
X 'ETA',,,' ,32060,'ALPH',,A',,32070,
X 'MUEL',,,' ,32510,'NUE',,,' ,34520,
X 'NUEP',,,' ,34550,'CHIF',,,' ,34610,
X 'CHIF',,D',,34620/
DATA KDR2/
* 'SGNC',,,' ,40022,'SGIL',,,' ,40021,
X 'SGIC',,,' ,40042,'SGIL',,Z',,40051,
X 'SGIC',,Z',,40052,'SGNI',,L',,40251,
X 'SGNI',,C',,40292,'LEGN',,L',,44631,
X 'LEGN',,C',,44632,'LEGI',,L',,44641,
X 'LEGI',,C',,44642,'LEGI',,LZ',,44651,
X 'LEGI',,CZ',,44652,'LGNL',,L',,44661,
X 'LGNL',,C',,44662,'CHII',,,' ,50040,
X 'CHII',,C',,50051,'SEDI',,C',,50053,
X 'CHI2',,N',,50160,'SED2',,N',,50163,
X 'CHI3',,N',,50170,'SED3',,N',,50173,
X 'SEDF',,,' ,54523,'CHIF',,Z',,54610,
X 'SEDF',,P',,54613,'CHIF',,CZ',,54620,
X 'SEDF',,D',,54623/
COMMON/INOUT/ NO,NI,NR1(123),NBX,N1X,N2X,N,NBXX,JX,IADIM,ITDIM,
* NR2(21)
DATA NO,NI,NBX,N1X,N2X,NBXX,JX,IADIM,ITDIM/6,5,3000,500,5000,
*12000,20000,20000,32560/
END

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SUBROUTINE PART1(NIS,MISO,KIF,KIF5,BK,NBT,JNT,X,Y,B,A)
COMMON/INOUT/NR1(2),NT,NR2(139),
*LABEL,LNER,ITAPE,MATP,EB,NI,SG,IF4,IF5
DIMENSION BK(1),NBT(1),JNT(1),X(1),Y(1),B(1),A(1)
CALL INITNL(MISO,321)
CALL IWREC(NBT,JNT,X,Y,B)
CALL IRREC(NBT,JNT,X,Y,B)
CALL FITAB1(NBT,JNT,X,Y)
CALL WRECS1(NBT,JNT,X,Y,B)
CALL WDA1(RK)
KIF=IF4
KIF5=IF5
CALL INGRID(MATP,LABEL,LRP,LEI,LNU,IF4,IF5)
IF(IF4.EQ.1)GOTO 11
IF(IF5.EQ.1)GOTO 12
3 CALL REGINE(MATP,LNER,LISRES,LISUNR,NIS,LFWRES,LFW,EB,LRP,
* NISO,BK,NBT,JNT,X,Y,B,A)
CALL DENISE(MATP,LRP,LISRES,LISUNR,NIS,LFW,LNU,EK,NBT,JNT,X,Y,B,A)
CALL SUZY(LNU,ITAPE,BK,BK,BK,BK,NBT,JNT,X,Y,B)
IF(IF4.EQ.-1)GOTO 14
11 CALL ISABEL(MATP,IF4,BK,BK,BK,NBT,JNT,X,Y,B,A,A,A)
14 IF(IF4.EQ.1.OR.IF5.EQ.-1)GOTO 13
12 CALL FILES(MATP,IF5,BK,BK,NBT,JNT,X,Y,B,B)

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```

13 CALL WDA3(IW)
RETURN
21 REWIND NT
STOP
END

SUBROUTINE INITNL(MISO,*)
LOGICAL*4 SALINE
COMMON/INOUT/NO,NI,NT,MODE,LEDIT1,LEDIT2,LEDITS,KEDAVE,ABIS(20),
* NOTTRL(25),BIB12,SALINE,
*NAMKDK,MATKDK,NTYP,NAMISO(10),MATISO(10),EBISO(10),ICOMB,NR1(15),
*EPSRES,SCREPS,LABEL,LNER,ITAPE,MATP,EB,NISC,IF4,IF5,FAICOD,MLSLSW,
*NR3(1),ELINEG
REAL*8 NAMKDK,NAMISO,NOTTRL,BLANC
DATA BLANC/' '/
COMMON/TAB/NR2(315),NCNAME(39),NRIF3,NRIF4
INTEGER BIB12,ELINEG
NAMELIST/INPUT/MODE,LABEL,LNER,LEDIT1,LEDIT2,LEDITS,KEDAVE,
* ITAPE,MATP,NAMKDK,MATKDK,EB,NISC,IF4,ICOMB,IF5,NOTTRL,
* BIB12,SALINE,EPSRES,SCREPS,FAICOD,MLSLSW,ELINEG
NAMELIST/ISO/NAMISO,MATISO,EBISO
NI=5
NC=6
NT=9
DO 20 I=1,39
20 NCNAME(I)=0
DO 25 I=1,25
25 NOTTRL(I)=BLANC
C-----MODE=DATA INPUT FORMAT DEFINED AS FOLLOWS---
C =1 BINARY TAPE, STANDARD ARRANGEMENT---
C =3 BCD CARD IMAGE TAPE---
C-----LABEL=IDENTIFICATION NUMBER OF THE INPUT TAPE,TREATED AS FOLLOWS
C =GREATER THAN ZERO---READ TPID AND CHECK LABEL
C =EQUAL ZERO---READ TPID, DO NOT CHECK LABEL
C =LESS THAN ZERO---DO NOT READ TPID OR CHECK LABEL
C-----LNER=0 CALCULATE ENERGIES USING 51 ENERGY POINTS PER RESONANCE
C >1 CALCULATE ENERGIES USING LNER ENERGY POINTS PER RESONANCE
C-----KEDAVE=0 OLD KEDAK VERSION.
C 1 NEW KEDAK VERSION.
C-----ICOMB=0 THE INTERPOLATION SCHEME IN COMBP IS CHOOSEN
C BY THE ROUTINE.
C 1 THE INTERPOLATION SCHEME USED IN COMBP IS
C ALWAYS LINEAR-LINEAR.
C-----BIB12=1,2,.....,9,10 NUMBER OF BLOCKS OF 400 WORDS FOR TABLE OF
C CONTENTS FOR FILE 12.
C-----MLSLSW=SWITCH FOR USING MULTILEVEL OR SINGLELEVEL BREIT-WIGNER
C FORMALISM IN RESOLVED RESONANCE CALCULATING SECTION.
C =0 THE ENDF/B RECOMMENDED FORMALISM IS USED.
C =1 SINGLE LEVEL FORMALISM IS USED INDEPENDENTLY FROM
C THE ENDF/B RECOMMENDED FORMALISM.
ICOMB=0
MODE=3

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LABEL=-1
LNER=0
LEDIT1=1
LEDIT2=1
LEDITS=0
KEDAVE=1
ITAPE=0
NISC=0
IF4=0
IF5=0
NRIF3=0
NRIF4=0
BIB12=1
SALINE=.FALSE.
EPSRES=1.0E-2
SCREPS=5.0E-3
FAICOD=10.0
MLSLSW=0
FLINEG=)
READ(NI,INPUT,END=21)
WRITE(NO,INPUT)
BIB12=MAXO(1,MINO(BIB12,10))
10 IF(MODE.EQ.1.OR.MODE.EQ.3)GOTO 7
WRITE(NO,6)MODE
6 FORMAT(1X,'MODE =',I4,' OUT OF RANGE')
STOP
7 IF(NISO.LE.1)GOTO 9
READ(NI,ISO)
WRITE(NO,I50)
9 MISO=NISO
RETURN
21 RETURN 1
END

SUBROUTINE INGRID(MATP,LABEL,LRP,LFI,LNU,IF4,IF5)
DIMENSION IOFFCS(9)
EQUIVALENCE (IOFFCS(1),MAT)
COMMON/INOUT/NO,NI,NT,MODE,LEDIT1,NR1(2),KEDAVE,NR2(72),NAMKDK,
*MATKDK,NR3(70),FLINEG
COMMON/RECS/MAT,ME,MT,C1,C2,L1,L2,N1,N2,NS
INTEGER ELINEG
REAL*8 NAMKDK
DO 1 I=1,9
1 IOFFCS(I)=0
NS=1
IF((MODE.LT.1).OR.(MODE.GT.3)) CALL ERROR(400)
REWIND NT
IF(LABEL.LT.0)GOTO 100
CALL RREC(6,NT,MODE,-1.0)
GOTO(120,130,120),MODE
120 LARE=MAT
GOTO 140

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130	LABEL=IABS(MF)	150	IF(LFI.EQ.1.AND.IF452.NE.1) WRITE(NO,1452)	740
140	IF(LABEL.EQ.0)GOTO 100	200	1452 FORMAT('0',120('*'))/' ',29('*'),' WARNING: FISSIONABLE MATERIAL. N	750
	IF(LABEL.NE.LABEL) CALL ERROR(401)	210	*0 SECTION 452 ( NUE ) FOUND. ',29('*'))/' ',120('*'))	760
100	WRITE(NO,450)	220	IF(LFI.NE.1.OR.IF452.EQ.0) GOTO 180	770
450	FORMAT('1')	230	CALL SEARCH(NT,MODE,MATP,1,452,LNT)	780
	CALL RAMANF(NO)	240	IF(LNT.EQ.1) CALL ERROR(202)	790
	CALL A8FORM(NO,'BRIGITTE')	250	LNU=L2	800
	CALL RAMEND(NO)	260	WRITE(NO,250)	810
	WRITE(NO,150) MATP,NAMKDK,MATKDK	270	250 FORMAT(1H1,'NUE DATA'/)	820
150	FORMAT('///' MATERIAL ',6X,I8/1X,	280	GCTO(190,200),LNU	830
	*'KEDAK NAME ',4X,A8/1X, 'KEDAK NUMBER ',4X,I6)	290	C DATA IS POLYNOMIAL COEFFICIENTS	840
	WRITE(NO,160) LABEL,MODE	300	190 CALL RREC(2,NT,MODE,-1.0)	850
160	FCRMT(1X,'ENDF/B TAPE ',I6,' IN MODE ',I2)	310	WRITE(NO,350)	860
	WRITE(NO,451)	320	350 FORMAT(1H0,'DATA IS POLYNOMIAL CCEFFICIENTS'/)	870
451	FORMAT(' NEGATIVE CROSS SECTION VALUES FOR SGT AND SGN WILL')	330	CALL WRECS4(0,4H NUE,4H )	880
	IF(FLINEG.NE.1) WRITE(NO,452)	340	CALL WREC(2,NO,4)	890
	IF(FLINEG.EQ.1) WRITE(NO,453)	350	GCTO 240	900
452	FCRMT('* ',51X,' N O T BE CHANGED.'//')	360	C LNU=2 DATA IS TABULAR	910
453	FORMAT('* ',51X,' BE D E L E T E D.'//')	370	200 CALL RREC(3,NT,MODE,-1.0)	920
	CALL RAMANF(NO)	380	WRITE(NO,260)	930
	CALL A8FORM(NO,NAMKDK)	390	260 FORMAT(1H0,'DATA IS TABULAR'/)	940
	CALL RAMEND(NO)	400	CALL WRECS2(0,1,452,0.0,0.0)	950
	WRITE(NO,450)	410	CALL WREC(3,NC,4)	960
	CALL RAMANF(NO)	420	CONTINUE	970
	IF(KEDAVE.EQ.0) CALL A8FORM(NO,' O L D '	430	WRITE(NO,251)	980
	IF(KEDAVE.NE.0) CALL A8FORM(NO,' N E W '	440	251 FORMAT(' NUE PROCESSED')	990
	CALL A8FORM(NO,'FORM OF '	450	180 RETURN	1000
	CALL A8FORM(NO,' KEDAK '	460	END	1010
	CALL RAMEND(NO)	470		
	CALL SEARCH(NT,MODE,MATP,1,451,LNT)	480		
	IF(LNT.EQ.1) CALL ERROR(201)	490		
	LRP=L1	500		
	LFI=L2	510		
	NC=N2	520	SUBROUTINE REGINE(MATP,LNER,LISRES,LISUNR,NIS,LFWRES,LFW,EB,	10
	CALL RREC(5,NT,MODE,-1.0)	530	* LRP,NISO,BK,NBT,JNT,X,Y,B,AH)	20
	IF452=0	540		30
	WRITE(NO,400)	550	C PROCESS RESONANCE DATA	40
400	FORMAT(1H1 //30X,'ENDF/B FILE 1'//30X,13('*'))//)	560	C	50
	CALL PRIZEI	570	COMMON/INOUT/NO,NI,NT,MODE,NR1(4), ABIS(20),NR2(52),NAMKDK,MATKDK,	60
	IF(LEDT1.NE.1)GOTO 165	580	*NR3(21),MATS0(10),FBISO(10),NR4(4),N,NBKK,XJ	70
	CALL WREC(5,NO,4)	590	COMMON/RECS/MAT,MF,MT,C1,C2,L1,L2,N1	80
	WRITE(NO,410)	600	COMMON/FLCT/ NR5(6),NDIM6,NPEMAX	90
410	FCRMT('//20X,'CONTENTS OF THE FILES'//1X,'FILE',3X,'MAT',3X,	610	DIMENSION AWRISO(20),NAM1(2),BK(1),NBT(1),JNT(1),X(1),Y(1),B(1),	100
	* 'NUMB. OF REC.'//)	620	*AH(1)	110
165	DO 170 I=1,NC	630	REAL*8 NAM8,NAMKDK	120
	CALL RREC(1,NT,MODE,-1.0)	640	LOGICAL*1 LNAM(8)	130
	IF(L2.EQ.452) IF452=1	650	EQUIVALENCE (NAMKDK,LNAM(1))	140
	IF(LEDT1.NE.1)GOTO 170	660	EQUIVALENCE (NAM8,NAM1(1))	150
	WRITE(NO,420) L1,L2,N1	670	NISMAX=10	160
420	FCRMT(14,4X,13,6X,14)	680	JRES=0	170
170	CONTINUE	690	WRITE(NO,20)	180
	IF(IF4.EQ.1.OR.IF5.EQ.1)GOTO 180	700	20 FORMAT(1H1,30X,'ENDF/B FILE 2'//31X,13('*'))//)	190
C		710	CALL PRIZEI	200
C	NUE	720	CALL SEARCH(NT,MODE,MATP,2,151,LNT)	210
C		730	IF(LNT.EQ.0)GOTO 300	220
			WRITE(NO,301)	230

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301 FORMAT(1X,'FILE 2 NOT PRESENT FOR THIS MATERIAL')
   NIS=1
   LFW=0
   LFWRES=0
   NER=0
   LISRES=1
   LISUNR=1
   BACKSPACE NT
   RETURN
300 AWR=C2
   NIS=N1
   IF(NIS.GT.NISMAX)CALL ERROR(210)
   WRITE(NO,1)NIS
   1 FORMAT(1H0,'NUMBER OF ISOTOPES IN FILE 2,151',I6)
   IF(NIS.EQ.1)GOTO 450
   IF(NIS.NE.NIS0)CALL ERROR(219)
C
C-----LCCP ON ISOTOPES
C
450 DC 103 MIS=1,NIS
   CALL RREC(1,NT,MODE,-1.0)
C   (Z,A)DESCRIPTION FOR AN ISOTOPE IN A MATERIAL
   ZAI=C1
C ABUNDANCE(WEIGHT FRACTION) OF AN ISOTOPE IN THIS MATERIAL
   ABN=C2
   ABIS(MIS)=ABN
C IF=1 FISSION WIDTHS ARE GIVEN
   LFW=L2
C NUMBER OF ENERGY RANGES FOR THIS ISOTOPE
   NER=N1
   IF(LRP.GT.0) GOTO 251
250 WRITE(NO,252)
252 FORMAT(1H0,'NO RESONANCE INFORMATION FOR THIS ISOTOPE')
   GOTO 7
251 IF(NER.EQ.1) WRITE(NO,5)
   5 FORMAT(1H0,'ONLY ONE ENERGY RANGE FOR RESONANCE PARAMETERS')
   IF(NER.LE.2) GOTO 7
   WRITE(NO,6)
   6 FORMAT(1H0,'MORE THAN TWO ENERGY RANGES---FATAL ERROR...')
   STCP
C
C-----LOOP ON ENERGY RANGES
C
7 LISRES=1
  LISUNR=1
  DO 2 I=1,NER
    JRES=JRES+1
    CALL RREC(1,NT,MODE,-1.0)
    EL=C1
    EF=C2
C====LRU=0 NO INFORMATION DATA FOR THE CORRESPONDING ENERGY RANGE
C   =1 RESOLVED RESONANCES
C   =2 UNRESOLVED RESONANCES
  LRU=L1
  LRF=L2

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240   IF(LRF.LE.2)GOTO 50
250   IF(LRF.EQ.3)WRITE(NO,30)
260   30 FORMAT(1H0,'REICH-MOORE MULTILEVEL PARAMETERS')
270   IF(LRF.EQ.4)WRITE(NO,40)
280   40 FORMAT(1H0,'ADLER-ADLER FORMALISM')
290   STOP
300   LRU=LRU+1
310   GOTO(10,3,4),LRU
320 C NO INFORMATION
330   10 CALL RREC(1,NT,MODE,-1.0)
340   SPI=C1
350   AF=C2
360   AWR=AWR
370   LISRES=1
380   LISUNR=1
390   GOTO 2
400 C RESONANCES ARE RESOLVED
410   3 LISRES=0
420   JXRES=JX/200
430   JXEV=10*JXRES
440   JXRED=JX-JXEV
450   CALL MARY(EL,EH,LRF,ZAI,ABN,LFWRES,NIS,LNER,MIS,LIS1,SPI,AWRI,AP,
460   *BK,NBT,JNT,X,Y,B,AH(1+JXEV),JXRED,AH(1),AH(1+JXRES),AH(1+2*JXRES),
470   *AH(1+3*JXRES),AH(1+4*JXRES),AH(1+5*JXRES),AH(1+6*JXRES),
480   *AH(1+7*JXRES),AH(1+8*JXRES),AH(1+9*JXRES),JXRES)
490   LISRES=LIS1
500   GOTO 2
510 C RESONANCES ARE NOT RESOLVED
520   4 LISUNR=0
530   NCTM6=6
540   NPEMAX=JX/45
550   CALL ANNICK(EL,EH,LRF,ZAI,ABN,LFW,NIS,MIS,LIS2,SPI,AWRI,AP,BK,
560   *NBT,JNT,X,Y,B,AH(1),AH(1+NPEMAX),AH(1+2*NPEMAX),AH(1+3*NPEMAX),
570   *AH(1+4*NPEMAX),AH(1+5*NPEMAX),AH(1+6*NPEMAX),AH(1+7*NPEMAX),
580   *AH(1+8*NPEMAX),AH(1+9*NPEMAX),AH(1+15*NPEMAX),AH(1+21*NPEMAX),
590   *AH(1+27*NPEMAX),AH(1+33*NPEMAX),AH(1+39*NPEMAX))
600   LISUNR=LIS2
610   2 CONTINUE
620   A=AMOD(ZAI,1000.)
630   AWRIS0(MIS)=A
640 C
650 C-----ISCT1 AND ISOT2
660 C
670   WRITE(NO,302)
680   302 FORMAT('1')
690   CALL XTEXT(4HISOT,NAMI(1))
700   CALL XTEXT(4H,NAMI(2))
710   CALL A9FORM(NO,NAMB)
720   IF(NIS.GT.1.AND.MIS.EQ.1)GOTO 60
730   IBCL=1
740   GOTO 61
750   60 IPCL=2
760   WRITE(NO,210)
770   210 FORMAT(1H0,'ISOT1 AND ISOT2 APPEAR TWO TIMES'/
780   * 1X,' -FOR THE WHOLE MATERIAL'/

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* 1X, ' -FOR THE FIRST ISOTOPE '
61 DC 103 IJ=1,IBCL
Z=FIX(ZAI/1000.0)
AWRI=AWRI
IF(NIS.GT.1.AND.MIS.EQ.1.AND.IJ.EQ.1)AWRI=AWR
ANEUT=1.0086654
BK(1)=AWRI*ANEUT
BK(2)=Z
BK(3)=SPI
N=3
WRITE(NO,102)
102 FORMAT(1H0,' ISOT1'/1X,5('*')//)
WRITE(NO,104)(BK(KB),KB=1,3)
104 FORMAT(1X,' ATOMIC WEIGHT',T35,1PE15.6/
* 1X,' ATOMIC NUMBER',T35,1PE15.6/
* 1X,' NUCLEAR SPIN OF GROUND STATE',T35,1PE15.6/)
MATK=MATKDK
IF(NIS-1)62,62,63
62 MMIS=1
GOTO 64
63 MMIS=MIS
IF(MIS.GT.1.OR.IJ.EQ.2)GOTO 108
GOTO 64
108 MATK=MATISO(MIS)
EB=ERISO(MIS)
MMIS=MIS+1
64 CALL WDA2(MATK ,MMIS,4HISOT,4H1 )
WL=455.197*(AWRI+1.0)/AWRI
WRITE(NO,105)
105 FORMAT(1H0,' ISOT2'/1X,5('*')//)
WRITE(NO,106)WL,AP,EB
106 FORMAT(1X,' REDUCED NEUTRON WAVE LENGTH',T35,1PE15.6,3X,
* '10**(-12)CM'/
* 1X,' NUCLEAR RADIUS',T35,1PE15.6,3X,
* '10**(-12)CM'/
* 1X,' BINDING ENERGY OF THE LAST NEUTRON',T38,1PE12.6,3X,
* 'EV'/)
BK(1)=WL
BK(2)=AP
BK(3)=EB
N=3
CALL WDA2(MATK ,MMIS,4HISOT,4H2 )
AWRI =AWRI
103 CONTINUE
C
C-----ISOT3
C
IF(NIS.EQ.1) GOTO 402
CALL XTEXT(4HISOT,NAM1(1))
CALL XTEXT(4H3 ,NAM1(2))
CALL A9FORM(NO,NAM8)
WRITE(NO,206)
206 FORMAT(1H0,' ISOT3'/1X,5('*')//)
WRITE(NO,107)(AWRISO(MIS),ABIS(MIS),MIS=1,NIS)
107 FORMAT(1X,' ATOMIC WEIGHTS',T25,' ISOTOPIC ABUNCEANCE'//)

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* (3X,1PE12.6,T26,1PE12.6))
I=1
DC 401 MIS=1,NIS
BK(I)=AWRISO(MIS)
BK(I+1)=ABIS(MIS)*100.0
I=I+2
IF(I.GT.NBKX) CALL ERROR(204)
401 CCNTINUE
N=I-1
CALL WDA2(MATKDK,1,4HISOT,4H3 )
GOTO 402
402 IF((MATKDK/1000)*1000.NE.MATKDK) GOTO 400
IAWR=AWRISO(1)+0.00001
WRITE(NO,109) NAMKDK,ABIS(1),LNAM(1),LNAM(2),IAWR,(LNAM(I),I=6,8)
109 FORMAT('OTHE MATERIAL ',A8,' CONSISTS OF ',2PF7.2,'% ',2A1,
*PI3,3A1,'.')
400 RETURN
END
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SUBROUTINE MARY(EL,EH,LRP,ZAI,ABN,LFWRES,NIS,LNER,MIS,LIS,SPI,AWRI
* ,AP,BK,NBT,JNT,ENS,SIGS,B,AH,JXRED,ER,AJ,GT,GN,GG,GF,AL,G,SER,PER,
*NDIM)
COMMON/INOUT/NO,NI,NT,MODE,LEDIT1,LEDIT2,NREF1(78),
*NAMISO(10),NREF2(22),N1X,N2X,N,NBKX,NREF4(21),MLSLSW
COMMON/RECS/NREF3(3),C1,C2,L1,L2,N1,N2
REAL*8 NAMISO
DIMENSION ER(1),AJ(1),GT(1),GN(1),GG(1),GF(1),AL(1),
* G(1),BK(1),NBT(1),JNT(1),B(1),SER(1),FER(1),
*SIGS(1),ENS(1),LLSTR(5),AH(1)
PI=3.14159265
PI4=4.0*PI
CON1=2.196771E-3
CON2=PI4/(CON1**2)
CALL RREC(1,NT,MODE,-1.0)
SPI=C1
AP=C2
C-----LIS=0 CALCULATE CROSS SECTION FROM RESONANCE PARAMETERS
C PLUS SMOOTH CONTRIBUTION FROM FILE 3
C =1 USE SMOOTH CROSS SECTION IN FILE 3 ONLY
LIS=L1
NLS=N1
WRITE(NO,1)
1 FORMAT(1H1,20X,'RESOLVED RESONANCE DATA')
IF(NIS.GT.1)WRITE(NO,360)NAMISO(MIS)
360 FORMAT('+',60X,'ISOTOPE',3X,A8)
CALL PRIZEI
IZ=0.001*ZAI
IA=ZAI-1000.*IZ
WRITE(NO,2)IZ,IA,ABN,EL,EH,SPI,AP,NLS
2 FORMAT(1H0,'Z-----',I11/
* 1X,'A-----',I11/
* 1X,'FRACTIONAL ABUNDANCE-----',1PE11.4/

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*          LX,'LOWER ENERGY LIMIT-----',E11.4/      24C
*          LX,'UPPER ENERGY LIMIT-----',E11.4/      350
*          LX,'NUCLEAR SPIN-----',E11.4/            36C
*          LX,'SPIN SCATTERING LENGTH (A+)--',E11.4/   37C
*          LX,'NUMBER OF L STATES-----',I11/         380
WRITE(NO,1000)                                          39C
1000 FORMAT(' FORMALISM RECOMMENDED BY ENDF/B TO CALCULATE THE RESONANC
*ES:')
IF(LRF,EQ.1) WRITE(NO,1001)
IF(LRF,EQ.2) WRITE(NO,1002)
1001 FORMAT('+',60X,' SINGLE LEVEL BREIT WIGNER.')
1002 FORMAT('+',60X,' MULTI LEVEL BREIT WIGNER.')
WRITE(NO,1003)
1003 FORMAT(' FORMALISM USED IN CALCULATING THE RESONANCES:')
IF(MLSLSW,GE.1) LRF=1
IF(LRF,EQ.1) WRITE(NO,1001)
IF(LRF,EQ.2) WRITE(NO,1002)
IF(NLS,GT.5) CALL ERROR(207)
NRSI=1
DC 3 I=1,NLS
J=1
CALL RREC(2,NT,MODE,-1.0,C,B)
AWRI=C1
C=2.196771E-3*AP*AWRI/(AWRI+1.0)
AM=C2
ARAT=AWRI/(AWRI+1.0)
RAD=0.1*(0.9+1.23*AWRI**0.33333333)
CCN3=CON1*ARAT*RAD
PHO=CON3*SQRT(E)
CCN4=CON2/(ARAT**2)
C          4PI/K**2 = CON4/E
C          RHOCF = C * SQRT(E)
L=L1
LLSTR(I)=L
NRS=N2
NRSF=NRSI+N2-1
WRITE(NO,5) L,NRS,AM
5 FORMAT(1H0,' L VALUE-----',I11/
*          LX,'NUMBER OF RESONANCES-----',I11/
*          LX,'SPIN SCATTERING LENGTH-----',1FE11.4//)
IF(LEDIT2,EQ.1) WRITE(NO,120)
120 FORMAT( 44X,'RESONANCE WIDTHS (EV)'/
*          6X,'ENERGY (EV)',2X,'J VALUE',8X,
*          'TOTAL',17X,'NEUTRON',8X,'RADIATION',6X,
*          'FISSION',8X,'L VALUE',8X,'G VALUE')
DC 4 K=NRSI,NRSF
IF(K,GT,NDIM) CALL ERROR(205)
ER(K)=B(J)
AJ(K)=B(J+1)
GT(K)=B(J+2)
GN(K)=B(J+3)
GG(K)=B(J+4)
GF(K)=B(J+5)
AL(K)=L
G(K)=(2.*AJ(K)+1.)/(2.*(2.*SPI+1.))
C          CALCULATE PENETRATION AND SHIFT FACTOR AT RESONANCE ENERGY ER(K).      89C
C          PHO=CON3*SQRT(ABS(ER(K)))
C          CALL FACTS(L,RHO,SER(K),PER(K))
C          J=J+6
C          IF(LEDIT2,EQ.1)
*          WRITE(NO,6) ER(K),AJ(K),GT(K),GN(K),GG(K),GF(K),AL(K),G(K)
6          FORMAT(1P8E15.4)
4          CONTINUE
NRSI=NRSF+1
3          CONTINUE
C-----NUMBER OF RESONANCES
NRES=NRSF
1000
C-----SORT RESONANCE DATA ON INCREASING ENERGY
IF(NLS,GT.1) CALL SORTRE(NRES,ER,AJ,GT,GN,GG,GF,AL,G,SER,PER)
C-----STORE RESONANCE DATA IN THE ARRAY BK.
CALL MARYD1(EL,EH,NRES,BK,ER,AL,AJ,G,GT,GN,GG,GF,NIS,MIS,LIS,
*          LFWRES,89)
C=====CALCULATE ENERGY POINTS USING NPTR POINTS PER RESONANCE
140 IF(LNER,LE.0) NPTR=51
IF(LNER,GT.1) NPTR=LNER
NPTR=MIN0(NPTR,200)
EUP=EH
CALL EPT(EL,EUP,NPTR,NSE,NRES,ER,GT,ENS,N2X)
C          CHECK IF THERE ARE POINTS WITH THE SAME VALUE.
C          IF YES, PUT ONLY ONE OF THIS POINTS.
CALL MARYD2(ENS,NSE)
WRITE(NO,300) NSE,(ENS(I),I=1,NSE)
300 FORMAT(1H1,I6,3X,' ENERGIES SELECTED BY EPT'// (1P8E15.5))
CALL PRIZEI
C          CALCULATE SGN, SGG AND IF LFWRES=1 SGF.
CALL MARGEL(LFWRES,LRF,C,CON3,CON4,PI*4,AP,NRES,NLS,
*          LLSTR,ER,GN,GG,GF,G,AL,SER,PER)
IF((LFWRES,EQ.1.AND.3*NSE,GT,JXRED).OR.
*          (LFWRES,NE.1.AND.2*NSE,GT,JXRED)) CALL ERROR(208)
DC 40 K=1,NSE
CALL MARGOT(ENS(K),AH(K),AH(NSE+K),SGF)
IF(LFWRES,EQ.1) AH(2*NSE+K)=SGF
40          CONTINUE
C          CHECK IF THERE ARE ENOUGH ENERGY POINTS FOR DESCRIBING
C          POINT RESONANCE CROSS SECTIONS. CALCULATE NBT AND JNT.
IF((6*NIX).GT,NBXX) CALL ERROR(206)
CALL MILLI(ENS,NSE,AH,BK(1),BK(1+NIX),BK(1+2*NIX),BK(1+3*NIX),
*          *BK(1+4*NIX),BK(1+5*NIX),LFWRES,NISGN,NISGG,NISGF,JXRED)
CALL MELA(LFWRES,BK,AH,NBT,JNT,SIGS,NISGN,NISGG,NISGF,NSE,MIS,
*          *EL,EH)
9          RETURN
END

```

C	SUBROUTINE SORTRE(NRES,ER,AJ,GT,GN,GG,GF,AL,G,SER,PER)	10	ENTRY MARGOT(E,SGN,SGG,SGF)	100
	SORTIERT RESONANZDATEN NACH AUFSTIEGENDEN ENERGIEN.	20	SGN8=0.0	110
	DIMENSION ER(1),AJ(1),GT(1),GN(1),GG(1),GF(1),AL(1),G(1),	20	SGG8=0.0	120
	*SER(1),PER(1)	40	SGF8=0.0	130
	NRI=NRES-1	50	RHO CF=C*SQRT(E)	140
	DO 8 IR=1,NRI	60	DO 50 LL=1,NLS	150
	IX=IR+1	70	LLS=LLSTR(LL)	160
	DO 8 I=IX,NRES	80	CALL FACPHI(LLS,RHO CF,PHIL)	170
	IF(ER(IR).LT.ER(I)) GOTO 9	90	LLSI=LLS+1	180
	T1=ER(IR)	100	SINP2(LLSI)=SIN(PHIL)**2	190
	T2=AJ(IR)	110	SIN2P(LLSI)=SIN(2.0*PHIL)	200
	T3=GT(IR)	120	CCS2P(LLSI)=CCS(2.0*PHIL)	210
	T4=GN(IR)	130	SPDT=0.0	220
	T5=GG(IR)	140	IF(ABS(PHIL).GT.5.0E-4.OR.LLS.GT.0)	230
	T6=GF(IR)	150	*SPDT=CON4*(2.0*FLOAT(LLS)+1.0)*SINP2(LLSI)/E	240
	T7=AL(IR)	160	IF(ABS(PHIL).LE.5.0E-4.AND.LLS.EQ.0)	250
	T8=G(IR)	170	*SPDT=PI*4*AP**2	260
	T9=SER(IR)	180	SGN8=SGN8+SPDT	270
	T10=PER(IR)	190	50 CONTINUE	280
	ER(IR)=ER(I)	200	CON5=CON4/E	290
	AJ(IR)=AJ(I)	210	RHCFE=CON3*SQRT(E)	300
	GT(IR)=GT(I)	220	DO 54 J=1,NRES	310
	GN(IR)=GN(I)	230	LLS=AL(J)+0.001	320
	GG(IR)=GG(I)	240	LLSI=LLS+1	330
	GF(IR)=GF(I)	250	IF(LLS) 55,56,57	340
	AL(IR)=AL(I)	260	55 CALL ERROR(209)	350
	G(IR)=G(I)	270	56 ERE=ER(J)	360
	SER(IR)=SER(I)	280	GN1=GN(J)*SQRT(E/ABS(ER(J)))	370
	PER(IR)=PER(I)	290	GCTO 22	380
	ER(I)=T1	300	57 CALL FACTS(LLS,RHO E,SLE,PLE)	390
	AJ(I)=T2	310	ERE=ER(J)+(SER(J)-SLE)*GN(J)/(2.0*PER(J))	400
	GT(I)=T3	320	GN1=GN(J)*PLE/PER(J)	410
	GN(I)=T4	330	22 GX=GF(J)+GG(J)	420
	GG(I)=T5	340	GT1=GN1+GX	430
	GF(I)=T6	350	XZ=2.0*(E-ERE)/GT1	440
	AL(I)=T7	360	SIG0=CON5*G(J)	450
	G(I)=T8	370	GN1GT1=GN1/3T1	460
	SER(I)=T9	380	CCMFAC=SIG0*GN1GT1/(1.0+XZ**2)	470
	PER(I)=T10	390	SIG1=CCMFAC*(GN1GT1*CCS2P(LLSI)-2.0*SINP2(LLSI)*GX/GT1	480
8	CONTINUE	400	*+XZ*SIN2P(LLSI))	490
	RETURN	410	SGN8=SGN8+SIG1	500
	EAD	420	SIGMA=CCMFAC/GT1	510
			SGG8=SGG8+SIGMA*GG(J)	520
			IF(LFWRES.EQ.1) SGF8=SGF8+SIGMA*GF(J)	530
			IF(LRF.NE.2) GOTO 54	540
			MULTILEVEL INTERFERENCE	550
			IF(J.EQ.1) GOTO 54	560
			JJ=J-1	570
			SML8=0.0	580
			DO 58 JS=1,JJ	590
C	SUBROUTINE MARCEL(LFWRES,LRF,C,CON3,CON4,PI*4,AP,NRES,NLS,	30	CHFK L-VALUE AND G-VALUE TO CALCULATE INTERFERENCE	600
	*LLSTR,FR,GN,GG,GF,G,AL,SER,PER)	40	ONLY FOR SAME SPIN STATES.	610
C	CALCULATES ELASTIC, RADIATIVE CAPTURE AND (IF LFWRES=1) FISSION	50	IF(G(JS).NE.G(J).OR.AL(JS).NE.AL(J)) GCTO 58	620
C	CROSS SECTIONS. DEPENDING ON LRF SINGLE LEVEL OR MULTI LEVEL	60	IF(LLS) 59,60,61	630
	BREIT WIGNER IS USED.	70	59 CALL ERROR(209)	640
	REAL*8 SGN8,SGG8,SGF8,SML8	80		
	DIMENSION LLSTR(5),SIN2P(5),SINP(5),COS2P(5),AL(1),ER(1),	90		
	*GN(1),GG(1),GF(1),G(1),SER(1),PER(1)			
	GCTO 333			

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60 ERES=ER(JS)
GNS=GN(JS)*SQRT(E/ABS(ER(JS)))
GOTO 23
61 ERES=ER(JS)+(SER(JS)-SLE)*GN(JS)/(2.0*PER(JS))
GNS=GN(JS)*PLE/PER(JS)
23 GTS=GNS+GF(JS)+GG(JS)
XS=2.0*(E-ERES)/GTS
SMLB=SMLB+GNS/GTS*(1.0+XZ*XS)/(1.0+XS**2)
58 CONTINUE
SGNB=SGNB+2.0*SMLB*COMFAC
54 CONTINUE
SGN=SGNB
SGG=SGGB
SGF=SGFB
333 RETURN
END

SUBROUTINE MILLI(ENS,NSE,AH,NBTN,JNTN,NBTG,JNTG,NBTF,JNTF,
*LFWRES,NISGN,NISGG,NISGF,JXRED)
CHECKS IF THERE ARE ENOUGH POINTS DESCRIBING POINT RESONANCE
CROSS SECTIONS. CALCULATES NBT AND JNT.
NSE MAY HAVE BEEN ENLARGED AFTER RETURN. IN NBT AND JNT THERE ARE
AFTER RETURN NI VALUES FOR INTERPOLATING THE POINT CROSS SECTION.
N MEANS SGN, G MEANS SGG AND F MEANS SGF.
IN AH YOU WILL FIND THE FOLLOWING DATA.
1 ... NSE SGN
NSE+1 ... 2*NSE SGG
2*NSE+1 ... 3*NSE SGF

COMMON/INOUT/ IOU, NR1(126), N2X, NR2(12), EPS
DIMENSION ENS(1), AH(1), NBTN(1), JNTN(1), NBTG(1), JNTG(1),
*NBTF(1), JNTF(1), EPSI(5), NIC(5), CST(3)
DATA CST/'SGN ','SGG ','SGF '/
LOGICAL*4 T
T=.FALSE.
IF(LFWRES.EQ.1) T=.TRUE.
IEND=2
IF(T) IEND=3
NISGN=1
NISGG=1
NISGF=1
NBTN(1)=2
JNTN(1)=2
NBTG(1)=2
JNTG(1)=2
IF(.NOT.T) GOTO 17
NBTF(1)=2
JNTF(1)=2
C LOOP FOR ALL ENERGY-INTERVALS.
17 K=2
1 E1=ENS(K-1)
E2=ENS(K)

E=(E2+E1)*.5
IF(.LT.E1) E=0.5*E2+0.5*E1
CALL MARROT(E,SGN,SGG,SGF)
CALCULATE FOR SGN,SGG AND SGF INTERPOLATION CODE AND EPSI
EPSSGF=0.0
DO 2 I=1,IEND
GOTO (3,4,5),I
3 IP=0
SIGMA=SGN
ICALT=JNTN(NISGN)
GOTO 6
4 IP=NSE
SIGMA=SGG
ICALT=JNTG(NISGG)
GOTO 6
5 IP=2*NSE
SIGMA=SGF
ICALT=JNTF(NISGF)
6 S1=AH(IP+K-1)
S2=AH(IP+K)
DO 7 IC=2,5
IF((IC.GE.4.AND.(S1.LE.C.O.OR.S2.LE.O.C)).OR.
*((IC.EQ.3.OR.IC.EQ.5).AND.((E2/E1.EQ.1.0).OR.
*(E.LT.E1.OR.F2.LT.E1)))) GOTO 23
CALL TERP1(E1,S1,E2,S2,E,YINT,IC)
EPSI(IC)=ABS((YINT-SIGMA)/SIGMA)
GOTO 7
23 EPSI(IC)=1.0E+50
7 CONTINUE
INCREASING ORDER OF EPSI IS MADE.
DO 18 IC=2,5
18 NIC(IC)=IC
DO 19 IC=2,4
ICP1=IC+1
DO 20 ICI=ICP1,5
IF(EPSI(IC).LE.EPSI(ICI)) GOTO 20
IP=NIC(IC)
NIC(IC)=NIC(ICI)
NIC(ICI)=IP
SIGMA=EPSI(IC)
EPSI(IC)=EPSI(ICI)
EPSI(ICI)=SIGMA
20 CONTINUE
19 CONTINUE
IF(EPSI(2).GT.EPS.AND.E.NE.E1) GOTO 11
IF(EPSI(2).GT.EPS.AND.E.EQ.E1) WRITE(IOU,1000) EPS,E1,E2,CST(I),
*EPSI(2)
1000 FERMAT('0***** WARNING IN SUBROUTINE MILLI. WISHED ACCURACY=',
*E5.4,' FOR INTERPOLATING POINTS'/
*' ***** IN THE ENERGY INTERVAL E1=',1PE15.7,' E2=',1PE15.7,
*' COULD NOT BE FULFILLED FOR'/
*' ***** CROSS-SECTION TYPE ',A4,'. BEST ACCURACY FOR THIS TYPE IN
* THIS INTERVAL IS=',1PE15.7,'.')
III=2
IF(K.EQ.2) GOTO 21

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DC 22 II=2,5
IF(NIC(II).NE.ICALT) GOTO 22
IF(II.EQ.2) GOTO 21
IF(EPSI(II).LE.EPS.AND.EPSI(II).LE.(6*EPSI(2)/II)) III=II
GOTO 21
22 CCNTINUE
21 ICODE=NIC(III)
GOTO (8,9,10),I
8 ICSGN=ICODE
GOTO 2
9 ICSGG=ICODE
GOTO 2
10 ICSGF=ICODE
2 CONTINUE
C EPSSGN,EPSSGG AND EPSSGF ARE LESS EPS. STORE INTERPOLATIONCODE
C SET K=K+1 AND BRANCH TO 1 IF(K.LE.NSE).
C FIRST INTERVAL?
IF(K.GT.2) GOTO 12
JNTN(1)=ICSGN
JNTG(1)=ICSGG
IF(.NOT.T) JNTF(1)=ICSGF
GOTO 13
12 CONTINUE
CALL MILLI1(NBTN,JNTN,NISGN,ICSGN,K,NSE)
CALL MILLI1(NBTG,JNTG,NISGG,ICSGG,K,NSE)
IF(T) CALL MILLI1(NBTF,JNTF,NISGF,ICSGF,K,NSE)
13 K=K+1
IF(K.LE.NSE) GOTO 1
GOTO 333
11 CCNTINUE
ENLARGE ENS BY POINT E. ENLARGE AH BY SIGMAS.
NSE=NSE+1
IF(NSE.GT.N2X) CALL RESERR(105,K,NSE,1)
IF((T.AND.(3*NSE).GT.JXRED).OR.
* (.NOT.T.AND.(2*NSE).GT.JXRED)) CALL ERROR(208)
C MAKING A GAP.
KK=IEND*NSE
IVER=IEND
IM=IEND-1
14 I1=KK-IVER
AH(KK)=AH(I1)
IF(KK.LE.NSE) ENS(KK)=ENS(I1)
15 KK=KK-1
IF(KK.NE.(I4*NSE+K)) GOTO 14
IF(IM.LE.0) GOTO 16
IVER=IVER-1
IM=IM-1
IF(IVER.LT.0) CALL ERRCR(777)
GOTO 15
16 ENS(K)=E
AH(K)=SGN
AH(NSE+K)=SGG
IF(T) AH(2*NSE+K)=SGF
GOTO 1
333 RETURN
END

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1010
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1100
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1120
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1170
1180
1190
1200
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SUBROUTINE MILLI1(NBT,JNT,N1,IC,K,NSE)
COMMON/INOUT/ NR1(126),N1X
DIMENSION JNT(1),NBT(1)
IF(JNT(N1).EQ.TC) GOTO 1
N1=N1+1
IF(N1.GT.N1X) CALL RESERR(103,K,NSE,1)
JNT(N1)=IC
1 NBT(N1)=K
333 RETURN
END
10
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100

SUBROUTINE RESERR(N,K,NSE,IFERR)
COMMON/INOUT/ IC
WRITE(ID,100)
100 FORMAT(' ***** ERROR IN RESOLVED RESONANCE CALCULATING SECTION.')
IF(IFERR.EQ.1) WRITE(ID,102) K,NSE
102 FORMAT(' ***** ',15,' POINTS OUT OF ',15,' PROCESSED.')
IF(IFERR.EQ.2) WRITE(ID,103) K,NSE
103 FORMAT(' ***** ',15,' RESONANCES OUT OF ',15,' PROCESSED.')
I=0
IF(N.EQ.103) I=1
IF(N.EQ.105) I=2
WRITE(ID,101) I
101 FORMAT(' ***** INCREASE N',11,'X OR TOTAL REGION.')
STOP 2
END
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SUBROUTINE MELA(LFWRES,IBK,AH,NBT,JNT,SIGS,NISGN,NISGG,NISGF,NSE,
*MIS,FL,EH)
COMMON/INOUT/ NO,NR1(5),LEDTIS,NR2(119),N1X,N2X
COMMON/RECS/ NR3(3),C1,NR4(3),N1,N2
DIMENSION IBK(1),AH(1),NBT(1),JNT(1),SIGS(1),TK(3)
REAL*8 T(6)
DATA T/'SCATTERI','NG','CAPTURE',' ',' ',' '
*'FISSION',' ','/','TK/'SCAT','CAPT','FISS'/'
K3=2
IF(LFWRES.EQ.1) K3=3
C1=0.0
DO 77 K2=1,K3
GOTO (70,71,72),K2
70 N1=NISGN
IP=0
GOTO 73
71 N1=NISGG
IP=2*N1X
10
20
30
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GOTO 73
72 N1=N1SGF
   IF=4*N1X
73 N2=NSE
   DC 74 I=1,N1
   NBT(I)=IBK(IP+I)
74 JAT(I)=IBK(IP+N1X+I)
   IF(N2.GT.N2X) CALL ERROR(105)
   DC 75 I=1,N2
75 SIGS(I)=AH((K2-1)*NSE+I)
   IF(LEDITS.NE.1) GOTO 101
   WRITE(NO,250) T(2*K2-1),T(2*K2)
250 FORMAT(' POINT WISE ',A8,A3,' CROSS SECTION CALCULATED FROM RESOLVE
 *D RESONANCE PARAMETER///')
   CALL WREC(3,NO,4)
101 CONTINUE
   CALL WRECS2(MIS,TK(K2),4H RES,EL,EH)
77 CONTINUE
   RETURN
   END

SUBROUTINE MARY01(EL,EH,NRES,BK,ER,AL,AJ,G,GT,GN,GG,GF,NIS,MIS,
 *LIS,LEWRES,*)
  DIMENSION NAM1(2),BK(1),ER(1),AL(1),AJ(1),G(1),GT(1),GN(1),GG(1),
 *GF(1)
  REAL *8 NAMISO,NAM8
  COMMON/INOUT/NO,NR1(4),LEDIT2,NR2(76),MATKDK,NR3(1),NAMISO(10),
 *MATISO(10),NR4(14),N,NBKK
  EQUIVALENCE (NAM8,NAM1(1))

  J=1
  DC 17 K=1,NRES
  IF(I.GT.NBKK) CALL ERROR(206)
  BK(J)=ER(K)
  BK(J+1)=AL(K)
  BK(J+2)=AJ(K)
  BK(J+3)=G(K)
  BK(J+4)=GT(K)
  BK(J+5)=GN(K)
  IF(ER(K).LT.0.0) BK(J+5)=GN(K)/SQRT(ABS(ER(K)))
  BK(J+6)=GG(K)
  BK(J+7)=GF(K)
  BK(J+8)=0.0
  BK(J+9)=0.0
  BK(J+10)=0.0
  J=J+11
17 CONTINUE
  N=J-1
  IF(LEDIT2.NE.1) GOTO 100
  WRITE(NO,21)
21 FORMAT(1H1,' RESONANCE DATA IN KEDAK FORMAT')
  IF(NIS.GT.1) WRITE(NO,360)NAMISO(MIS)

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19C
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210
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23C
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25C
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27C
28C
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30C
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35C
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38C
1C
20
3C
4C
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6C
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11C
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13C
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16C
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200
21C
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230
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26C
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300
31C

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360 FORMAT(1H+,60X,' ISOTOPE ',3X,A8)
   CALL XTEXT(4HRES ,NAM1(1))
   CALL XTEXT(4H ,NAM1(2))
   CALL ABFORM(NC,NAM8)
   CALL LUCY(2,BK)
100 CONTINUE
   WRITE(NO,350)
350 FORMAT(1H1,' KEDAK REACTION TYPE RANGRES')
   IF(NIS.GT.1) WRITE(NO,360)NAMISO(MIS)
   CALL XTEXT(4HRANG,NAM1(1))
   CALL XTEXT(4HRES ,NAM1(2))
   CALL ABFORM(NC,NAM8)
   MMIS=1
   MATK=MATKDK
   IF(NIS.EQ.1) GOTO 301
   MATK=MATISO(MIS)
   MMIS=MIS+1
301 CALL WDA2(MATK ,MMIS,4HRES ,4H )
C====RANGRES
  BK(1)=EL
  BK(2)=EH
  BK(3)=NRES
  BK(4)=1.0
  N=4
  WRITE(NO,351)(BK(K),K=1,4)
351 FORMAT(' LOWER ENERGY BOUNDARY OF THE RESOLVED RESONANCE REGION',
 *T58,IPE15.6,' EV'/
 * ' UPPER ENERGY BOUNDARY OF THE RESOLVED RESONANCE REGION',
 *T58,IPE15.6,' EV'/
 * 1X,' NUMBER OF RESONANCES',T58,CPF15.0/
 * 1X,' CONSISTENCY TEST PARAMETER',T58,F15.0)
  CALL WDA2(MATK,MMIS,4HRANG,4HRES )
  IF(LIS.NE.0) RETURN 1
C====DETERMINE IF THE FISSION WIDTHS ARE ZERO
  10 LEWRES=1
  DC 130 IP=1,NRES
  IF(GF(IP).GT.1.E-6) GOTO 333
130 CONTINUE
  LEWRES=0
  WRITE(NO,150)
150 FORMAT(1H0,' NO FISSION WIDTHS FOR THIS ISOTOPE')
333 RETURN
   END

SUBROUTINE MARY02(ENS,NSE)
  CHECKS IF THERE ARE POINTS WITH THE SAME VALUE.
  IF YES, ALL BUT ONE OF THESE VALUES ARE DELETED.
  DIMENSION EVS(1)
  I=2
  1 IF(ENS(I).EQ.ENS(I-1)) GOTO 2
  I=I+1

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5 IF(I,LE,NSE) GOTO 1
333 RETURN
2 IA=I+1
IF(IA,GT,NSE) GOTO 4
DO 3 II=IA,NSE
3 ENS(II-1)=ENS(II)
4 NSF=NSE-1
GOTO 5
END

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SUBROUTINE FACTS(L,RHO,SE,PE)
C*****
C* CALCULATES PENETRATION AND SHIFT FACTORS *
C* LIBERATED FROM GREGSON ET. AL. *
C*****
LL=L+1
GC TO(10,20,30,40,50),LL
C-----S-WAVE
10 SE=0.0
PE=RHO
GO TO 100
C-----P-WAVE
20 R2=RHO*RHO
DEN=1.0+R2
PE=R2*RHO/DEN
SE=-1.0/DEN
GO TO 100
C-----D-WAVE (L=2)
30 R2=RHO*RHO
R4=R2*R2
DEN=3.0*R2+R4+9.0
PE=R4*RHO/DEN
SE=-{(18.0+3.0*R2)/DEN}
GO TO 100
C-----F-WAVE (L=3)
40 R2=RHO*RHO
R4=R2*R2
R6=R4*R2
DEN=225.0+45.0*R2+6.0*R4+R6
PE=R6*RHO/DEN
SE=-{(675.0+90.0*R2+6.0*R4)/DEN}
GO TO 100
C-----G-WAVE (L=4)
50 R2=RHO*RHO
R4=R2*R2
R6=R4*R2
R8=R4*R4
DEN=11025.0+1575.0*R2+135.0*R4+10.0*R6+R8
PE=R8*RHO/DEN
SE=-{(44100.0+4725.0*R2+270.0*R4+10.0*R6)/DEN}
100 RETURN
END

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SUBROUTINE FACPHI(L,RHO,PHI)
C*****
C* CALCULATES PHASE SHIFT *
C*****
LL=L+1
GC TO (10,20,30,40,50),LL
C-----S-WAVE
10 PHI=RHO
GC TO 100
20 PHI=RHO-ATAN(RHO)
GC TO 100
30 R2=RHO*RHO
PHI=RHO-ATAN(3.0*RHO/(3.0-R2))
32 IF((PHI/RHO).LT.0.000001) PHI=0.0
GC TO 100
40 R2=RHO*RHO
PHI=RHO-ATAN((15.0*RHO-RHO*R2)/(15.0-6.0*R2))
GC TO 32
50 R2=RHO*RHO
R4=R2*R2
TCP=105.0*RHO-10.0*R2*RHO
BOT=105.0-45.0*R2+R4
PHI=RHO-ATAN(TOP/BOT)
GC TO 32
100 RETURN
END

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SUBROUTINE ANNICK(EL,EH,LR,F,ZAI,ABN,LFW,NIS,MIS,LIS,SPI,AWRI,AP,
*BK,NBT,JNT,BK2,Y,B,SIGS,E,SIGG,SIGF,SIGT,ESC,XT,YT,TEMP,
*C,CX,GNO,GG,GF,AJ)
COMMON/INOUT/NO,NI,NT,MCDE,LEDIT1,LEDIT2,NRI(76),
*MATKDK,NTYP,NAMISO(10),MATISO(10),NR2(11),NBX,N1X,N2X,N,NBXX,
*NR3(10),EPS
COMMON/RECS/MAT,MF,MT,C1,C2,L1,L2,N1,N2
REAL*8 NAMISO,NAMB
COMMON/F LCT/ZYN,ZYF,U,V,RHON,RHOV,NDIM6,NPEMAX
DIMENSION SIGS(NPEMAX),E(NPEMAX),SIGG(NPEMAX),SIGF(NPEMAX),
*SIGT(NPEMAX),D(NDIM6,NPEMAX),GX(NDIM6,NPEMAX),
*GNO(NDIM6,NPEMAX),GG(NDIM6,NPEMAX),GF(NDIM6,NPEMAX),
*AMUX(6),AMUN(6),AMUG(6),AMUF(6),AJ(NDIM6,NPEMAX),INTS(18),
*ESC(NPEMAX),XT(NPEMAX),YT(NPEMAX),TEMP(NPEMAX),
*BK2(1),BK(1),NAM1(2),NBT(1),JNT(1),Y(1),B(1)
EQUIVALENCE (NAMB,NAM1(1))
NB=1
NB2=1
INT=0
LS=1
N=1

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ICIMBK=N8KX	220	11 CONTINUE	770
JLI=1	22C	AMUN(J)=B(JS+2)	78C
DO 31 K=1,NPEMAX	240	AMUX(J)=0.0	790
SIGS(K)=0.0	250	AMUG(J)=0.0	80C
SIGG(K)=0.0	260	AMUF(J)=0.0	81C
SIGF(K)=0.0	270	INTS(JLI)=5	820
SIGT(K)=0.0	28C	JLI=JLI+1	83C
31 CONTINUE	290	JS=JS+6	840
WRITE(N0,1)	300	10 CONTINUE	850
1 FORMAT(1H1,30X,'UNRESOLVED RESONANCE DATA')	310	GCTO 12	86C
IF(NIS.GT.1)WRITE(N0,110)NAMISO(MIS)	320	C-----FISSION-WIDTHS GIVEN	870
110 FORMAT(1H+,60X,' ISOTOPE',3X,A8)	33C	C-----ONLY FISSION-WIDTHS ARE ENERGY-DEPENDENT	880
CALL PRIZEI	340	3 IF(LS.GT.1)GOTO 41	89C
CALL INITSC(NPEMAX)	350	LFCHG=1	900
CALL RREC(1,NT,MODE,-1.0)	360	PACKSPACE NT	910
SPI=C1	370	CALL RREC(2,NT,MODE,-1.0,B)	92C
A=C2	38C	NE=N1	930
LIS=L1	390	IF(NE.GT.NPEMAX)CALL ERROR(229)	94C
C	400	NLS=N2	950
C BEGINNING OF LOOP ON L STATES	410	DC 13 K=1,NE	96C
C	420	F(K)=B(K)	97C
39 IF(LRF.EQ.2)GOTO 4	43C	13 CONTINUE	980
IF(LFW)2,2,3	440	ISTGF=1	990
C-----FISSION-WIDTHS NOT GIVEN	450	41 CALL RREC(1,NT,MODE,-1.0,B)	1000
C----- ALL PARAMETERS ARE ENERGY INDEPENDENT	46C	AWRI=C1	1010
2 IF(LS.GT.1)GOTO 40	470	C=2.19685E-3*A*AWRI/(AWRI+1.0)	102C
LFCHG=0	48C	L=L1	1030
NLS=N1	490	NJS=N1	1040
C-----CALCULATES ENERGIES USING EQUAL LETHARGY	500	IF(NJS.GT.6)CALL ERROR(600)	1050
E(1)=EL	510	DO 17 J=1,NJS	1060
NPUNR=20	520	CALL RREC(2,NT,MODE,-1.0,B)	1070
DE=NPUNR-1	53C	DC 18 K=1,NE	1080
DE=ALOG(EH/EL)/DE	540	D(J,K)=B(1)	1090
DE=EXP(DE)	550	AJ(J,K)=B(2)	1100
DO 5 K=2,NPUNR	560	GNO(J,K)=B(4)	111C
E(K)=E(K-1)*DE	570	GG(J,K)=B(5)	1120
5 CONTINUE	58C	GX(J,K)=0.0	113C
E(NPUNR)=EH	590	GF(J,K)=B(K+6)	1140
ISTGF=0	600	18 CONTINUE	1150
40 CALL RREC(2,NT,MODE,-1.0,B)	610	AMUN(J)=B(3)	116C
AWRI=C1	620	AMUX(J)=0.0	1170
C=2.19685E-3*A*AWRI/(AWRI+1.0)	63C	AMUG(J)=0.0	1180
L=L1	640	AMUF(J)=L2	1190
NJS=N2	650	INTS(JLI)=5	1200
IF(NJS.GT.6)CALL ERROR(600)	66C	JLI=JLI+1	121C
JS=1	670	17 CONTINUE	1220
NE=NPUNR	680	GCTO 12	1230
DC 10 J=1,NJS	69C	C-----ALL ENERGY-DEPENDENT PARAMETERS	1240
DO 11 K=1,NE	700	4 IF(LS.GT.1)GOTO 42	1250
D(J,K)=B(JS)	71C	LFCHG=2	126C
AJ(J,K)=B(JS+1)	720	NLS=N1	1270
GNO(J,K)=B(JS+3)	730	ISTGF=1	1280
GG(J,K)=B(JS+4)	740	42 CALL RREC(1,NT,MODE,-1.0)	1290
GF(J,K)=0.0	750	AWRI=C1	1300
GX(J,K)=0.0	76C	C=2.19685E-3*A*AWRI/(AWRI+1.0)	131C

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L=L1
NJS=N1
IF(NJS.GT.6)CALL ERROR(600)
DC 22 J=1,NJS
CALL RREC(2,NT,MODE,-1.0,B)
NE=N2
IF(NE.GT.NPEMAX)CALL ERROR(220)
JK=1
DC 23 K=1,NE
E(K)=B(JK+6)
D(J,K)=B(JK+7)
AJ(J,K)=C1
GNO(J,K)=B(JK+9)
GG(J,K)=B(JK+10)
GX(J,K)=B(JK+8)
GF(J,K)=B(JK+11)
JK=JK+6
23 CONTINUE
AMUN(J)=B(4)
AMUX(J)=B(3)
AMUG(J)=B(5)
AMUF(J)=B(6)
INTS(JLI)=L1
JLI=JLI+1
22 CONTINUE
C
C====PRINTING OF UNRESOLVED RESONANCE DATA
C
12 IF(IWT.GE.1)GOTO 25
WRITE(ND,24)EL,EH,SPI,NLS
24 FORMAT(1H0,'LOWER ENERGY LIMIT-----',1PE11.4/
* 1X,'UPPER ENERGY LIMIT-----',E11.4/
* 1X,'SPIN SCATTERING LENGTH-----',E11.4/
* 1X,'NUMBER OF L STATES-----',I11//)
IWT=IWT+1
25 IF(LEDIT2.NE.1)GOTO 100
WRITE(ND,26)L
26 FORMAT(1H1,10X,'L =',I3/11X,6('='))//)
DC 27 J=1,NJS
JINT=AJ(J,1)
WRITE(ND,28)JINT
28 FORMAT(//' J = ',I3/1X,7('-')//)
WRITE(ND,32) INTS(JLI-1-NJS+J)
32 FORMAT(' INTERPOLATION-CODE=',I1,'.')
WRITE(ND,29)AMUN(J),AMUX(J),AMUG(J),AMUF(J)
29 FORMAT(1X,'NUMBER OF DEGREES OF FREEDOM',
*1X,'NEUTRON WIDTH-----',F5.1/
*30X,'COMPETITIVE WIDTH-',F5.1/
*30X,'RADIATION WIDTH---',F5.1/
*30X,'FISSION WIDTH-----',F5.1//)
WRITE(ND,30)
30 FORMAT(3X,'ENERGY (EV)',6X,'MEAN LEVEL',5X,'AV NEUTRON',5X,
* 'AV RADIAT.',5X,'AV FISSION',5X,'AV COMPET.'/20X,
*'SPACING',10X,4('WIDTH',10X))
WRITE(ND,50)(E(K),D(J,K),GNO(J,K),GG(J,K),GF(J,K),GX(J,K)),

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* K=1,NE)
50 FORMAT(1X,1P6E15.5)
27 CCNTINUE
C
C====CALCULATION OF POINT-WISE CROSS-SECTIONS
C
100 T=C/A
AWR=AWRI
CCNST=19.7392088/(T*T)
PIFOUR=12.566371
ASS=AWR*1.008665
CRT=ASS**(.1.0/3.0)
TERM=0.123*CRT+.08
APEN=TERM
CPEN=C/A*APEN
LI=L+1
DO 33 J=1,NJS
MU=IFIX(AMUN(J))
IF(LFCHG.EQ.0)GOTO 300
N1=1
DO 301 IE=1,LFCHG
DO 302 K=1,NE
XT(K)=E(K)
IF(IE.EQ.2)GOTO 303
YT(K)=GF(J,K)
GOTO 302
303 YT(K)=GNO(J,K)
302 CONTINUE
NBT(1)=NE
JNT(1)=INTS(JLI-1-NJS+J)
N2=NE
CALL CHSINT(EPS,N1,N2,NBT,JNT,XT,YT,N1X,NPEMAX,8304)
GTO 305
304 CALL ERROR(221)
305 CALL GENER(N2,XT,TEMP,ESC,ND,NCU)
IF(NCU.GT.NPEMAX)CALL ERROR(222)
301 CONTINUE
NU=IFIX(AMUF(J))
LAMBDA=IFIX(AMUX(J))
DO 34 K=1,NE
GJ=(2.0*AJ(J,K)+1.0)/(4.0*SPI+2.0)
E2=SQRT(E(K))
EJ=E(K)
WAVE=C/A*E2
RHO=C*E2
RHOPEN=CPEN*E2
GOTO (35,36,37),LI
35 VL=E2*AMUN(J)
PS=RHO
GOTO 38
36 VL=E2*RHOPEN**2/(1.0+RHOPEN**2)*AMUN(J)
PS=RHO-ATAN(RHO)
GOTO 38
37 VL=E2*RHOPEN**4/(RHOPEN**4+3.0*RHOPEN**2+9.0)*AMUN(J)
PS=RHO-ATAN(3.0*RHO/(3.0-RHO**2))

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187C
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1890
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38 SIGPCT=(SIN(PS)/WAVE)**2
SIGPOT=SIGPCT*PIFOUR*(2.0*L+1.0)
GA=GNO(J,K)*VL
GAMMA=GG(J,K)
GALPHA=GN
GBETA=GF(J,K)
DIFF=GX(J,K)
GT=GN+GAMMA+GBETA+DIFF
TERG=(CONST*GJ*GN*GAMMA)/(EJ*D(J,K))
TERF=(CONST*GJ*GN*GN)/(EJ*D(J,K))
TERF=(CONST*GJ*GN*GBETA)/(EJ*D(J,K))
CALL GNRL(GALPHA,GBETA,GAMMA,MU,NU,LAMBDA,GS,DIFF,1)
CALL GNRL(GALPHA,GBETA,GAMMA,MU,NU,LAMBDA,GC,DIFF,2)
CALL GNRL(GALPHA,GBETA,GAMMA,MU,NU,LAMBDA,GFF,DIFF,3)
GS=GS*TERG
GC=GC*TERG
GFF=GFF*TERF
CCRR=(CONST*GJ*2.0*GN*SIN(PS)*SIN(PS))/(EJ*D(J,K))
GS=GS-CORR
SIGG(K)=SIGG(K)+GC
SIGS(K)=SIGS(K)+GS
SIGF(K)=SIGF(K)+GFF
IF(ISTGF.EQ.0)GOTO 80
IF(NB2.GT.N2X) CALL ERRCR(105)
BK2(NB2)=E(K)
BK2(NB2+1)=FLCAT(L)
BK2(NB2+2)=AJ(J,K)
BK2(NB2+3)=AMUF(J)
BK2(NB2+4)=GF(J,K)
BK2(NB2+5)=GG(J,K)
BK2(NB2+6)=GN
ZYN=AMUN(J)
ZYF=AMUF(J)
U=GN/GG(J,K)/AMUN(J)
V=GF(J,K)/GG(J,K)
RHGV=1.
RHON=0.
JINT=AJ(J,1)
IF(V.NE.0.0.AND.ZYF.EQ.0.0) WRITE(ND,1000) L,JINT,K,V
1000 FCRMAT(' ***** WARNING IN SUBROUTINE ANNICK. NUMBER OF DEGREES OF
*FREEDOM FOR FISSION WIDTH*/
*' ***** IN THE UNSOLVED RESONANCE REGION FOR L=',I2,' J=',I2,
*' K=',I3,' IS EQUAL ZERC AND*/
*' ***** V=GF(J,K)/GG(J,K) HAS THE VALUE=',1PE15.7/
*' ***** THE STATISTICAL FLUCTUATION FACTORS ARE SET TO ZERC.')
IF(V.NE.0.0.AND.ZYF.NE.0.0) GOTO 200
PEER=0.
XEER=0.
RG=0.
RF=0.
GOTO 220
200 CALL ERF(PEER)
R+CV=0.
CALL ERF(XEER)
RF=(V*PEER*(ZYN+ZYF+2.)+ZYF*(XEER-1.))/(V*ZYN-L*ZYF)

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242C
2430
2440
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247C
2480
2490
250C
2510
252C
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2540
2550
2560
257C
2580
2590
2600
2610
262C
2630
2640
2650
2660
267C
2680
2690
2700
2710
272C
2730
274C
2750
2760
277C
2780
2790
2800
2810
282C
2830
284C
2850
286C
287C
2880
289C
2900
2910
292C
2930
2940
2950
2960

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RHCN=1.
CALL ERF(RG)
220 BK2(NB2+7)=PEER
BK2(NB2+8)=XEER
BK2(NB2+9)=RF
BK2(NB2+10)=RG
NB2=NB2+11
80 IF(J.GT.1)GOTO 34
SIGS(K)=SIGS(K)+SIGPOT
34 CONTINUE
IF(NR.GT.NBKX) CALL ERRCR(204)
BK(NB)=FLCAT(L)
BK(NB+1)=AJ(J,1)
BK(NB+2)=GG(J,1)
BK(NB+3)=D(J,1)
BK(NB+4)=GNO(J,1)
BK(NB+5)=GNO(J,1)/D(J,1)
BK(NB+6)=AMUF(J)
BK(NB+7)=AMUN(J)
NE=NB+8
33 CONTINUE
LS=LS+1
C
C END OF LOOP ON L STATES
C
IF(NLS.GT.3) CALL ERRCR(601)
IF(LS.LE.NLS)GOTO 39
CALL ANNI01(NE,SIGT,SIGG,SIGS,SIGF,NIS,MIS,E,NB,BK,ISTGF,NB2,
*NB2,LFCHG,NCU,ESC,IDI MBK,INTS,NBT,JNT,Y,EL,EH,LFW)
RETURN
END
SUBROUTINE ANNI01(NE,SIGT,SIGG,SIGS,SIGF,NIS,MIS,E,NB,BK,ISTGF,
*NB2,BK2,LFCHG,NCU,ESC,IDI MBK,INTS,NBT,JNT,Y,EL,EH,LFW)
DIMENSION SIGT(1),SIGG(1),SIGS(1),SIGF(1),E(1),NAM1(2),BK(1),
*BK2(1),ESC(1),INTS(1),NBT(1),JNT(1),Y(1)
REAL *8 NAMISO,NAMB
EQUIVALENCE (NAMB,NAM1(1))
COMMON/INOUT/ NO,NI,NT,MODE,LEDIT1,LEDIT2,LEDITS,NR1(75),
*MATKDK,NTYP,NAMISO(10),MATISO(10),NR2(11),NBX,N1X,N2X,N,NBKX
COMMON/RECS/ MAT,ME,MT,C1,C2,L1,L2,N1,N2
C
C CALCULATION OF TOTAL CROSS-SECTION
C
DO 44 K=1,NE
STGT(K)=SIGG(K)+SIGS(K)+SIGF(K)
44 CONTINUE
M+IS=1
MATK=MATKOK
IF(NIS.EQ.1)GOTO 105
MATK=MATISO(MIS)
M+IS=MIS+1

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2980
2990
3000
3010
302C
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307C
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3090
3100
3110
312C
3130
314C
3150
3160
3170
3180
319C
3200
3210
322C
3230
3240
3250
3260
327C
10
20
30
4C
50
6C
7C
80
9C
100
11C
120
13C
14C
150
16C
170
180
19C
200

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105	IF(LEDITS.NE.1)GOTO 101	210	CONTINUE	760
	WRITE(NO,43)	220	WRITE(NO,317)(ESC(NCU),NCU=1,NCU)	770
43	FORMAT(1H1,'POINT-WISE CROSS SECTIONS FROM UNRESOLVED PARAMETERS'	230	FORMAT('0ENERGY POINTS ADDED BY CHGINT IN THE STGF DATA TYPE'//	780
*	////1X,'ENERGY',13X,	240	* (1X,1P6E15.5))	790
*	1X,'SCATTERING',5X,'CAPTURE',8X,'FISSION',8X,'TCTAL'//)	250	C INSERTION OF THE NEW ENERGY POINTS IN BK ARRAY	800
	WRITE(NO,60)(E(K),SIGS(K),SIGG(K),SIGF(K),SIGT(K),K=1,NE)	260	NBL=N/NE	810
60	FORMAT(1X,1P5F15.5)	270	DO 318 INCU=1,NCU	820
101	CALL XTEXT(4H ,NAME2)	280	IF((N+NBL).GT.IDIMBK)CALL ERROR(226)	830
	N=NB-1	290	DO 319 IK=1,N,NBL	840
C		300	IF(ESC(INCU).EQ.BK(IK))GOTO 318	850
C	PRINTING OF KEDAK DATA TYPE ST	310	IF(ESC(INCU).LT.BK(IK))GOTO 320	860
C		320	319 CONTINUE	870
	IF(LEDIT2.NE.1)GOTO 102	330	CALL ERROR(225)	880
	WRITE(NO,91)	340	DO 321 IK1=IK,N	890
81	FORMAT(1H1,'KEDAK DATA TYPE ST')	350	NCRT=N-1K1+1K	900
	IF(NIS.GT.1)WRITE(NO,110)NAMISO(MIS)	360	321 BK(NCRT+NBL)=BK(NCRT)	910
110	FORMAT(1H+,60X,'ISOTOPE',3X,A8)	370	N=N+NBL	920
	CALL XTEXT(4HST ,NAM1(1))	380	KF=NBL/11	930
	CALL XTEXT(4H ,NAM1(2))	390	DO 323 NJK=1,KF	940
	CALL ABFORM(NO,NAM8)	400	BK(IK)=ESC(INCU)	950
	CALL LUCY(3,BK)	410	DO 322 JK=4,10	960
102	CALL XTEXT(4HST ,NAME1)	420	IK2=IK+JK	970
	CALL WDA2(MATK,MMIS , NAME1,NAME2)	430	CALL TERP1(BK(IK-NBL),BK(IK2-NBL),BK(IK+NBL),BK(IK2+NBL),	980
C		440	* BK(IK),BK(IK2),INTS(NJK))	990
C	REORDERING OF DATA IN STGF: E,L,J	450	322 CONTINUE	1000
C		460	IK=IK+11	1010
	IF(ISTGF.EQ.0)GOTO 83	470	323 CONTINUE	1020
	N=NB2-1	480	318 CONTINUE	1030
	K=1	490	GOTO 312	1040
	DO 251 IE=1,NE	500	310 WRITE(NO,311)	1050
	K2=(IE-1)*11+1	510	311 FORMAT('0NO ENERGY POINTS ADDED BY CHGINT IN STGF')	1060
252	IF((K+11).GT.NBKX) CALL ERROR(204)	520		1070
	DO 250 IK=1,11	530	C PRINTING OF KEDAK DATA TYPE STGF	1080
	I=IK-1	540	C	1090
250	BK(K+I)=BK2(K2+I)	550	312 IF(LEDIT2.NE.1)GOTO 103	1100
	K=K+11	560	WRITE(NO,94)	1110
	K2=K2+11*NF	570	84 FORMAT(1H1,'KEDAK DATA TYPE STGF')	1120
	IF(K2-N)252,252,251	580	IF(NIS.GT.1)WRITE(NO,110)NAMISO(MIS)	1130
251	CONTINUE	590	CALL XTEXT(4HSTGF,NAM1(1))	1140
C		600	CALL ABFORM(NO,NAM8)	1150
C	INSERTION IN STGF OF ENERGY POINTS ADDED BY CHGINT	610	CALL LUCY(2,BK)	1160
C		620	103 CALL XTEXT(4HSTGF,NAME1)	1170
	IF(LFCHG.EQ.0.OR.ISTGF.EQ.0)GOTO 312	630	CALL WDA2(MATK,MMIS , NAME1,NAME2)	1180
	IF(NCU.EQ.NE)GOTO 310	640	83 N2=NE	1190
C	SELECTION IN THE ARRAY ESC OF POINTS ADDED BY CHGINT	650	N1=1	1200
	DO 313 IE=1,NE	660	NBT(1)=N2	1210
	DO 314 INCU=1,NCU	670	JNT(1)=INTS(1)	1220
	IF(E(IE).GE.ESC(INCU)*0.99999.AND.E(IE).LE.ESC(INCU)*1.00001)	680	BK2 AND X HAVE THE SAME STARTING ADDRESS.	1230
*	GOTO 315	690	CALL MAGGY(E,SIGS,NE,BK2,Y)	1240
314	CONTINUE	700	CALL WRECS2(MIS,4HSCAT,4HSTAT,EL,EH)	1250
	GOTO 313	710	CALL MAGGY(E,SIGG,NE,BK2,Y)	1260
315	NCU=NCU-1	720	CALL WRECS2(MIS,4HCAPT,4HSTAT,EL,EH)	1270
	IF(NCU.EQ.0)CALL ERROR(224)	730	IF(LFW.NE.1)GOTO 70	1280
	DO 316 IESC=1,NCU	740	CALL MAGGY(E,SIGF,NE,BK2,Y)	1290
316	ESC(IESC)=ESC(IESC+1)	750	CALL WRECS2(MIS,4HFISS,4HSTAT,EL,EH)	1300

70 CONTINUE	1310	SUBROUTINE GNRL (GALPHA, GBETA, GAMMA, MU, NU, LAMBDA, S, DF, ID)	10
RETURN	1320	COMMON/XBAR/XX(4,10)	20
END	1330	S=0.0	30
		IF (GALPHA) 1000, 1000, 1001	40
		1001 IF (GAMMA) 1000, 1000, 1002	50
		1002 IF (GBETA) 1000, 1003, 1004	60
		1003 IF (DF) 1000, 1005, 1006	70
		1005 DO 100 J=1, 10	80
	10	XJ=XX(MU, J)	90
	20	GO TO (200, 201, 202), ID	100
	30	200 S=S+((XJ*XJ)/(GALPHA*XJ+GAMMA))	110
	40	GO TO 100	120
	50	201 S=S+(XJ/(GALPHA*XJ+GAMMA))	130
	60	202 CONTINUE	140
	70	100 CONTINUE	150
	80	S=S/10.0	160
	90	GO TO 1000	170
	100	1006 DO 101 J=1, 10	180
	110	XJ=XX(MU, J)	190
	120	DO 102 K=1, 10	200
	130	XK=XX(LAMBDA, K)	210
	140	GO TO (300, 301, 302), ID	220
	150	300 S=S+((XJ*XJ)/(GALPHA*XJ+GAMMA+DF*XK))	230
	160	GO TO 102	240
	170	301 S=S+(XJ/(GALPHA*XJ+GAMMA+DF*XK))	250
		302 CONTINUE	260
		102 CONTINUE	270
		101 CONTINUE	280
		S=S/100.0	290
		GO TO 1000	300
	10	1004 IF (DF) 1000, 1007, 1008	310
	20	1007 DO 103 J=1, 10	320
	30	XJ=XX(MU, J)	330
	40	DO 104 K=1, 10	340
	50	XK=XX(NU, K)	350
	60	GO TO (400, 401, 402), ID	360
	70	400 S=S+((XJ*XJ)/(GALPHA*XJ+GBETA*XK+GAMMA))	370
	80	GO TO 104	380
		401 S=S+(XJ/(GALPHA*XJ+GBETA*XK+GAMMA))	390
		GO TO 104	400
		402 S=S+((XJ*XK)/(GALPHA*XJ+GBETA*XK+GAMMA))	410
		104 CONTINUE	420
		103 CONTINUE	430
		S=S/100.0	440
		GO TO 1000	450
	10	1008 DO 105 J=1, 10	460
	20	XJ=XX(MU, J)	470
	30	DO 106 K=1, 10	480
	40	XK=XX(NU, K)	490
	50	DO 107 L=1, 10	500
	60	XL=XX(LAMBDA, L)	510
	70	GO TO (500, 501, 502), ID	520
		500 S=S+((XJ*XJ)/(GALPHA*XJ+GBETA*XK+GAMMA+DF*XL))	530
		GO TO 107	540
		501 S=S+(XJ/(GALPHA*XJ+GBETA*XK+GAMMA+DF*XL))	550
			550

SUBROUTINE ERF(EER)	10
COMMON/FLCT/ZYN, ZYF, U, V, RHON, RHOF	20
SW=0.05	30
SUMME=0.	40
HU=0.	50
HC=SW	60
DC 1 J=1, 9999	70
CALL GAUSS(HU, HO, B)	80
HL=HO	90
HO=HO+SW	100
IF (SQRT(B*B)-1.E-6) 6, 6, 3	110
3 SUMME=SUMME+B	120
1 CONTINUE	130
6 SUMME=SUMME*(1.+2./ZYN)**RHON	140
EER=SUMME	150
RETURN	160
END	170

SUBROUTINE GAUSS (HU, HO, B)	10
D=(HO-HU)*0.5	20
F=(HO+HU)*0.5	30
B=(BINT(9.061799E-1*D+E)*2.369269E-1+BINT(5.384693E-1*D+E)	40
1*4.786287E-1+BINT(E)*5.688889E-1+BINT(-5.384693E-1*D+E)	50
2*4.786287E-1+BINT(-9.061799E-1*D+E)*2.369269E-1)*D	60
RETURN	70
END	80

FUNCTION BINT(X)	10
COMMON/FLCT/ZYN, ZYF, U, V, RHON, RHOF	20
Z=ZYN*0.5+1.+RHON	30
Y=ZYF*0.5+RHOF	40
BINT=EXP(-X)/((1.+2.*U*X/ZYN)**Z*(1.+2.*V*X/ZYF)**Y)	50
RETURN	60
END	70



GO TO 107	560	CALL STORE(3,1000,LOF)	100
502 S=S+((XJ*XK)/(GALPHA*XJ+GBETA*XK+GAMMA+DF*XL))	57C	IF(LOF.EQ.1)CALL ERROR(408)	200
107 CONTINUE	580	IF(LFW.NE.1)GOTO1	210
106 CONTINUE	59C	CALL RRECS(0,4HSGF,NAME2,EL,EH,LNTR)	220
105 CONTINUE	600	IF(LNTR.NE.0)CALL ERROR(777)	230
S=S/1000.0	610	CALL STORE(3,1001,LOF)	240
1000 RETURN	62C	IF(LOF.EQ.1)CALL ERROR(408)	250
END	630	6 CALL COMB(ADD,CON,1000,1001,EL,EH,EPS,LCF,LNT)	260
		IF(LOF.EQ.0)GOTO 5	270
		EPS=EPS+EPSINC	280
		IF(EPS-EPSMAX)6,6,7	290
		7 CALL ERROR(409)	300
		5 IF(LNT.EQ.1)CALL ERROR(410)	310
FUNCTION ADD(YA,YB,CON)	10	CALL STORE(3,1000,LOF)	320
DIMENSION CON(20)	20	IF(LOF.EQ.1)CALL ERROR(408)	330
ADD=YA+YB	30	1 CONTINUE	340
RETURN	40	CALL FETCH(1000,LNT)	350
END	50	IF(LNT.EQ.1)CALL ERROR(410)	360
		WRITE(ND,10)EPS	370
		10 FORMAT(140,'EPS=',1PE12.4,' FOR CALCULATING SGT'//)	380
		RETURN	390
		END	400
FUNCTION SUB(YA,YB,CON)	10		
SUB=YA-YB	20		
RETURN	30		
END	40		
REAL FUNCTION MULT(YA,YB,CON)	10	SUBROUTINE MYRIAM(NIS,LISRES,LISUNR,NAME,EL1,EH1,EPS,ITEST,Y)	10
MULT=YA*YB	20	EXTERNAL ADD	20
RETURN	30	DIMENSION CON(20),Y(1)	30
END	40	EL1=1.0E8	40
		FH1=1.0E-6	50
		IFIRST=)	60
		DO 10 MIS=1,NIS	70
		1 IF(ITEST.EQ.1)GOTO 5	80
		IF(LISRES.NE.0)RETURN	90
		CALL RRECS(MIS,NAME,4HRES,EL,EH,LNTR)	100
		IF(LNTR.NE.0)CALL ERROR(777)	110
		CALL BETSY(MIS,Y)	120
		EL1=AMIN1(EL,EL1)	130
		FH1=AMAX1(FH,FH1)	140
		IF(IFIRST)4,4,3	150
		3 CALL STORE(3,1001,LOF)	160
		IF(LOF.EQ.1)CALL ERROR(407)	170
		CALL COMB(ADD,CON,1000,1001,EL1,EH1,EPS,LOF,LNT)	180
		IF(LOF.EQ.1)CALL ERROR(403)	190
		IF(LNT.EQ.1)CALL ERROR(404)	200
		CALL STORE(3,1000,LOF)	210
		IF(LOF.EQ.1)CALL ERROR(407)	220
		GOTO 9	230
		4 CALL STORE(3,1000,LCF)	240
		IF(LOF.EQ.1)CALL ERROR(407)	250
		IFIRST=1	260
		GOTO 9	270
		5 IF(LISUNR.NE.0)RETURN	280
		CALL RRECS(MIS,NAME,4HSTAT,EL,EH,LNTR)	290
SUBROUTINE MARION(LFW,EL,EH,EPS,NAME2)	10		
COMMON/INOUT/NO	20		
EXTERNAL ADD	30		
DATA EPSMAX/2.E-2/,EPSINC/1.E-3/	40		
CALL RRECS(0,4HSGN,NAME2,EL,EH,LNTR)	50		
IF(LNTR.NE.0)CALL ERROR(777)	60		
CALL STORE(3,1000,LCF)	70		
IF(LOF.EQ.1)CALL ERROR(408)	80		
CALL RRECS(0,4HSGG,NAME2,EL,EH,LNTR)	90		
IF(LNTR.NE.0)CALL ERROR(777)	100		
CALL STORE(3,1001,LOF)	110		
IF(LOF.EQ.1)CALL ERROR(408)	120		
3 CALL COMB(ADD,CON,1000,1001,EL,EH,EPS,LCF,LNT)	130		
IF(LOF.EQ.0)GOTO 2	140		
EPS=EPS+EPSINC	150		
IF(EPS-EPSMAX)3,3,4	160		
4 CALL ERROR(409)	170		
2 IF(LNT.EQ.1)CALL ERROR(410)	180		

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IF(LNTR.NE.0) CALL ERROR(777)
CALL BETSY(MIS,Y)
EL1=AMINI(EL,FL1)
EH1=AMAX1(EH,EH1)
IF(IFIRST)8,8,7
7 CALL STORE(3,1001,LOF)
IF(LOF.EQ.1)CALL ERROR(407)
CALL COMB(ADD,CON,1000,1001,FL1,EH1,EPS,LOF,LNT)
IF(LOF.EQ.1)CALL ERROR(403)
IF(LNT.EQ.1)CALL ERROR(404)
CALL STORE(3,1000,LOF)
IF(LOF.EQ.1)CALL ERROR(407)
GOTO 9
8 CALL STORE(3,1000,LOF)
IF(LOF.EQ.1)CALL ERROR(407)
IFIRST=1
9 CONTINUE
10 CONTINUE
CALL FETCH(1000,LNT)
IF(LNT.EQ.1)CALL ERROR(404)
RETURN
END

```

```

300
31C
320
33C
34C
350
36C
37C
380
390
400
41C
420
43C
440
45C
46C
470
480
490
500
51C

```

```

SUBROUTINE SOPHIE(MATP,LISRES,LISUNR,LFW,LNU,EPS,BK,XIN,YIN,NCMAX,
*NBT,JNT,X,Y,B,TEMP)
C-----CHANGED OCTOBRE 1973 BY E.STEIN.
LOGICAL*4 SALINE,SECONC,LODZMS
COMMON/RECS/MAT,MF,MT,C1,C2,L1,L2,N1,N2
COMMON/INDUT/NO,NI,NT,MODE,LEDIT1,LEDIT2,LEDAVE,NR1(20),
*NOTR1(50),BIB12,SALINE,NR2(2),MATKCK,NR3(42),NEX,NLX,N2X
COMMON/TAB/ENDTAB(25),KEDTAB(50),KEDR(240),NCNAME(39),NRIF3,NRIF4
EQUIVALENCE (KEDR(1),KCR1(1)),(KEDR(124),KCR2(1))
INTEGER ENDTAB,BLANK,SGIZ,CZ,BIB12
EXTERNAL ADD,SUB,MULT
DIMENSION IABSP(8),CON(1),TEMP(1),NAM1(2),XIN(1),YIN(1),BK(1),
*KCR1(123),KCR2(1),NBT(1),JNT(1),X(1),Y(1),B(1)
DATA BLANK/' ',SGIZ/'SGIZ'/,CZ/'C '/
REAL*8 NAM8
EQUIVALENCE (NAM1(1),NAM8,NAMEK1),(NAMEK2,NAM1(2))
WRITE(NO,6)
6 FORMAT(1H1,20X,'FILE 3 - SMOOTH CROSS SECTIONS'//21X,32('*')//)
WRITE(NO,21)
21 FORMAT(1X,'PRINTING AFTER COMBINATION WITH FILE 2 DATA, BEFORE CHG
*INT'//)
CALL PRIZEI
NF=10
CALL INITSC(NCMAX)
REWIND NF
IFSGIP=0
IFSGIC=0
IFSGA=0
IFSGI=0

```

```

10
20
30
4C
5C
6C
70
8C
9C
100
11C
120
13C
14C
150
16C
170
18C
19C
200
21C
220
230
24C
250
26C
270
280
29C

```

```

DO 5 I=1,25
MTP=FNDTAB(I)
NAMEK1=KEDTAB(2*I-1)
NAMEK2=KEDTAB(2*I)
DO 80 INTRL=1,50,2
IF(NOTR1(INTRL).NE.NAMEK1)GOTO 8C
IF(NOTR1(INTRL+1).NE.NAMEK2)GOTO 8D
WRITE(NO,85)NAMEK1,NAMEK2
85 FORMAT('0 DATA TYPE ',2A4,' NOT TRANSLATED. (USER'S REQUEST)')
GOTO 51
80 CONTINUE
1 CALL SEARCH(NT,MCDF,MATP,3,MTP,LNT)
IF(LNT.NE.1)GOTO 3
WRITE(NO,2) MTP
2 FORMAT(1H0,'DATA TYPE NUMBER',15,' NOT INCLUDED IN THE ENDF/B FILE
*')
BACKSPACE NT
MT=MTP
GOTO 51
3 CALL RREC(3,NT,MODE,-1.0,NBT,JNT,X,Y)
IF(MT.EQ.27)IFSGA=1
EL=X(1)
EH=X(N2)
IF(MT.NE.18)GOTO 7
IF(LNU.NE.1)GOTO 7
DO 18 J=1,6
IF(Y(J).GT.1.E-6)GOTO 8
18 CONTINUE
8 IF(J-1)9,9,10
9 ENU=X(1)
GOTO 11
10 ENU=X(J-1)
11 ENUMAX=X(N2)
7 IF(LISRES.EQ.1.AND.LISUNR.EQ.1)GOTO 4
IF(MT.GT.2.AND.MT.NE.18.AND.MT.NE.102)GOTO 4
CALL STORE(3,1000,LOF)
IF(LOF.EQ.1)CALL ERROR(500)
IF(MT.EQ.1)CALL XTEXT(4HSGT ,NAME1)
IF(MT.EQ.2)CALL XTEXT(4HSGN ,NAME1)
IF(MT.EQ.18)CALL XTEXT(4HSGF ,NAME1)
IF(MT.EQ.102)CALL XTEXT(4HSGG ,NAME1)
IF(LFW.NE.1.AND.MT.EQ.18)GOTO 4
DO 20 NA=1,2
IF(NA.EQ.1)CALL XTEXT(4HPRES ,NAME2)
IF(NA.EQ.2)CALL XTEXT(4HSTAT,NAME2)
CALL RRECS(3,NAME1,NAME2,EL1,EH1,LNTR,NBT,JNT,X,Y)
IF(LNTR.EQ.1)GOTO 20
CALL STORE(3,1001,LOF)
IF(LOF.EQ.1)CALL ERROR(500)
EL=AMINI(EL,FL1)
EH=AMAX1(EH,EH1)
CALL COMB(ADD,CON,1000,1001,EL,EH,EPS,LOF,LNT,NET,JNT,X,Y)
IF(LOF.EQ.1)CALL ERROR(501)
IF(LNT.EQ.1)CALL ERROR(502)
CALL STORE(3,1000,LOF)

```

```

300
31C
320
330
34C
350
36C
370
380
390
400
41C
420
430
440
450
46C
470
480
490
500
51C
520
530
540
550
56C
570
580
590
600
61C
620
63C
640
650
66C
670
68C
690
700
710
720
73C
740
750
760
770
78C
790
80C
810
82C
83C
840

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      IF(LOF.EQ.1)CALL ERROR(500)
20 CONTINUE
   4 IF(KEDAVE.EQ.0.AND.MT.EQ.4)GOTO 86
      CALL WRECS2(0,MF,MT,0.0,0.0)
86 IF(LEDITS.NE.1)GOTO 50
      WRITE(NO,22)MT,NAMEK1,NAMEK2
22 FORMAT(1H1,'ENDF/B REACTION TYPE',I4,20X,'KEDAK REACTION TYPE',
* 3X,2A4/)
      CALL PRIZEI
      CALL WREC(3,NO,4)
C-----INELASTIC CROSS-SECTION
   50 IF(KEDAVE.NE.0.OR.MT.NE.4)GOTO 51
C-----ENERGY GRID FROM REACTION 4
      IFSGI=1
      CALL WREC(3,NF,2)
      CALL CHGINT(EPS,N1,N2,NBT,JNT,X,Y,N1X,N2X,&70)
      IF(N2.GT.NCMAX)GOTO 76
      CALL GENER(N2,X,TEMP,XIN,ND,NCU)
      IF(NCU.GT.NCMAX)CALL ERROR(506)
      NKI=NCU
      GOTO 5
C-----INELASTIC EXCITATION CROSS-SECTIONS   MT=51 TO 53
   51 IF(MT.NE.29)GOTO 5
      IN=50
      DO 81 INTRL=1,50,2
      IF(NOTTRL(INTRL).NE.SGIZ)GOTO 81
      IF(NOTTRL(INTRL+1).NE.BLANK)GOTO 81
      WRITE(NO,85)SGIZ,BLANK
      GOTO 65
   81 CONTINUE
      SECOND=.FALSE.
      INANF=51
      NAMEK1=SGIZ
      NAMEK2=BLANK
      DO 53 IN=51,90
      CALL SEARCH(NT,MODE,MATP,3,IN,LNT)
      IF(LNT.NE.1)GOTO 55
      WRITE(NO,2)IN
      BACKSPACE NT
      IF(.NOT.SALINE.OR.SECOND) GOTO 65
      GOTO 53
   55 IF(SECOND) GOTO 551
      SECOND=.TRUE.
      INANF=IN
551 CALL RREC(3,NT,MODE,-1.0,NBT,JNT,X,Y)
      IF(KEDAVE.EQ.0) GOTO 552
      EXC=ABS(C2)
      NRIF3=NRIF3+1
      CALL WRECS2(0,MF,IN,EXC,0.0)
      GOTO 553
552 IF(N2.GT.NCMAX)CALL ERROR(510)
      IFSGIP=1
553 IF(LEDITS.NE.1) GOTO 400
      WRITE(NO,22) IN,NAMEK1,NAMEK2
      CALL PRIZEI

```

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85C
860
870
880
890
900
910
920
930
940
950
960
970
980
990
1000
1010
1020
1030
1040
1050
1060
1070
1080
1090
1100
1110
1120
1130
1140
1150
1160
1170
1180
1190
1200
1210
1220
1230
1240
1250
1260
1270
1280
1290
1300
1310
1320
1330
1340
1350
1360
1370
1380
1390

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```

      CALL WREC(3,NO,4)
400 IF(KEDAVE.NE.0) GOTO 53
      CALL WREC(3,NF,2)
      CALL CHGINT(EPS,N1,N2,NBT,JNT,X,Y,N1X,N2X,&70)
      IF(N2.GT.NCMAX)GOTO 76
      CALL GENER(N2,X,TEMP,XIN,ND,NCU)
      IF(NCU.GT.NCMAX)CALL ERROR(506)
      NKI=NCU
   53 CONTINUE
      IN=91
C-----INELASTIC CROSS-SECTION TO THE CONTINUUM   MT=91
   65 INFN=IN-1
      DC 82 INTRL=1,50,2
      IF(NOTTRL(INTRL).NE.SGIZ)GOTO 82
      IF(NOTTRL(INTRL+1).NE.CZ)GOTO 82
      WRITE(NO,85)SGIZ,CZ
      GOTO 5
   82 CONTINUE
      IN=91
      CALL SEARCH(NT,MODE,MATP,3,91,LNT)
      IF(LNT.NE.1)GOTO 66
      WRITE(NO,2)IN
      BACKSPACE NT
      GOTO 5
   66 CALL RREC(3,NT,MODE,-1.0,NBT,JNT,X,Y)
      IFSGIC=1
      IF(LEDITS.NE.1) GOTO 401
      NAMEK2=CZ
      WRITE(NO,22) IN,NAMEK1,NAMEK2
      CALL PRIZEI
      CALL WREC(3,NO,4)
401 IF(KEDAVE.EQ.0) GOTO 554
      CALL WRECS2(0,MF,IN,0.0,0.0)
      GOTO 5
554 CALL WREC(3,NF,2)
      CALL CHGINT(EPS,N1,N2,NBT,JNT,X,Y,N1X,N2X,&70)
      IF(N2.GT.NCMAX)GOTO 76
      CALL GENER(N2,X,TEMP,XIN,ND,NCU)
      IF(NCU.GT.NCMAX)CALL ERROR(506)
      NKI=NCU
   5 CONTINUE
C
C
      IF(KEDAVE.NE.0) GOTO 74
      REWIND NF
      IF(IFSGI.EQ.0)GOTO 71
C
C-----REREAD TOTAL INELASTIC CROSS-SECTION FROM TEMP. FILE NF
C
      CALL RREC(3,NF,2,-1.0,NBT,JNT,X,Y)
   61 DO 62 KI=1,NKI
   62 CALL ITABI(XIN(KI),YIN(KI))
      DO 63 KI=1,NKI
      X(KI)=XIN(KI)
      Y(KI)=YIN(KI)

```

```

1400
1410
1420
1430
1440
1450
1460
1470
1480
1490
1500
1510
1520
1530
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1560
1570
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1590
1600
1610
1620
1630
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1700
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1940

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63	CONTINUE	1950	MT=MTP	2500
	N2=NKI	1960	IF(LNT.NE.0)GOTO 31	2510
	N1=1	1970	WRITE(NO,33) KEDTAB(2*IABSP(I)-1),KEDTAB(2*IABSP(I))	2520
	NBT(1)=NKI	1980	33 FORMAT(1X,2A4)	2530
	JNT(1)=2	1990	32 EL=AMINI(EL,X(I))	2540
	IF(Y(1).LE.1.0E-20) Y(1)=0.0	2000	EF=AMAX1(EH,X(N2))	2550
	IF(MT.NE.4)GOTO 72	2010	IF(SECOND)GOTO 45	2560
	CALL WRECS2(0,MF,MT,0.0,0.0)	2020	CALL STORE(3,1000,LOF)	2570
	GOTO 71	2030	IF(LOF.EQ.1)CALL ERROR(500)	2580
72	IF(IN.EQ.91)GOTO 68	2040	SECOND=.TRUE.	2590
	EXC=ABS(C2)	2050	GOTO 31	2600
	NRIF3=NRIF3+1	2060	45 CALL STORE(3,1001,LOF)	2610
	CALL WRECS2(0,MF,IN,EXC,0.0)	2070	IF(LOF.EQ.1)CALL ERROR(500)	2620
	GOTO 73	2080	CALL COMB(ADD,CON,1000,1001,EL,EF,EPS,LOF,LNT,NBT,JNT,X,Y)	2630
68	CALL WRECS2(0,MF,IN,0.0,0.0)	2090	IF(LOF.EQ.1)CALL ERROR(501)	2640
	GOTO 74	2100	IF(LNT.EQ.1)CALL ERROR(502)	2650
C		2110	CALL STORE(3,1000,LOF)	2660
C-----	REREAD PARTIAL INEL. CROSS-SECTIONS FROM TEMP. FILE NF	2120	IF(LOF.EQ.1)CALL ERROR(500)	2670
C		2130	31 CCNTINUE	2680
71	IF(IFSGIP.EQ.0)GOTO 75	2140	IF(SECOND) GOTO 23	2690
	DO 73 IN=INANF,INFN	2150	N1=1	2700
	CALL RREC(3,NF,2,-1.0,NBT,JNT,X,Y)	2160	N2=2	2710
	GOTO 61	2170	NBT(1)=2	2720
73	CONTINUE	2180	JNT(1)=2	2730
C		2190	X(1)=1.0E-5	2740
C-----	REREAD INEL. CROSS-SECTION TO THE CONTINUUM FROM TEMP. FILE NF	2200	Y(1)=0.0	2750
C		2210	X(2)=20.0E+6	2760
75	IF(IFSGIC.EQ.0)GOTO 74	2220	Y(2)=0.0	2770
	IN=91	2230	CALL STORE(3,1000,LOF)	2780
	CALL RREC(3,NF,2,-1.0,NBT,JNT,X,Y)	2240	IF(LOF.EQ.1) CALL ERROR(500)	2790
	GOTO 61	2250	WRITE(NO,1001)	2800
74	IF(IFSGA.EQ.1)GOTO 34	2260	1001 FORMAT('OTHE CROSS-SECTION TYPE SGA HAS BEEN SET TO 0.0 IN THE WHO	2810
C-----	ABSORPTION CROSS SECTION MT=27 SGA	2270	*LE ENERGY RANGE ( 1.0E-5 TO 20.0E+6 ) .'/	2820
	NAMEK1=KEDTAB(23)	2280	* THIS HAS BEEN DONE, BECAUSE NO CROSS-SECTIONS NECESSARY TO CALCU	2830
	NAMEK2=KEDTAB(24)	2290	*LATE SGA HAVE BEEN FOUND.')	2840
	DO 84 INTRL=1,50,2	2300	23 MT=27	2850
	IF(NOTTRL(INTRL).NE.NAMEK1)GOTO 84	2310	MF=3	2860
	IF(NOTTRL(INTRL+1).NE.NAMEK2)GOTO 84	2320	IF(LFDITS.NE.1)GOTO 35	2870
	WRITE(NO,85)NAMEK1,NAMEK2	2330	WRITE(NO,22)MT,NAMEK1,NAMEK2	2880
	GOTO 34	2340	CALL PRIZEI	2890
84	CONTINUE	2350	CALL WREC(3,NO,4)	2900
	IABSP(1)=15	2360	35 CALL WRECS2(0,MF,MT,0.0,0.0)	2910
	IABSP(2)=7	2370	C-----	2920
	DO 30 I=3,8	2380	TRANSPORT CROSS SECTION MT=201 SGTR	2930
30	IABSP(I)=I+13	2390	34 NAMEK1=KEDTAB(43)	2940
	EL=1.E8	2400	NAMEK2=KEDTAB(44)	2950
	EH=1.E-10	2410	DC 83 INTRL=1,50,2	2960
	WRITE(NO,1000)	2420	IF(NOTTRL(INTRL).NE.NAMEK1)GOTO 83	2970
1000	FORMAT('///OSGA WILL BE CALCULATED BY USING THE FOLLOWING CROSS-SE	2430	IF(NOTTRL(INTRL+1).NE.NAMEK2)GOTO 83	2980
	*CTIONS:')	2440	WRITE(NO,85)NAMEK1,NAMEK2	2990
	SECOND=.FALSE.	2450	GOTO 38	3000
	DC 31 I=1,8	2460	83 CCNTINUE	3010
	MTP=ENDTAB(IABSP(I))	2470	RN=0.0	3020
	RN=0.0	2480	CALL RRECS(0,3,2,RN,RN,LNT,NBT,JNT,X,Y)	3030
	CALL RRECS(0,3,MTP,RN,RN,LNT,NBT,JNT,X,Y)	2490	IF(LNT.EQ.0)GOTO 36	3040
			WRITE(NO,37)KEDTAB(3),KEDTAB(43)	

37	FORMAT(//IX,A4,2X,' NOT PRESENT. ',A4,' CANNOT BE CALCULATED'//)	3050	14	Y(K)=B(1)+B(2)*X(K)	3600
	GOTO 38	3060		GOTO 13	3610
36	EL=X(1)	3070	15	ET=X(K)	3620
	EH=X(N2)	3080		DO 16 I=2,NB	3630
	CALL STORE(3,1000,LOF)	3090		Y(K)=Y(K)+B(I)*ET	3640
	IF(LOF.EQ.1)CALL ERROR(500)	3100		ET=ET*X(K)	3650
	RN=0.0	3110	16	CONTINUE	3660
	CALL RRECS(0,3,251,RN,RN,LNT,NBT,JNT,X,Y)	3120	13	CONTINUE	3670
	IF(LNT.EQ.0)GOTO 39	3130		N1=1	3680
	WRITE(NO,37)KEDTAB(49),KEDTAB(43)	3140		NBT(1)=2.00	3690
	GOTO 38	3150		JNT(1)=2	3700
39	CALL STORE(3,1001,LCF)	3160		N2=200	3710
	IF(LOF.EQ.1)CALL ERROR(500)	3170		CALL WRECS2(0,1,452,0.0,0.0)	3720
	EL=AMINI(EL,X(1))	3180		RETURN	3730
	EH=AMAXI(EH,X(N2))	3190	70	CALL ERROR(505)	3740
	CALL COMB(MULT,CON,1000,1001,EL,EH,EPS,LOF,LNT,NBT,JNT,X,Y)	3200		STOP	3750
	IF(LOF.EQ.1)CALL ERROR(501)	3210	76	CALL ERROR(506)	3760
	IF(LNT.EQ.1)CALL ERROR(502)	3220		STCP	3770
	CALL STORE(3,1001,LCF)	3230		END	3780
	IF(LOF.EQ.1)CALL ERROR(500)	3240			
	RN=0.0	3250			
	CALL RRECS(0,3,1,RN,RN,LNT,NBT,JNT,X,Y)	3260			
	IF(LNT.EQ.0)GOTO 40	3270			
	WRITE(NO,37)KEDTAB(1),KEDTAB(43)	3280			
	GOTO 38	3290			
40	CALL STORE(3,1000,LOF)	3300		SUBROUTINE ELIMZ(N1,N2,NBT,JNT,X,Y)	10
	IF(LOF.EQ.1)CALL ERROR(500)	3310		DIMENSION X(1),Y(1),NBT(1),JNT(1)	20
	EL=AMINI(EL,X(1))	3320		C*****	30
	EH=AMAXI(EH,X(N2))	3330		C* THIS SUBROUTINE SUPPRESSES THE POINTS WHERE Y=0.0,	40
	CALL COMB(SUB,CON,1000,1001,EL,EH,EPS,LOF,LNT,NBT,JNT,X,Y)	3340		C* EXCEPT THE POINT PRECEDING THE THRESHOLD OF THE CROSS-SECTION	50
	IF(LOF.EQ.1)CALL ERROR(501)	3350		C*****	60
	IF(LNT.EQ.1)CALL ERROR(502)	3360		IF(N1.GT.1.OR.JNT(1).GT.2.OR.NBT(1).NE.N2)CALL ERROR(777)	70
	MT=201	3370		C	80
	MF=3	3380		C=====ELIMINATE ZERO POINTS AT THE END OF ARRAYS X AND Y	90
	IF(LEDITS.NE.1)GOTO 41	3390		C	100
	WRITE(NO,22)MT,NAMEK1,NAMEK2	3400		DO 1 I1=1,N2	110
	CALL PRIZEI	3410		I=I1	120
	CALL WREC(3,NO,4)	3420		IF(Y(N2-I+1).NE.0.0)GOTO 2	130
41	CALL WRECS2(0,MF,MT,0.0,0.0)	3430		1 CONTINUE	140
C-----	ETA CROSS SECTION MT=206 ETA	3440		C-----ALL POINTS ARE ZERO. SAVE ONLY FIRST AND LAST CNES.	150
38	CONTINUE	3450		7 X(2)=X(N2)	160
C-----	NUE DATA TYPE MT=452	3460		N2=2	170
	IF(LNU.NE.1)RETURN	3470		GOTO 3	180
	CALL RRECS4(0,4H NUE,4H ,LNT,B)	3480		C-----CROP END OF ARRAYS	190
	IF(LNT.EQ.1)RETURN	3490		C-----THE FIRST POINT EQUAL ZERO AT THE END OF THE ARRAY AND THE LAST	200
	NB=N1	3500		C-----POINT EQUAL ZERO AT THE END OF THE ARRAY WILL BE SAVED.	210
	X(1)=ENU	3510		2 IF(I.LE.3)GOTO 4	220
	DE=ALOG(ENMAX/ENU)/199	3520		N2=N2-I+3	230
	DE=EXP(DE)	3530		X(N2)=X(N2-I-3)	240
	DO 12 K=2,200	3540		C	250
	X(K)=X(K-1)*DE	3550		C=====ELIMINATE ZERO POINTS AT THE BEGINNIG OF ARRAYS X AND Y.	260
12	CONTINUE	3560		C	270
	DC 13 K=1,200	3570		4 DO 5 I1=1,N2	280
	Y(K)=B(1)	3580		I=I1	290
	IF(NB-2)13,14,15	3590		IF(Y(I).NE.0.0)GOTO 6	300
				5 CONTINUE	310
				GOTO 7	320

```

C-----CROP LEADER ZEROS.
  6 IF(I.LE.2) GOTO 3
    I=I-2
    N2=N2-1
    DC 8 I2=1,N2
    X(I2)=X(I2+I)
  8 Y(I2)=Y(I2+I)
C
C=====FINISHED
C
  3 NBT(1)=N2
    RETURN
    END

SUBROUTINE SUZY(LNU,ITAPE,BK,A,TAB,ITAB,NBT,JNT,X,Y,B)
LOGICAL*4 SALINE,SECOND
COMMON/INOUT/NO,NI,NT,MODE,NR1(3),KEDAVE,NR2(7),SALINE,NR3(2),
*MATKDK,NR4(42),NBX,NIX,NZX,N,NBXX,NR5(10),EPS,SCREPS
COMMON/RECS/MAT,MF,MT,C1,C2,L1,L2,N1,N2,NS
COMMON/TAB/ENDTAB(25),KEDTAB(50),KEDR(240),NCNAME(39),NRIF3,NRIF4
EQUIVALENCE (KEDR(1),KDR1(1)),(KEDR(124),KDR2(1))
INTEGER ENDTAB
REAL*8 AM241,NAM8
EQUIVALENCE (NAM1(1),NAME1,NAM8)
DIMENSION A(1),TAB(1),ITAB(1),NAM1(2),BK(1),NBT(1),JNT(1),X(1),
*Y(1),KDR1(123),KDR2(1),B(1),SCRCLA(10),NUMSCR(10)
CALL IRRECS(NBT,JNT,X,Y,B)
NTP=15
NS=1
DATA EPSMAX/.02/,EPSINC/.001/
MF=3
IADIM=NBXX
C-----REACTION TYPES FROM FILE 3
DO 1 I=1,26
  PSI=EPS
  IF(ITAPE.EQ.1)CALL WCONT(NTP,MODE,MAT,MF,NS)
  IF(I.LE.25) MT=ENDTAB(I)
  IF(I.EQ.26) MT=91
21 RN=0.0
  CALL RRECS(O,MF,MT,RN,RN,LNT,NBT,JNT,X,Y)
  IF(LNT.EQ.1)GOTO 1
  IF(KEDAVE.EQ.0.AND.(MT.EQ.4.OR.MT.EQ.91)) GOTO 132
  CALL CHGINT(PSI,N1,N2,NBT,JNT,X,Y,NIX,NZX,820)
  CALL ELIMZ(N1,N2,NBT,JNT,X,Y)
132 CALL ADDPNT(X,Y,NBT,JNT,MT)
  IF(KEDAVE.EQ.0.AND.(MT.EQ.4.OR.MT.EQ.91)) GOTO 14
  XL=X(1)
  XH=X(N2)
  CALL CROP(XL,XH,SCREPS,.TRUE.,EPS,SCRCLA,NUMSCR,LOF,NBT,JNT,X,Y)
  IF(LOF.NE.0) CALL ERROR(777)
  GOTO 14
20 PSI=PSI+EPSINC

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IF(PSI-EPSMAX)21,21,22
22 CALL ERROR(850)
14 IF(I.EQ.26) GOTO 40
NAME1=KEDTAB(2*I-1)
NAME2=KEDTAB(2*I)
GOTO 41
40 CALL XTEXT(4HSGIZ,NAME1)
CALL XTEXT(4HC ,NAME2)
41 CONTINUE
WRITE(NO,2)MT,PSI,SCREPS,EPS
2 FORMAT(1H1,'REACTION TYPE',I4,' AFTER CHGINT',8X,
*'PRECISION FPS=',F5.4,5X,'EPSMIN=',F5.4,' EPSMAX=',F5.4)
IF(KEDAVE.EQ.0.AND.(MT.EQ.4.OR.MT.EQ.91)) GOTO 134
WRITE(NO,133) (SCRCLA(J),J=1,10),(NUMSCR(J),J=1,10)
133 FORMAT(' SCRATCHCLASS (IN %) 0.00',10(' < <=',2PF4.2)/
*' NUMBER OF POINTS ',10I9)
IF(KEDAVE.EQ.0.AND.(MT.EQ.4.OR.MT.EQ.91)) GOTO 134
CALL DZMSCH(X,Y,NBT,JNT,NDZ,EPS,SCRCLA,NUMSCR)
IF(NDZ.GT.0) WRITE(NO,30) NDZ,(SCRCLA(J),J=1,9),(NUMSCR(J),J=1,10)
30 FORMAT(' TOTALLY',I7,' POINTS HAVE HAD X-VALUES (COMPARED WITH OT-
*ER POINTS) CLOSER THAN (RELATIVELY) 5.0E-6. '
*' THESE POINTS WERE SCRATCHED AND THE NEIGHBOURING X- AND Y-VALUES
* HAVE BEEN REPLACED BY AVERAGE VALUES. '
*' THESE POINTS COULD HAVE BEEN PREDICTED WITH THE FOLLOWING PRECIS
*IONS: '
*' CLASSES (IN %) 0.0',9(' < <=',2PF4.1), ' < <=UNENDLICH' /
*' NUMBER OF POINTS',10I9)
134 IF(MT.EQ.1.OR.MT.EQ.2) CALL SUPNEG(MT,N1,N2,NBT,JNT,X,Y)
CALL PRI ZEI
NAM1(2)=NAME2
CALL ABFORM(NO,NAM8)
C1=0.0
C2=0.0
L1=0
L1=0
CALL WREC(3,NO,4)
IF(ITAPE.EQ.1)CALL WREC(3,NTP,3)
K=1
DC 3 J=1,N2
BK(K)=X(J)
BK(K+1)=Y(J)
K=K+2
3 CONTINUE
N=N2*2
CALL WDA2(MATKDK,1,NAME1,NAME2)
1 CONTINUE
C-----INELASTIC EXCITATION CROSS SECTIONS
NW=NRIF3*4+1
IA=NW
NINIT=1
NWS=NW
NRS=1
SECOND=.FALSE.
CALL XTEXT(4HSGIZ,NAME1)
CALL XTEXT(4H ,NAME2)

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DO 7 MT=51,90
PS1=EPS
131 RN=0.0
CALL RRECS(0,MF,MT,EXC,RN,LNT,NBT,JNT,X,Y)
IF(LNT.EQ.0)GOTO 12
IF(SECOND) GOTO 13
IF(SALINE) GOTO 7
GOTO 8
130 PS1=PS1+EPSINC
IF(PS1.LE.EPSMAX) GOTO 131
CALL ERROR(850)
12 IF(SECOND) GOTO 121
SECOND=.TRUE.
121 IF(KEDAVE.EQ.0) GOTO 135
CALL CHGINT(PS1,N1,N2,NBT,JNT,X,Y,N1X,N2X,&130)
CALL ELIMZ(N1,N2,NBT,JNT,X,Y)
135 CALL ADDPNT(X,Y,NBT,JNT,MT)
IF(KEDAVE.EQ.0) GOTO 123
XL=X(1)
XF=X(N2)
CALL CROP(XL,XH,SCREPS,.TRUE.,EPS,SCRCLA,NUMSCR,LOF,NBT,JNT,X,Y)
IF(LOF.NE.0) CALL ERROR(777)
123 WRITE(NO,2) MT,PS1,SCREPS,EPS
IF(KEDAVE.EQ.0) GOTO 136
WRITE(NO,133) (SCRCLA(J),J=1,10),(NUMSCR(J),J=1,10)
CALL DZMSCH(X,Y,NBT,JNT,NDZ,EPS,SCRCLA,NUMSCR)
IF(NDZ.GT.0) WRITE(NO,30) NDZ,(SCRCLA(J),J=1,9),(NUMSCR(J),J=1,10)
136 CALL PRIZEI
CALL XTEXT(4H 00 ,NAME1(2))
IH1=MT/10
NAME1(2)=NAME1(2)+IH1*65536+(MT-IH1*10)*256
CALL ABFORM(NO,NAM8)
C1=0.0
C2=-EXC
L1=0
L2=0
CALL WREC(3,NO,4)
DO 9 J=1,N2
A(IA)=X(J)
A(IA+1)=Y(J)
IA=IA+2
IF(IA.GE.IADIM)CALL ERRCR(515)
9 CONTINUE
TAB(NINIT)=EXC
ITAB(NINIT+1)=N2
ITAB(NINIT+2)=NRS
ITAB(NINIT+3)=NWS
NINIT=NINIT+4
NRS=(IA-1)/880+1
NWS=MOD(IA-1,880)+1
7 CCNTINUE
13 IA=IA-1
CALL XTEXT(4H ,NAME2)
N=IA
CALL WDA2(MATKDK,1,NAME1,NAME2)

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940
950
960
97C
980
990
1000
1010
102C
1030
104C
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107C
1080
109C
1100
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1120
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114C
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1170
1180
119C
1200
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132C
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137C
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1400
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147C
1480

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WRITE(NO,10)NAME1,NAME2
10 FORMAT(1H1,2A4,' REACTION TYPE TABLE'//)
NWM1=NW-1
WRITE(NO,11)(A(I),I=1,NWM1)
11 FORMAT(1X,1PE15.5,3I8)
C-----NUE DATA TYPE
8 PS1=EPS
28 RN=0.0
CALL RRECS(0,1,452,RN,RA,LNT,NBT,JNT,X,Y)
IF(LNT.EQ.0)GOTO 5
IF(ITAPE.EQ.1)CALL WCONT(NTP,MODE,MAT,0,NS)
GOTO 6
5 CALL CHGINT(PS1,N1,N2,NBT,JNT,X,Y,N1X,N2X,&23)
XL=X(1)
XF=X(N2)
CALL CROP(XL,XH,SCREPS,.TRUE.,EPS,SCRCLA,NUMSCR,LOF,NBT,JNT,X,Y)
IF(LOF.NE.0) CALL ERROR(777)
GOTO 24
23 PS1=PS1+EPSINC
IF(PS1-EPSMAX)28,28,25
25 CALL ERROR(851)
24 CALL XTEXT(4HNUE ,NAME1)
CALL XTEXT(4H ,NAME2)
MF=3
MT=452
WRITE(NO,2) MT,PS1,SCREPS,EPS
WRITE(NO,133) (SCRCLA(J),J=1,10),(NUMSCR(J),J=1,10)
CALL DZMSCH(X,Y,NBT,JNT,NDZ,EPS,SCRCLA,NUMSCR)
IF(NDZ.GT.0) WRITE(NO,30) NDZ,(SCRCLA(J),J=1,9),(NUMSCR(J),J=1,10)
CALL PRIZEI
NAME1(2)=NAME2
CALL ABFORM(NO,NAM8)
C1=0.0
C2=0.0
L1=0
L2=0
CALL WREC(3,NO,4)
K=1
DO 4 J=1,N2
BK(K)=X(J)
BK(K+1)=Y(J)
K=K+2
4 CCNTINUE
N=N2*2
CALL WDA2(MATKDK,1,NAME1,NAME2)
IF(ITAPE.EQ.1)CALL WREC(3,NTP,3)
6 IF(ITAPE.EQ.1)CALL WCONT(NTP,MODE,C,C,NS)
RETURN
END
149C
1500
1510
1520
1530
154C
1550
1560
1570
1580
159C
1600
161C
1620
1630
164C
1650
166C
1670
1680
1690
1700
171C
1720
1730
174C
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176C
1770
1780
179C
1800
181C
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184C
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1870
188C
189C
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191C
1920
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194C
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196C
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SUBROUTINE DZMSCH(X,Y,NBT,JNT,NDZ,EPS,SCRCLA,NUMSCR) 10
DIMENSION X(1),Y(1),NBT(1),JNT(1),SCRCLA(10),NUMSCR(10) 20
COMMON/RECS/ NR1(7),N1,N2 30
C N PUNKTE, DEREN X-WERTE AUF 6 STELLEN UEBEREINSTIMMEN, 40
C WERDEN FOLGENDERMASSEN ZUSAMMENGEFASST: 50
C XNEU=1/N*(X1+X2+...XN); YNEU=1/N*(Y1+Y2+...+YN) . 60
C 70
C IF(N1.GT.1.OR.JNT(1).GT.2.OR.NBT(1).NE.N2) CALL ERROR(777) 80
C BILDUNG DER SCRATCHCLASSEN. 90
SCRCLA(1)=EPS 100
HI1=(1.0-EPS)/9.0 110
DO 1 I=2,9 120
1 SCRCLA(I)=SCRCLA(1)+(I-1)*HI1 130
DO 2 I=1,10 140
2 NUMSCR(I)=0 150
NDZ=0 160
I=1 170
305 IF((I+1).GT.N2) GOTO 320 180
J=I+1 190
307 IF(J.GT.N2) GOTO 306 200
IF((X(J)-X(I))/X(J).GE.5.0E-6) GOTO 306 210
J=J+1 220
GOTO 307 230
306 IF(J.EQ.(I+1)) GOTO 308 240
C MIT WELCHER GENAUIGKEIT WAREN DIE PUNKTE I+1,...,J-1 ZWISCHEN 250
C DEN PUNKTEN I UND MIN(J,N2) ZU INTERPOLIEREN. (SIEHE CROP). 260
IP=I+1 270
JM1=J-1 280
MINJN2=MIN(J,N2) 290
DO 10 K=IP1,JM1 300
CALL TERP1(X(I),Y(I),X(MINJN2),Y(MINJN2),X(K),Y(2) 310
EPSAKT=ABS((Y(K)-YP)/Y(K)) 320
DO 11 KK=1,10 330
IF(KK.EQ.10) GOTO 12 340
IF(EPSAKT.LE.SCRCLA(10-KK)) GOTO 11 350
12 NUMSCR(11-KK)=NUMSCR(11-KK)+1 360
GOTO 10 370
11 CCNTINUE 380
10 CONTINUE 390
C MITTELWERT DER PUNKTE Y(I), Y(I+1), ..... Y(J-1) BILDEN 400
M=J-I 410
NDZ=NDZ+(M-1) 420
X1=0.0 430
Y1=0.0 440
DO 309 K=1,M 450
X1=X1+X(I+K-1) 460
309 Y1=Y1+Y(I+K-1) 470
X(I)=1.0/M*X1 480
Y(I)=1.0/M*Y1 490
C DELETEN DER PUNKTE X(I+1),Y(I+1), X(I+2),Y(I+2), ... X(J-1),Y(J-1) 500
C IF((J-1).NE.N2) GOTOC 310 510
C J=1=N2 520
N2=I 530
NBT(1)=N2 540
GOTO 320 550

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310 N=I+1 560
M=M-1 570
L=N2-M 580
DO 311 K=N,L 590
X(K)=X(K+M) 600
311 Y(K)=Y(K+M) 610
N2=N2-M 620
NBT(1)=N2 630
308 I=I+1 640
GOTO 305 650
320 CCNTINUE 660
RETURN 670
ENC 680

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SUBROUTINE ADDPNT(X,Y,NBT,JNT,MT) 10
C BEI REAKTIONEN MIT EINER SCHWELLE WERDEN DAS VORHANDENSEIN 20
C VON ZWEI NULLEN, EINE BEI 1.0E-5 UND EINE BEI DER SCHWELLE, 30
C NOTETIGENFALLS DURCH ERZEUGUNG SICHERGESTELLT. 40
LOGICAL*4 LT1 50
DIMENSION X(1),Y(1),NBT(1),JNT(1) 60
COMMON/RECS/ NR1(7),N1,N2 70
COMMON/INDUT/ NO,NR2(126),N2X 80
C NUR LINEARE ODER KONSTANTE INTERPOLATION UND EIN ENERGIEBEREICH 90
C ERLAUBT. 100
IF(N1.NE.1.OR.NBT(1).NE.N2.OR.JNT(1).GT.2) CALL ERROR(777) 110
C SCHWELLRREAKTION? 120
IF((ABS(X(1)-1.0E-5)/X(1).LE.1.0E-6).OR.(ABS(Y(1)).GT.1.0E-20)) 130
*RETURN 140
C Y(1).GT.0? 150
LT1=.FALSE. 160
IF(Y(1).EQ.0.0) GOTOC 2 170
LT1=.TRUE. 180
Y(1)=0.0 190
WRITE(NO,100) MT 200
100 FCRMAT('0$$$ FOR ENDF/B TYPE MF=3 MT=',I3) 210
WRITE(NO,101) 220
101 FCRMAT('+',31X,' Y(1) HAS BEEN SET TO ZERO.') 230
2 N2=N2+1 240
IF(N2.GT.N2X) CALL ERROR(105) 250
NBT(1)=N2 260
DO 1 I=2,N2 270
X(N2+2-I)=X(N2+1-I) 280
1 Y(N2+2-I)=Y(N2+1-I) 290
X(1)=1.0E-5 300
Y(1)=0.0 310
IF(.NOT.LT1) WRITE(NO,100) MT 320
IF(LT1) WRITE(NO,102) 330
102 FCRMAT(' $$$ ADDITIONALLY') 340
WRITE(NO,103) 350
103 FCRMAT('+',31X,' A POINT X=1.0E-5EV, Y=0.0BARN HAS BEEN ADDED.') 360
RETURN 370
ENC 380

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SUBROUTINE SUPNEG(MT,N1,N2,NBT,JNT,X,Y)
C FUER MT=1 UND MT=2 WERDEN NEGATIVE Y-BEREICHE
C DIAGNOSTIZIERT UND BESEITIGT.
INTEGER*4 ELINEG
DIMENSION X(1),Y(1),NBT(1),JNT(1),TEXT(2)
LOGICAL*4 NEG,PREVAL
COMMON/INOUT/ NO,NRI(152),ELINEG
DATA TEXT/'SGT ','SGN '/
IF(MT.NE.1.AND.MT.NE.2) CALL ERROR(777)
NEG=.FALSE.
PREVAL=.FALSE.
I=1
C LOOP UEBER ALLE ENERGIEPUNKTE.
4 IF(Y(I).GE.0.0) GOTO 2
IF(.NOT.PREVAL) INEGAN=I
IF(.NOT.NEG) WRITE(NO,100) TEXT(MT)
100 FORMAT('NEGATIVE VALUES FOUND FOR CROSS-SECTION-TYPE ',A4,'.')
IF(.NOT.PREVAL) WRITE(INC,101)
101 FORMAT('NUMBER ENERGY ( EV ) CROSS SECTION ( BARN )')
WRITE(NO,102) I,X(I),Y(I)
102 FORMAT(' ',16,2X,1PE15.7,4X,1PE15.7)
NEG=.TRUE.
PREVAL=.TRUE.
GOTO 1
C BESEITIGEN DER NEGATIVEN BEREICHE VON INEGAN BIS INEGEN
2 CONTINUE
IF(.NOT.PREVAL) GOTO 1
INEGEN=I-1
IF(INEGEN.LT.INEGAN) CALL ERROR(777)
C BESEITIGUNG ERFOLGT NUR AUF WUNSCH.
IF(ELINEG.NE.1) GOTO 1
IF(INEGAN.LE.1) CALL ERROR(777)
WRITE(NO,103) X(INEGAN),X(INEGEN),X(INEGAN-1),Y(INEGAN-1),
*X(INEGEN+1),Y(INEGEN+1)
103 FORMAT(' THE PREVIOUS VALUES FROM ',1PE15.7,' EV TO ',
*1PE15.7,' EV WILL BE SCRATCHED. '/
*' POINT BEFORE INTERVAL:',1PE15.7,' EV, ',1PE15.7,' BARN. '/
*' POINT AFTER INTERVAL:',1PE15.7,' EV, ',1PE15.7,' BARN. ')
C DIE PUNKTE INEGAN+1 BIS N2 WERDEN AUF INEGAN BIS
C N2-(INEGEN-INEGAN+1) GELEGT.
IANZ=N2-INEGEN
DO 3 II=1,IANZ
X(INEGAN+II-1)=X(INEGEN+II)
3 Y(INEGAN+II-1)=Y(INEGEN+II)
IF(N1.NE.1.OR.JNT(1).GT.2.OR.NBT(1).NE.N2) CALL ERROR(777)
N2=N2-(INEGEN-INEGAN+1)
NBT(1)=N2
PREVAL=.FALSE.
1 I=I+1
IF(I.LE.N2) GOTO 4
333 RETURN
END

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SUBROUTINE ISABEL(MATP,IF4,ITAB,TAB,CLEG,NBT,JNT,X,Y,B,
*A,ITAB2,TAB2)
COMMON/INOUT/NO,NI,NT,MODE,NRI(78),MATKDK,NR2(45),N,NBKKX,JX,IADIM
COMMON/RECS/MAT,MF,MT,C1,C2,L1,L2,N1,N2
COMMON/TAB/ENDTAB(25),KEDTAB(50),KEDR(240),NCNAME(39),NRIF3,NRIF4
EQUIVALENCE (KEDR(1),KDR1(1)),(KEDR(124),KCR2(1))
INTEGER ENDTAB
DIMENSION CLEG(1),ITAB(1),TAB(1),ITAB2(1),TAB2(1),
* NBT(1),JNT(1),X(1),Y(1),B(1),A(1),ISATZ(400),NAUX(2),
*KDR1(123),KDR2(1)
REAL*4 LEGP
REAL*8 NAMB
EQUIVALENCE (NAUX(1),NAMB)
IDLEG=NBKKX
NXMAX=NBKKX
NF=10
NF1=11
NINIT2=1
REWIND NF
WRITE(NO,1)
1 FORMAT(1H1,20X,'FILE 4 - ANGULAR DISTRIBUTIONS OF SECONDARY NEUT
*RCNS'//21X,54('*')//)
CALL PRIZEI
CALL SEARCH(NT,MODE,MATP,4,2,LNT)
IF(LNT.EQ.0)GOTO 4D
WRITE(NO,2)
2 FORMAT(1H0,'ELASTIC ANGULAR DISTRIBUTIONS NOT INCLUDED IN THE ENDF
*/B FILE')
BACKSPACE NT
GOTO 3
4C LVT=L1
LTT=L2
IF(LTT.EQ.2)GOTO 4
C
C=====LEGFCRE POLYNOMIAL COEFFICIENTS. LTT=1
C
IF(LVT.EQ.0)GOTO 5
C TRANSFORMATION MATRIX GIVEN LVT=1
CALL RREC(2,NT,MODE,-1.0,B)
GOTO 6
C TRANSFORMATION MATRIX NOT GIVEN LVT=0
5 CALL RREC(1,NT,MODE,-1.0)
6 LCT=L2
CALL PREC(4,NT,MODE,-1.0,C,NBT,JNT)
NE=N2
IF(LCT-1)9,9,10
C LCT=1 LAB SYSTEM
9 IKL=25
INCL=9
IKS=1
INCS=1

```

	GCTO 11	520	X(NPMAX/2+1)=0.0	1070
C	LCT=2 C.M. SYSTEM	530	DX=1.0/(NPMAX/2)	1080
10	IKL=28	540	DC 45 K=2,NPMAX	1090
	INCL=10	550	IF(K.EQ.(NPMAX/2+1))GOTO 45	1100
	IKS=4	560	X(K)=X(K-1)+DX	1110
	INCS=2	570	45 CCNTINUE	1120
11	NCNAME(INCL)=NE	580	X(NPMAX)=1.	1130
	NCNAME(INCS)=NE	590	NRS=(IA-1)/880+1	1140
	IKL2=IKL+2	600	NWS=MOD(IA-1,880)+1	1150
	NAUX(1)=KDR2( IKL )	610	DC 46 K=1,NPMAX	1160
	NAUX(2)=KDR2( IKL+1 )	620	Y(K)=LEGP(X(K),N1,B)/6.283184	1170
	CALL A8FORM(NC,NAM8)	630	A(IA)=X(K)	1180
	NAUX(1)=KDR2( IKS )	640	A(IA+1)=Y(K)	1190
	NAUX(2)=KDR2( IKS+1 )	650	IA=IA+2	1200
	CALL A8FORM(NO,NAM8)	660	IF(IA.GE.IADIM) CALL ERROR(799)	1210
	WRITE(NO,13)(KDR2(I),I=IKL,IKL2),NE	670	46 CCNTINUE	1220
13	FORMAT(1H0,'THE LEGENDRE POLYNOMIAL COEFFICIENTS FOR REACTION TYPE	680	N2=NPMAX	1230
	*,2X,2A4,2X,I6,2X,'ARE GIVEN AT',I6,' ENERGY PCINTS'/)	690	N1=1	1240
	CALL A8FORM(NO,'LEGENDRE')	700	NBT(1)=NPMAX	1250
	CALL A8FORM(NO,'COEFFIC.')	710	JNT(1)=2	1260
	NINIT=1	720	CALL WREC(3,NO,4)	1270
	NWORD=NE*5+1	730	GOTO 22	1280
	NWORD2=NE*4+1	740	7 CCNTINUE	1290
	NRIF4=NRIF4+2	750	GOTO 36	1300
	NLEGI=NWORD	760		1310
	IA=NWORD2	770	C	1320
	DO 7 I=1,NE	780	C====TABULATED DISTRIBUTION PROBABILITIES LTT=2	1330
	CALL RREC(2,NT,MODE,-1.0,B)	790	C	1340
	E=C2	800	C	1350
61	IF(B(N1).GT.0.)GOTO 60	810	4 IF(LVT.EQ.0)GOTO 16	1360
	IF(N1.EQ.1)GOTO 60	820	TRANSFORMATION MATRIX GIVEN LVT=1	1370
	N1=N1-1	830	CALL RREC(2,NT,MODE,-1.0,B)	1380
	GCTO 61	840	GOTO 17	1390
60	NWL=MOD(NLEGI-1,880)+1	850	C TRANSFORMATION MATRIX NCT GIVEN LVT=0	1400
	NRL=(NLEGI-1)/880+1	860	16 CALL RREC(1,NT,MODE,-1.0)	1410
	DC 8 N=1,N1	870	17 LCT=L2	1420
	IF(NLEGI.GE.IDLEG)CALL ERROR(700)	880	CALL RREC(4,NT,MODE,-1.0,NBT,JNT)	1430
	CLEG(NLEGI)=N	890	NE=N2	1440
	CLEG(NLEGI+1)=B(N)	900	IF(LCT-1)18,18,19	1450
8	NLEGI=NLEGI+2	910	18 IKS=1	1460
	TAB(NINIT)=E	920	INCS=1	1470
	TAB(NINIT+1)=N1	930	GOTO 20	1480
	ITAB(NINIT+2)=N1	940	19 IKS=4	1490
	ITAB(NINIT+3)=NRL	950	INCS=2	1500
	ITAB(NINIT+4)=NWL	960	20 NCNAME(INCS)=NE	1510
	NINIT=NINIT+5	970	IKS2=IKS+2	1520
	WRITE(NO,15)	980	NAUX(1)=KDR2( IKS )	1530
15	FCRMT(1H0//5X,' ENERGY', 7X,'LEGENDRE COEFFICIENTS'/)	990	NAUX(2)=KDR2( IKS+1 )	1540
	WRITE(NO,14)E,(B(N),N=1,N1)	1000	CALL A8FORM(NC,NAM8)	1550
14	FORMAT(1X,1P7E15.5/(16X,1P6E15.5))	1010	WRITE(NO,21)(KDR2(I),I=IKS,IKS2),NE	1560
	WRITE(NO,52)	1020	21 FCRMT(1H0,'THE ANGULAR DISTRIBUTION PROBABILITIES FOR REACTION TY	1570
52	FORMAT(1H0)	1030	*PE',2X,2A4,2X,I6,2X,' ARE GIVEN AT',I6,' ENERGY PCINTS'/)	1580
C	CALCUL DES SIGMA A PARTIR DES POLYNOMES	1040	CALL A8FORM(NO,' POINT')	1590
	X(1)=-1.0	1050	CALL A8FORM(NO,' WISE')	1600
	NPMAX=81	1060	NWORD2=NE*4+1	1610

NRIF4=NRIF4+1	1620	ISATZ(IW+4)=IR	2170
NPMAX=0	1630	ISATZ(IW+5)=NLEGI	2180
DC 41 I=1,NF	1640	IW=IW+6	2190
CALL RREC(3,NT,MODE,-1.0,NBT,JNT,X,Y)	1650	IF(IW.GE.390)CALL ERROR(701)	2200
IF(NPMAX.GE.N2)GOTO 42	1660	WRITE(NF'I'IR)(CLEG(I),I=1,NLEGI)	2210
NPMAX=N2	1670	IR=IR+NLEGI/400+1	2220
IF(N2.GT.NXMAX/2) CALL ERROR(700)	1680	WRITE(NO,51)KDR2(IKL),KCR2(IKL+1)	2230
DO 25 KI=1,N2	1690	51 FORMAT(1H1,2A4,'REACTION TYPE TABLE' /)	2240
25 TAB(KI)=X(KI)	1700	NWMI=NWORD-1	2250
42 CCNTINUE	1710	WRITE(NO,53)(CLEG(I),I=1,NWMI)	2260
CALL WREC(3,NF,2)	1720	53 FORMAT(1X,1P2E15.5,3I8)	2270
41 CCNTINUE	1730	37 ISATZ(IW+2)=KDR2(IKS)	2280
REWIND NF	1740	ISATZ(IW+3)=KDR2(IKS+1)	2290
IA=NWORD2	1750	ISATZ(IW+4)=IR	2300
WRITE(NO,28)	1760	ISATZ(IW+5)=IA	2310
28 FORMAT('1')	1770	ISATZ(IW)=MATKDK	2320
DC 24 I=1,NE	1780	ISATZ(IW+1)=1	2330
CALL RREC(3,NF,2,-1.0,NBT,JNT,X,Y)	1790	IW=IW+6	2340
E=C2	1800	IF(IW.GE.390)CALL ERROR(701)	2350
DO 26 J=1,NPMAX	1810	WRITE(NF'I'IR)(A(I),I=1,IA)	2360
CALL ITAB1(TAB(J),TAB(J+NXMAX/2))	1820	IR=IR+IA/400+1	2370
26 CONTINUE	1830	ISATZ(399)=IR	2380
DO 27 J=1,NPMAX	1840	ISATZ(400)=IW	2390
X(J)=TAB(J)	1850	WRITE(NF'I'1)(ISATZ(I),I=1,400)	2400
Y(J)=TAB(J+NXMAX/2)/6.283184	1860	NM1=NWORD2-1	2410
27 CCNTINUE	1870	WRITE(NO,51)KDR2(IKS),KDR2(IKS+1)	2420
N2=NPMAX	1880	WRITE(NO,43)(A(I),I=1,NM1)	2430
N1=1	1890	43 FORMAT(1X,1PE15.5,3I8)	2440
NBT(1)=NPMAX	1900	GOTO 3	2450
JNT(1)=2	1910	C EXTENDED DO LOOP OF 7 AND 24	2460
CALL WREC(3,NC,4)	1920	22 TAB2(NINIT2)=E	2470
NRS=(IA-1)/880+1	1930	ITAB2(NINIT2+1)=N2	2480
NWS=MOD(IA-1,880)+1	1940	ITAB2(NINIT2+2)=NRS	2490
DO 29 J=1,NPMAX	1950	ITAB2(NINIT2+3)=NWS	2500
A(IA)=X(J)	1960	NINIT2=NINIT2+4	2510
A(IA+1)=Y(J)	1970	IF(LTT-1)7,7,24	2520
IA=IA+2	1980	C	2530
IF(IA.GE.IADIM)CALL ERRCR(703)	1990	C=====INELASTIC ANGULAR DISTRIBUTIONS	2540
29 CONTINUE	2000	C	2550
GOTO 22	2010	3 CONTINUE	2560
24 CCNTINUE	2020	RETURN	2570
36 NLEGI=NLEGI-1	2030	END	2580
IA=IA-1	2040		
IF(IF4.EQ.1)GOTO 47	2050		
READ(NF'I'1)(ISATZ(I),I=1,400)	2060		
IR=ISATZ(399)	2070		
IW=ISATZ(400)	2080		
GOTO 48	2090		
47 IR=2	2100	SUBROUTINE ITAB1(XP,YP)	10
IW=1	2110	C-----CHANGED OCTOBRE 1973 BY E.STEIN.	20
48 IF(LTT.EQ.2)GOTO 37	2120	COMMON/RECS/MAT,MF,MT,C1,C2,L1,L2,N1,N2	30
ISATZ(IW)=MATKDK	2130	DIMENSION NBT(1),JNT(1),X(1),Y(1)	40
ISATZ(IW+1)=1	2140	YP=0.0	50
ISATZ(IW+2)=KDR2(IKL)	2150	IF(XP.LT.X(1))GOTO 50	60
ISATZ(IW+3)=KDR2(IKL+1)	2160	DO 10 N=2,N2	70
		IF(XP.LT.X(N))GOTO 20	80
		10 CONTINUE	90

```

IF(XP.GT.X(N2))GOTO 50
IF(X(N2-1).NE.X(N2)) GOTO 1
YP=Y(N2)
GOTO 50
1 N=N2
20 CC 30 M=1,N1
IF(N.LE.NBT(M))GOTO 40
30 CONTINUE
CALL ERROR(900)
40 I=JNT(M)
CALL TERP1(X(N-1),Y(N-1),X(N),Y(N),XP,YP,I)
50 RETURN
ENTRY EITAB1(NBT,JNT,X,Y)
GOTO 50
END

```

```

SUBROUTINE WCONT(NT,MODE,MAT,MF,NS)
C====WRITE END RECORD ON TAPE NT
IF(NT.LE.0)GOTO 20
MT=0
C1=0.0
C2=0.0
L1=0
L2=0
N1=0
N2=0
GOTO(1,1,2),MODE
C-----MODE=1
1 WRITE(NT,MAT,MF,MT,C1,C2,L1,L2,N1,N2)
GOTO 20
C-----MODE=3
2 WRITE(NT,3) MAT,MF,MT,NS
3 FORMAT(66X,I4,I2,I3,I5)
20 RETURN
END

```

```

SUBROUTINE BETSY(MIS,Y)
COMMON/RECS/NR2(8),N2
DIMENSION Y(1)
COMMON/INOUT/NR1(8),ABIS(20)
DO 1 I=1,N2
Y(I)=Y(I)*ABIS(MIS)
1 CONTINUE
RETURN
END

```

```

100 PFAL FUNCTION LEGP(X,NL,CON)
110 DIMENSION CON(1)
120 LEGP=0.5+1.5*CON(1)*X
130 TF(NL.LF.1)GOTO 30
140 C1=1.0
150 C2=X
160 DO 20 N=2,NL
170 AL=N
180 C2=((2.)*AL-1.0)**X*C2-(AL-1.0)*C1/AL
190 LEGP=LEGP+(AL+0.5)*CON(N)*C2
200 C1=C2
210 C2=C3
220 RETURN
230
240

```

```

SUBROUTINE RDAB
10 DIMENSION ISATZ(400),XSATZ(400),BK(1)
20 EQUIVALENCE (ISATZ(1),XSATZ(1))
30 COMMON/TAB/ENDTAB(25),KEDTAB(50),KEDR(177),NCNAME(20),NRIF3,NRIF4
40 DIMENSION KDR1(117),KDR2(60)
50 INTEGER ENDTAB
60 EQUIVALENCE (KEDR(1),KDR1(1)),(KEDR(118),KDR2(1))
70 COMMON/INOUT/NR1(83),NTYP,NR2(44),N
80 ENTRY RDAB(BK)
90 ND=6
100 NF=11
110 L1=1
120 N4=400
130 READ(NF,L1)(ISATZ(I),I=1,N4)
140 RETURN
150 ENTRY RDA4(MATKDK,MIS,NAME1,NAME2,LNT,IR)
160 L4=1
170 GOTO 20
180 ENTRY RDA2(MATKDK,MIS,NAME1,NAME2,LNT)
190 L4=0
200 LNT=0
210 IW=1
220 IF(ISATZ(IW).NE.MATKDK)GOTO 1
230 IF(ISATZ(IW+1).NE.MIS)GOTO 1
240 IF(ISATZ(IW+2).NE.NAME1)GOTO 1
250 IF(ISATZ(IW+3).NE.NAME2)GOTO 1
260 GOTO 5
270 1 IW=IW+6
280 IF(IW-360)3,2,2
290 3 IR=ISATZ(IW+4)
300 N=ISATZ(IW+5)
310 IF(L4.EQ.1)RETURN
320 30 READ(NF,IR)(BK(I),I=1,N)
330 RETURN
340 2 LNT=1
350 RETURN
360 ENTRY RDA3(MATKDK)
370

```

```

NK=41
IK=1
NTYP=0
13 IF(ISATZ(IW).NE.MATKDK)GOTO 10
IK=1
DO 11 ITK=1,NK
IF(ISATZ(IW+2).NE.KEDR(IK))GOTO 12
IF(ISATZ(IW+3).NE.KEDR(IK+1))GOTO 12
NTYP=NTYP+1
GOTO 10
12 IK=IK+3
11 CCNTINUE
10 IW=IW+6
IF(IW-360)13,13,14
14 RETURN
END

SUBROUTINE FILE5(MATP,IF5,BK,IBK,NBT,JNT,X,Y,B,TEMP)
COMMON/INOUT/NO,NI,NT,MODE,NR1(78),MATKDK,NR2(42),NBX,N1X,N2X,N,
*NRKX,NR3(10),EPS
COMMON/RECS/MAT,MF,MT,C1,C2,L1,L2,N1,N2
COMMON/TAB/ENDTAB(25),KEDTAB(50),KEDR(240),NCNAME(39),NRIF3,NRIF4
EQUIVALENCE (KEDR(1),KDR1(1)),(KEDR(124),KDR2(1))
INTEGER ENDTAB
REAL*8 NAM8
EQUIVALENCE (NAUX(1),NAM8)
DIMENSION NUMTYP(4),KEDTYP(4),ENERG(300),TEMP(1),IBK(1),BK(1),
*NBT(1),JNT(1),X(1),NAUX(2),KDR1(123),KDR2(1),Y(1),B(1)
DATA NUMTYP/16,17,18,91/
DATA KEDTYP/61,67,70,55/
NF=10
C-----MAXIMUM NUMBER OF POINTS ALLOWED FOR THE PRIMARY ENERGY GRID
NCMAX=300
WRITE(ND,1)
1 FORMAT('1',20X,'FILE 5 - ENERGY DISTRIBUTIONS OF SECONDARY NEUTRON
*S'//21X,51('*')//')
CALL PRIZEI
IPAGE=0
DO 2 NUM=1,4
CALL INITSC(NCMAX)
NEN=0
REWIND NF
CALL SEARCH(NT,MODE,MATP,5,NUMTYP(NUM),LNT)
IF(LNT.EQ.0)GOTO 3
WRITE(ND,4)NUMTYP(NUM)
4 FCRMAT('0ENERGY DISTRIBUTIONS FOR REACTION TYPE',I3,' NOT INCLUDED
*IN THE ENDF/B FILE'//)
BACKSPACE NT
GOTO 2
3 NK=N1
CALL WREC(1,NF,2)
C====REREAD PROBABILITY AND TEMPERATURE ARRAYS FROM ENDF/B FILE

```

```

380 C-----GENERATE AN ENERGY GRID DESCRIBING EVERY DISTRIBUTION WITH A PRECI 360
390 C-----STORE DATA ON FILE NF IN ENDF/B FORMAT 370
400 DO 5 NKN=1,NK 380
410 DO 8 NKN2=1,2 390
420 CALL RREC(3,NT,MODE,-1.0,NBT,JNT,X,Y) 400
430 CALL WREC(3,NF,2) 410
440 IF(NKN2.EQ.2)GOTO 6 420
450 LF=L2 430
460 IF(LF.EQ.3.OR.LF.EQ.7.OR.LF.EQ.9.OR.LF.EQ.10)GOTO 6 440
470 WRITE(ND,7)LF,NUMTYP(NUM) 450
480 7 FCRMAT('0LF EQUALS',I4,'. THIS CASE IS NOT FORSEEN IN THE CODE.'// 460
490 *1X,'REACTION TYPE',I4,' NOT TRANSLATED'//) 470
500 GOTO 2 480
510 C-----LF=3,7,9,10 490
520 6 CALL CHGINT(FPS,N1,N2,NBT,JNT,X,Y,N1X,N2X,830) 500
530 IF(N2.GT.NCMAX) CALL ERROR(751) 510
CALL GENER(N2,X,TEMP,FENERG,ND,NCU) 520
IF(NCU.GT.NCMAX)CALL ERROR(751) 530
IF(LF.EQ.3)GOTO 5 540
IF(LF.EQ.10)GOTO 11 550
8 CCNTINUE 560
GOTO 5 570
11 CALL RREC(2,NT,MODE,-1.0,C,B) 580
CALL WREC(2,NF,2) 590
5 CONTINUE 600
ENDFILE NF 610
REWIND NF 620
C====REREAD ENDF DATA FROM FILE NF AND CONSTRUCT ARRAY BK (DATA IN 630
C-----KEDAK FORMAT) 640
IF(IPAGE.EQ.1)WRITE(ND,31) 650
31 FCRMAT('1') 660
K2=KEDTYP(NUM) 670
NAUX(1)=KDR2(K2) 680
NAUX(2)=KDR2(K2+1) 690
CALL ABFORM(ND,NAM8) 700
IPAGE=1 710
WRITE(ND,2000)(ENERG(I),I=1,NCU) 720
2000 FCRMAT('//0ENERGY GRID GENERATED'//(1P8E15.5)) 730
C-----LENGTH OF TYPE TABLE: NLT 740
NLT=4*NCU 750
C-----INDEX FOR WRITING IN RELOCATABLE PART: ILT 760
C-----INDEX FOR WRITING DATA: IA 770
ILT=1 780
CALL RREC(1,NF,2,-1.0) 790
NK=N1 800
NKM4=4*NK 810
C-----LOOP ON SPECTRUM TYPES 820
DO 12 NKN=1,NK 830
CALL RREC(3,NF,2,-1.0,NBT,JNT,X,Y) 840
LF=L2 850
IA=NLT+1+4*(NKN-1) 860
C-----LOOP ON INCIDENT ENERGIES 870
DO 14 I=1,NCU 880
IF((IA+3).GT.NBKX) CALL ERROR(752) 890
IF(NKN.GT.1) GOTO 36 900

```

```

BK(ILT)=ENERG(I)
IRK(ILT+1)=NK
IRK(ILT+2)=(IA-1)/890+1
IRK(ILT+3)=MCO(IA-1,880)+1
ILT=ILT+4
C====SFT KEDAK DATA TYPE NUMBER OF THE DISTRIBUTION
C-----LF=3
36 IF(LF.NE.3)GOTO 22
BK(IA)=4
GOTO 35
C-----LF=7
32 IF(LF.NE.7)GOTO 33
BK(IA)=2
GOTO 35
C-----LF=9
33 IF(LF.NE.9)GOTO 34
BK(IA)=1
GOTO 35
C-----LF=10
34 BK(IA)=3
C-----INTERPOLATION IN THE PROBABILITY ARRAY
35 CALL ITAB1(ENERG(I),BK(IA+1))
IF(LF.NE.3)GOTO 15
BK(IA+2)=C2
BK(IA+3)=0.0
GOTO 14
15 IF(LF.NE.7.AND.LF.NE.9)GOTO 14
C-----LF=7 OR 9: U
BK(IA+3)=C1
14 IA=IA+NKM4
IF(LF.EQ.3)GOTO 12
IF(LF.EQ.10)GOTO 16
C-----READ THE TEMPERATURE ARRAY
CALL RREC(3,NF,2,-1.0,NBT,JNT,X,Y)
GOTO 17
C-----READ THE WATT FISSION SPECTRUM PARAMETERS
16 CALL RREC(2,NF,2,-1.0,B)
C-----LCCP ON INCIDENT ENERGIES
17 IA=IA-4*NCU*NK
DO 18 I=1,NCU
IF(LF.EQ.10)GOTO 19
C-----LF=7 OR 9: INTERPOLATE TEMPERATURE
CALL ITAB1(ENERG(I),BK(IA+2))
GOTO 18
C-----LF=10
C DIE BEIDEN FOLGENDEN STATEMENTS SIND VERMUTLICH F A L S C H .
19 BK(IA+2)=B(1)
BK(IA+3)=B(2)
18 IA=IA+NKM4
12 CCNTINUE
NCNAME(KEDTYP(NUM)/3+1)=NCU
WRITE(N0,20)KDR2(K2),KDR2(K2+1)
20 FORMAT('0',2A4,' REACTICN TYPE AND PARAMETER TABLE'/)
IA=IA-1-4*(NK-1)
ILT=ILT-1

```

91C  
920  
930  
94C  
950  
96C  
970  
980  
99C  
1000  
101C  
1020  
103C  
1040  
1050  
106C  
1070  
108C  
109C  
1100  
111C  
1120  
113C  
114C  
1150  
1160  
1170  
1180  
1190  
1200  
121C  
1220  
123C  
1240  
1250  
126C  
1270  
128C  
1290  
1300  
131C  
1320  
1330  
1340  
1350  
136C  
1370  
138C  
1390  
1400  
141C  
1420  
143C  
1440  
1450

```

DO 40 I=1,ILT,4
IP3=I+3
WRITE(N0,21)(BK(II),II=I,IP3)
21 FORMAT(1X,10F15.5,3I8)
IPA=IRK(I+3)
DO 41 J=1,NK
IPE=IPA+3
WRITE(N0,22)(BK(II),II=IPA,IPE)
22 FORMAT(10X,1P4E15.5)
41 IPE=IPE+1
40 CONTINUE
N=IA
CALL WDA2(MATKDK,1,KDR2(K2),KDR2(K2+1))
NRIF4=NRIF4+1
2 CCNTINUE
RETURN
30 CALL ERROR (750)
STOP
END

```

```

SUBROUTINE CHGINT(EPS,N1,N2,NBT,JNT,X,Y,N1X,N2X,*)
SUBROUTINE CHGINT J.DANIELS NOV. 1970
MODIFIED MAY 1972
C
C DIMENSION NBT(1),JNT(1),X(1),Y(1)
COMMON/RFXR(300),YR(300)
DC 1 I=1,N1
IF(I-1)95,96,95
96 J=1
GO TO 97
95 J=NBT(I-1)
97 LLL=NBT(I)-1
JNTX=JNT(I)
IF(JNTX.LE.0.OR.JNTX.GT.5)CALL ERROR(621)
IF(JNTX.EQ.1)GO TO 3
IF(JNTX.EQ.2) GOTO 1
28 N=2
X1=X(J)
Y1=Y(J)
X2=X(J+1)
Y2=Y(J+1)
IF(ABS(X1-X2)/X1.LE.1.0E-06)GO TO 170
CALL DAN I(X1,Y1,X2,Y2,JNTX,N,EPS)
IF(N.EQ.2) GOTO 170
KKK=J+N-1
N2OLD=N2
N2=N2+N-2
IF(N2.GT.N2X)RETURN 1
25 DC 7 KK=KKK,N2
IX1=N2-KK+KKK
IX2=N2OLD-KK+KKK
X(IX1)=X(IX2)
7 Y(IX1)=Y(IX2)

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1470  
148C  
1490  
1500  
1510  
1520  
153C  
1540  
1550  
1560  
1570  
158C  
1590  
1600  
1610  
1620  
163C  
1640

1C  
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3C  
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6C  
70  
8C  
90  
100  
11C  
120  
13C  
140  
150  
16C  
170  
18C  
190  
200  
21C  
220  
23C  
240  
25C  
26C  
270  
280  
290  
300  
31C  
320

DO 8 K=1,N	23C	IF(Y1.EQ.Y2)GOTO 1010	130
X(J+K-1)=XR(K)	340	DC 100 M=1,300	14C
Y(J+K-1)=YR(K)	350	600 XP=(XR(J)+XR(J+1))*0.5	150
8 CONTINUE	360	GOTO(1010,1010,501,502,503),JNTX	16C
DO 6 K=I,N1	370	500 GOTO 1010	170
6 NBT(K)=NBT(K)+N-2	38C	501 IF(XR(J))50,50,70	180
170 J=J+N-1	390	70 IF(XR(J+1))50,50,80	190
LLL=LLL+N-2	400	90 IF(XP)90,90,101	200
IF(J.LE.LLL)GO TO 28	410	90 CALL ERROR(134)	21C
3 JNT(I)=2	420	101 XVER=XR(J+1)/XR(J)	220
1 CONTINUE	42C	XVERM1=XVER-1.0	23C
C	440	IF(XVERM1.GT.1.0F-3.0R.XVERM1.LT.0) GOTO 2101	240
C=====ELIMINATE DISCONTINUITIES.	450	AKOR=0.125*XVERM1*XVER/(1.0-XVERM1*(C.5-0.33333333*XVERM1))	250
C	460	YP=0.5*(YR(J)+YR(J+1))-AKOR*(YR(J+1)-YR(J))	26C
L=N2-1	470	GOTO 700	270
I=1	48C	2101 YP=YR(J)+ALOG(XP/XR(J))*(YR(J+1)-YR(J))/ALOG(XVER)	280
303 K=1	490	GO TO 700	290
IF(X(I).NE.X(I+1))GO TO 302	500	502 IF(YR(J))50,50,120	300
X(I)=X(I)*0.99999	510	120 IF(YR(J+1))50,50,130	31C
IF(I.GT.(L-1)) GOTO 304	520	130 YP=SQRT(YR(J+1)*YR(J))	320
K=2	53C	GO TO 700	330
IF(X(I+1).NE.X(I+2))GO TO 302	540	503 IF(YR(J))501,501,150	340
K=3	550	150 IF(YR(J+1))501,501,160	350
X(I+2)=X(I+2)+1.0E-5*X(I+2)	56C	160 IF(XR(J))130,130,170	36C
Y(I+1)=Y(I)+(X(I+1)-X(I))*(Y(I+2)-Y(I))/(X(I+2)-X(I))	570	170 IF(XR(J+1))130,130,180	370
302 I=I+K	58C	180 IF(XP)90,90,190	38C
IF(I.GT.L)GOTO 304	590	190 XVER=XR(J+1)/XR(J)	390
GO TO 303	600	XVERM1=XVER-1.0	40C
304 CONTINUE	610	IF(XVERM1.GT.1.0E-3.0R.XVERM1.LT.0) GOTO 2190	410
C INTERPOLATIONSTABELLE VEREINFACHEN.	620	AKOR=0.125*XVERM1*XVER/(1.0-XVERM1*(C.5-0.33333333*XVERM1))	420
C DC 2 M=1,N1	63C	YP=SQRT(YR(J+1)*YR(J))*EXP(-AKOR*ALOG(YR(J+1)/YR(J)))	43C
IF(JNT(M).NE.2) CALL ERRCR(777)	640	GOTO 700	440
2 CONTINUE	650	2190 YP=YR(J)*EXP(ALOG(XP/XR(J))*ALOG(YR(J+1)/YR(J))/ALOG(XVER))	450
NBT(I)=NBT(N1)	66C	700 YLIN=0.5*(YR(J)+YR(J+1))	46C
N1=1	670	IF(YP.NE.0.)GOTO 800	47C
C FALLS Y(I)<=1.0E-20 IST, SO WIRD Y(I)=0.0 GESETZT.	680	TEST=.0	48C
IF(Y(I).LE.1.0E-20) Y(I)=0.0	690	GOTO 801	490
RETURN	70C	800 TEST=ABS((YLIN-YP)/YP)	50C
END	71C	801 IF (TEST-EPS) 1001,1001,1002	510
		1002 N=N+1	520
		KKK=N	53C
		L=J+2	540
		65 XR(KKK)=XR(KKK-1)	550
		YR(KKK)=YR(KKK-1)	560
		KKK=KKK-1	570
		IF(KKK-L)66,65,65	58C
		66 XR(KKK)=XP	590
		YR(KKK)=YP	600
		GO TO 600	610
		1001 IF(XR(M+1)-X2)1009,1010,1010	620
		1009 J=J+1	63C
		100 CONTINUE	640
		CALL ERROR(620)	650
		50 YP=0.5*(YR(J)+YR(J+1))	66C
		GO TO 700	670
SUBROUTINE DANI(X1,Y1,X2,Y2,JNTX,N,EPS)	10		
C SUBROUTINE DANI J.DANIELS NOVEMBER 1970	20		
C CHANGED BY E. STEIN IN SEPTEMBER 1973.	30		
C INTERPOLATION OF VALUES FROM ENDF LIBRARY	4C		
C COMMON/RE/XR(300),YR(300)	50		
XR(1)=X1	60		
YR(1)=Y1	70		
XR(2)=X2	80		
YR(2)=Y2	9C		
N=2	100		
J=1	110		
IF(X1.EQ.X2)GOTO 1010	120		

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1010 RETURN
END

SUBROUTINE COMB(OPER,CON,MA,MB,XL,XH,EPS,LOF,LNT)
C=====COMBINE TWO TABI FUNCTIONS=====
C THE SUBROUTINE WILL PRODUCE A TABI FUNCTION IN THE COMMON
C RECS AND IN THE ARRAYS NBT,JNT,X, AND Y.
C ACCURATE TO EPS FOR X IN THE RANGE (XL,XH). IF FA(X) IS THE
C TABI RECORD IDENTIFIED BY MA, AND FB(X) IS THE TABI RECORD
C IDENTIFIED BY MB, THE ANSWER FC(X)=FA(X) * FB(X), WHERE THE
C OPERATION * IS DEFINED IN THE FUNCTION S.R. OPER WHICH MAY
C ALSO USE CONSTANTS CON(N).
C MA = IDENT OF TABI RECORD IN DENSE STORAGE
C MB = IDENT OF TABI RECORD IN DENSE STCRAGE
C XL = LOWER LIMIT OF X REQUIRED
C XH = UPPER LIMIT OF X REQUIRED
C LOF= OVERFLOW INDICATOR NORMALLY 0. IF LCF=1, ANSWER WILL
C NOT FIT IN EXPANDED STORAGE
C EPS= RELATIVE ERROR CRITERICK
C LNT= 0, NDRMAL EXIT.
C = 1, RECORD MA OR MB NOT IN STORAGE
C-----ERROR STOP 310, XL.GE.XH
C ERROR STOP 311, MA OR MB IS ZERO
DIMENSION CON(20),NBT(1),JNT(1),X(1),Y(1)
LOGICAL*4 LODZMS
COMMON/RECS/NR1(7),N1,N2
COMMON/INOUT/ NR2(141),SCREPS
EXTERNAL OPER
C=====INITIALIZE
100 IF(XL.GE.XH)CALL ERROR(310)
IF((MA.EQ.0).OR.(MB.EQ.0))CALL ERROR(311)
N1=0
N2=0
CALL LRIDS(MA,JA,LNT)
IF(LNT.GT.0)GOTO400
CALL LRIDS(MB,JB,LNT)
IF(LNT.GT.0)GOTO400
C-----GET STARTING POINTS
NA=0
CALL IPDS(JA,NA,XL,YA1,IA1)
NA=NA+1
CALL FPDS(JA,NA,XA,YA2,IA2)
NB=0
CALL IPDS(JB,NB,XL,YB1,IB1)
NB=NB+1
CALL FPDS(JB,NB,XB,YB2,IB2)
X1=XL
C=====TEST FOR SMALLEST OF XA, XB, XH
200 IF(XA.GT.XB)GOTO220
IF(XA.EQ.XB)GOTO230
C-----XA.LT.XB
210 IF(XA.GE.XH)GOTO300
680 X2=XA
690 CALL IPDS(JB,NB-1,X2,YB3,IB3)
CALL COMBP(X1,YA1,YB1,X2,YA2,YB3,IA1,IB3,OPER,CON,EPS,LOF,
*NBT,JNT,X,Y)
IF(LOF.GT.0)GOTO400
X1=X2
YA1=YA2
IA1=IA2
NA=NA+1
YB1=YB3
IB1=IB3
CALL FPDS(JA,NA,XA,YA2,IA2)
GOTO200
C-----XA.GT.XB
220 IF(XB.GE.XH)GOTO300
X2=XB
CALL IPDS(JA,NA-1,X2,YA3,IA3)
CALL COMBP(X1,YA1,YB1,X2,YA3,YB2,IA3,IB1,OPER,CON,EPS,LOF,
*NBT,JNT,X,Y)
IF(LOF.GT.0)GOTO400
X1=X2
YB1=YB2
IB1=IB2
NB=NB+1
YA1=YA3
IA1=IA3
CALL FPDS(JB,NB,XB,YB2,IB2)
GOTO200
C-----XA.EQ.XB
230 IF(XA.GE.XH)GOTO300
X2=XA
CALL COMBP(X1,YA1,YB1,X2,YA2,YB2,IA1,IB1,OPER,CON,EPS,LOF,
*NBT,JNT,X,Y)
IF(LOF.GT.0)GOTO400
X1=X2
YA1=YA2
IA1=IA2
NA=NA+1
YB1=YB2
IB1=IB2
NP=NB+1
CALL FPDS(JB,NB,XB,YB2,IB2)
GOTO200
C=====X2.GT.XH
300 CALL IPDS(JA,NA-1,XH,YA2,IA2)
CALL IPDS(JB,NB-1,XH,YB2,IB2)
CALL COMBP(X1,YA1,YB1,XH,YA2,YB2,IA1,IB1,OPER,CON,EPS,LOF,
*NBT,JNT,X,Y)
XLC=X(1)
XFC=X(N2)
CALL CRDP(XLC,XHC,SCREPS,.FALSE.,SCREPS,DUMMY,DUMMY,LOF,NBT,JNT,X,
*Y)
C=====FINISHED
400 RETURN

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ENTRY ICOMB(NBT,JNT,X,Y)
GOTO 400
END
1050
1060
1070

SUBROUTINE ERROR(N)
C=====PRINT ERROR MESSAGE=====
COMMON/RECS/MAT,MF,MT
COMMON/INOUT/INOUT
10 WRITE(NOUT,20) N
20 FORMAT(11H0ERROR STOP,I6)
WRITE(NOUT,30) MAT,MF,MT
30 FORMAT(5H0MAT=,I5,4H MF=,I3,4H MT=,I4)
WRITE(NOUT,50)
50 FORMAT(///' ***DUMP OF THE ARRAYS NBT,JNT,X, AND Y.***'//)
CALL WREC(3,NOUT,4)
STOP 1
END
10
20
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130

SUBROUTINE MAGGY(XD,YD,A,X,Y)
DIMENSION XD(1),YD(1),X(1),Y(1)
DO 1 I=1,N
X(I)=XD(I)
Y(I)=YD(I)
1 CONTINUE
RETURN
END
10
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SUBROUTINE SEARCH(NT,MODE,MATP,MFP,MTP,LNT)
C=====SEARCH TAPE NT FOR A PARTICULAR SECTION=====
C NT - LOGICAL TAPE NUMBER
C MCCE - TAPE MODE(1,2,3)
C MATP - DESIRED MATERIAL NUMBER(SEE NOTE 2)
C MFP - DESIRED FILE NUMBER(SEE NOTE 2)
C MTP - DESIRED SECTION NUMBER(SEE NOTE 2)
C LNT = 0, SECTION LOCATED, HEAD RECORD READ AND IN /RECS/
C = 1, SECTION NOT LOCATED BECAUSE NOT ENCOUNTERED, OR
C BECAUSE DESIRED RECORD WOULD OCCUR PRIOR TO
C CURRENT POSITION OF TAPE(SEE NOTE 1).
C NOTE 1 - SEARCH IS IN FORWARD DIRECTION ONLY
C NOTE 2 - ZERO VALUES OF MATP, MFP, MTP ARE READ AS (ANY).
C THUS, MATP,MFP,0 MEANS ANY SECTION IN MATERIAL MATP,
C FILE MFP. 0,MFP,0 MEANS ANY SECTION, FILE MFP
C FOR ANY MATERIAL.
C-----ERROR STOP 317, MODE NOT IN RANGE 1-3
COMMON/RECS/MAT,MF,MT
LNT=0
10
20
30
40
50
60
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80
90
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C-----TEST MODE
IF((MODE.LT.1).OR.(MODE.GT.3))CALL ERROR(317)
GOTO(100,200,100),MODE
200
210
220
C-----MCDES 1 AND 3
230
100 CALL RREC(1,NT,MODE,-1.0)
IF(MAT.LT.0)GOTO300
IF(MAT.EQ.0)GOTO100
IF(MF.EQ.0)GOTO100
IF(MT.EQ.0)GOTO100
IF(MATP.EQ.0)GOTO110
IF(MATP-MAT)300,110,130
300
110 IF(MFP.EQ.0)GOTO120
IF(MFP-MF) 300,120,130
310
120 IF(MTP.EQ.0)GOTO400
IF(MTP-MT) 300,400,130
320
330
130 GOTO(140,140,150),MODE
340
350
140 READ(NT)MAT,MF,MT
360
370
GOTO160
150 READ(NT,155)MAT,MF,MT
380
155 FORMAT(66X,I4,I2,I3)
390
160 IF(MT.NE.0)GOTO130
400
GOTO100
410
C-----MCDE 2
420
200 CALL RREC(1,NT,2,-1.0)
IF(MF.LT.0)GOTO300
IF(MF.EQ.0)GOTO200
IF(MAT.EQ.0)GOTO200
IF(MT.EQ.0)GOTO200
IF(MFP.EQ.0)GOTO210
IF(MFP-MF)300,210,230
430
440
450
460
470
480
490
500
210 IF(MATP.EQ.0)GOTO220
IF(MATP-MAT)300,220,230
510
520
220 IF(MTP.EQ.0)GOTO400
IF(MTP-MT)300,400,230
530
540
230 READ(NT)MF,MAT,MT
550
IF(MT.NE.0)GOTO230
560
GOTO200
570
C-----SECTION NOT THERE
300 LNT=1
580
C-----FINISHED
590
400 RETURN
600
END
610

SUBROUTINE TERP1(X1,Y1,X2,Y2,X,Y,I)
C-----CHANGED OCTOBER 1973 BY E.STEIN.
10
20
C=====INTERPOLATE ONE POINT=====
C (X1,Y1) AND (X2,Y2) ARE THE END POINTS OF THE LINE
C (X,Y) IS THE INTERPOLATED POINT
C I IS THE INTERPOLATION CODE
C IF XA.LE.0 OR XB.LE.0 OR XP/XA.EQ.0 AND XB/XA.EQ.0
C DETECTED, THE INTERPOLATION CODE IS AUTOMATICALLY
C CHANGED FROM LOG TO LINEAR
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SUBROUTINE WREC(JT,NT,MODE)
C====WRITE ONE RECCRD ON ENDF/B TAPE=====
C JT = RECCRD TYPE, 1-CONT, 2-LIST, 3-TAB1, 4-TAB2, 5-HCL LIST
C 6-TPID
C NT = OUTPUT TAPE NUMBER. IF NT.LE.0, WRITE IS IGNORED.
C MODE=1 BINARY TAPE, STANDARD ARRANGEMENT.
C =2 BINARY TAPE, ALTERNATE ARRANGEMENT.
C =3 BCD CARD IMAGE TAPE FOR PUNCHING.
C =4 EXPANDED AND INTERPRETED PRINT TAPE.
C-----ERROR STOP 108, JT OUT OF RANGE 1-6
C ERROR STOP 109, MODE OUT OF RANGE 1-4
C DIMENSION NBT(1),JNT(1),X(1),Y(1),P(1)
C COMMON/RECS/MAT,MF,MT,C1,C2,L1,L2,N1,N2
C-----PRELIMINARY TESTS
100 IF(NT.LE.0)GOTO900
IF(JT.LT.1.OR.JT.GT.6) CALL ERROR(108)
IF(MODE.LT.1.OR.MODE.GT.4) CALL ERROR(109)
GOTO(200,300,400,500),MODE
C-----BINARY TAPE, STANDARD ARRANGEMENT
200 GOTO(210,220,230,240,220,245), JT
210 WRITE(NT)MAT,MF,MT,C1,C2,L1,L2,N1,N2
GOTO250
220 WRITE(NT)MAT,MF,MT,C1,C2,L1,L2,N1,N2,(B(N),N=1,N1)
GOTO250
230 WRITE(NT)MAT,MF,MT,C1,C2,L1,L2,N1,N2,(NBT(N),JNT(N),N=1,N1),
1(X(N),Y(N),N=1,N2)
GOTO250
240 WRITE(NT)MAT,MF,MT,C1,C2,L1,L2,N1,N2,(NBT(N),JNT(N),N=1,N1)
GO TO 250
245 WRITE (NT) MAT,MF,MT,(B(N),N=1,17)
250 GOTO900
C-----BINARY TAPE, ALTERNATE ARRANGEMENT
300 GOTO(310,320,330,340,320,345), JT
310 WRITE(NT)MF,MAT,MT,C1,C2,L1,L2,N1,N2
GOTO350
320 WRITE(NT)MF,MAT,MT,C1,C2,L1,L2,N1,N2,(B(N),N=1,N1)
GOTO350
330 WRITE(NT)MF,MAT,MT,C1,C2,L1,L2,N1,N2,(NBT(N),JNT(N),N=1,N1),
1(X(N),Y(N),N=1,N2)
GOTO350
340 WRITE(NT)MF,MAT,MT,C1,C2,L1,L2,N1,N2,(NBT(N),JNT(N),N=1,N1)
GO TO 350
345 WRITE (NT) MF,MAT,MT,(B(N),N=1,17)
350 GOTO900
C-----BCD CARD IMAGE TAPE
400 GOTO(420,430,440,450,460,470), JT
420 CALL PUCONT(NT)
GOTO900
430 CALL PULIST(NT,B)
GOTO900
440 CALL PUTAB1(NT,NBT,JNT,X,Y)

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GOTO900
450 CALL PUTAB2(NT,NBT,JNT)
GOTO900
460 CALL PUHOL(NT,B)
GOTO900
470 CALL PUTPID(NT,B)
GC TO 900
C-----EXPANDED AND INTERPRETED PRINT
500 GOTO(520,530,540,550,560,570), JT
520 CALL PRCONT(NT)
GOTO900
530 CALL PRLIST(NT,B)
GOTO900
540 CALL PRTAB1(NT,NBT,JNT,X,Y)
GOTO900
550 CALL PRTAB2(NT,NBT,JNT)
GOTO900
560 CALL PRHOL(NT,B)
GC TO 900
570 CALL PRTPID(NT,B)
C-----FINISHED
900 RETURN
ENTRY IWREC(NBT,JNT,X,Y,B)
GOTO 900
END

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SUBROUTINE WRECS
INTEGER BIB12
COMMON/INOUT/NO,NR1(77),BIB12
COMMON/RECS/MAT,MF,MT,C1,C2,L1,L2,N1,N2
DIMENSION ISATZ(400),XSATZ(400),NBT(1),JNT(1),X(1),Y(1),B(1)
EQUIVALENCE (ISATZ(1),XSATZ(1))
ENTRY WRECS1(NBT,JNT,X,Y,B)
K12=1
NF=12
IBIB12=1
IR=IBIB12+1
IW=1
DO 1 I=1,400
ISATZ(I)=0
1 CONTINUE
DO 4 IB=1,IBIB12
4 WRITE(NF*IB)(ISATZ(I),I=1,400)
GOTO 2000
ENTRY WRECS2(MIS,NAME1,NAME2,EL,EH)
ASSIGN 1000 TO LABEN
IF(IW.GE.400) GOTC 3
1000 READ(NF*IBIB12)(ISATZ(I),I=1,400)
ISATZ(IW)=MIS
ISATZ(IW+1)=NAME1
ISATZ(IW+2)=NAME2
XSATZ(IW+3)=EL

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XSATZ(IW+4)=FH      270
ISATZ(IW+5)=IR      280
ISATZ(IW+6)=N1      290
ISATZ(IW+7)=N2      300
WRITE(NF'IBIB12)(ISATZ(I),I=1,400) 310
IW=IW+9             320
WRITE(NF'IR)(NBT(J),JNT(J),J=1,N1),(X(N),Y(N),A=1,N2) 330
IR=IR+(2*N1+2*N2)/400+1 340
GOTO 2000           350
ENTRY WRECS4(MIS,NAME1,NAME2) 360
ASSIGN 104 TO LABEN 370
IF(IW.GE.400) GOTO 3 380
104 READ(NF'IBIB12)(ISATZ(I),I=1,400) 390
ISATZ(IW)=MIS       400
ISATZ(IW+1)=NAME1   410
ISATZ(IW+2)=NAME2   420
ISATZ(IW+3)=0        430
ISATZ(IW+4)=0        440
ISATZ(IW+5)=IR      450
ISATZ(IW+6)=N1      460
ISATZ(IW+7)=0        470
WRITE(NF'IBIB12)(ISATZ(I),I=1,400) 480
IW=IW+9             490
WRITE(NF'IR)(B(N),N=1,N1) 500
IR=IR+N2/400+1     510
2000 RETURN         520
3 IBIB12=IBIB12+1  530
I=I+1              540
IF(IBIB12.LE.BIB12) GOTO LABEN,(1000,104) 550
WRITE(NO,2)        560
2 FORMAT(1X,'TABLE OF CONTENTS OF FILE 12 IS FULL.'/ 570
1' REDUCE THE NUMBER OF REACTION TYPES TO BE TRANSLATED OR INCREASE 580
2 BIB12. THEN START THE JOB AGAIN.') 590
DO 11 I=1,IBIB12  600
WRITE(NO,10) IB    610
10 FORMAT('TABLE OF CONTENTS OF FILE 12 PART ',I2,','//'+',38('_')// 620
14X,'MIS NAME',6X,'MF MT',14X,'EL',18X,'EH',12X,'IR N1 N2 630
2'//) 640
READ(NF'IB) (ISATZ(I),I=1,400) 650
DO 11 I=1,50      660
WRITE(NO,12) (ISATZ(J+8*(I-1)),J=1,3), (ISATZ(J+8*(I-1)),J=2,3), 670
1 (ISATZ(J+8*(I-1)),J=4,8) 680
12 FORMAT(' ',I6,2X,2A4,I4,I6,4X,2E20.7,3I6) 690
11 CONTINUE       700
STOP              710
END               720

SUBROUTINE CROP(XL,XH,EPSSMIN,LOSSTE,EPSSMAX,SCRCLA,NUMSCR,LOF,NBT, 10
*JNT,X,Y) 20
C====CROP A TAB1 RECORD===== 30
C A TAB1 RECORD IN /RECS/ AND ARRAYS NBT,JNT,X, AND Y IS CROPPED 40
C ONLY THE PORTION BETWEEN X=XL AND X=XH IS RETAINED. PORTIONS OF 50

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C THE RANGE NOT DEFINED BY THE ORIGINAL TAB1 RECORD ARE SET TO ZERO. 60
C UNNECESSARY POINTS ARE ELIMINATED IF THEY CAN BE 70
C PREDICTED BY INTERPOLATION BETWEEN ADJACENT POINTS TO 80
C RELATIVE ACCURACY OF EPS. 90
C THIS EPS IS EQUAL EPSSMIN IF LOSSTE IS .FALSE. IN THIS CASE 100
C THE VARIABLES EPSSMAX,SCRCLA AND NUMSCR HAVE NO MEANING. 110
C IF LOSSTE IS .TRUE., EPS HAS ALSO NORMALLY THE VALUE EPSSMIN. 120
C IT HAS THE VALUE EPSSMAX IF TWO ADJACENT POINTS FULLFILL THE 130
C CONDITION (X2-X1)/X1<5.0E-6. 140
C IN THE CASE OF LOSSTE EQUAL .TRUE. THE NUMBER OF POINTS WHICH 150
C HAVE BEEN ELIMINATED, WILL BE COUNTED. FIRST SCRATCH-CLASSES 160
C WERE BUILT IN THAT WAY: IF EPSSMIN IS LESS EPSSMAX FIVE INTERVALS 170
C ARE MADE FROM 0.0 UP TO EPSSMIN AND ANOTHER FIVE FROM EPSSMIN TO 180
C EPSSMAX. IF EPSSMIN IS EQUAL EPSSMAX TEN INTERVALS ARE MADE FROM 190
C 0.0 TO EPSSMIN (ORE EPSSMAX). LATER THE INTERPOLATION ACCURACY 200
C OF THE CROPPED POINTS WILL BE COMPUTED AND THE NUMBER OF THIS 210
C POINTS ADDED UP IN THE CORRECT ARRAY ELEMENT NUMSCR. THE ARRAY 220
C ELEMENT NUMSCR(I) GIVES THE NUMBER OF CROPPED POINTS THAT COULD 230
C HAVE BEEN PREDICTED WITH AN RELATIVE ERROR IN THE RANGE 240
C SCRCLA(I-1) < EPS <= SCRCLA(I). (SCRCLA(0) IS EQUAL 0.0) 250
C LOF=0, NORMAL RETURN 260
C =1, RECORD TOO LARGE, REPEAT WITH LARGER EPSSMIN. 270
C-----ERROR STOP 312, TAB1 RECORD INCORRECT 280
C ERROR STOP 313, NBT, JNT TABLES TOO LARGE 290
C DIMENSION NBT(1),JNT(1),X(1),Y(1),SCRCLA(10),NUMSCR(10) 300
C LOGICAL*4 LOSSTE 310
C COMMON/RECS/MAT,MF,MT,C1,C2,L1,L2,N1,N2 320
C COMMON/INOUT/NRI(125),NBX,NIX,N2X 330
C LOF=0 340
C IF(.NOT.LOSSTE) GOTO 100 350
C IF(EPSSMIN.GT.EPSSMAX) CALL ERROR(777) 360
C BILDUNG DER SCRATCHKLASSEN. 370
C IF(EPSSMIN.EQ.EPSSMAX) GOTO 310 380
C SCRCLA(5)=EPSSMIN 390
C H1=SCRCLA(5)/5.0 400
C DO 311 I=1,4 410
C 311 SCRCLA(I)=I*H1 420
C SCRCLA(10)=EPSSMAX 430
C H1=(SCRCLA(10)-SCRCLA(5))/5.0 440
C DO 312 I=6,9 450
C 312 SCRCLA(I)=SCRCLA(5)+(I-5)*H1 460
C GOTO 315 470
C 310 H1=EPSSMIN/10.0 480
C SCRCLA(10)=EPSSMIN 490
C DO 313 I=1,9 500
C 313 SCRCLA(I)=I*H1 510
C NULLSETZEN DER ZAEHLER. 520
C DO 314 I=1,10 530
C 314 NUMSCR(I)=0 540
C ENDE DER BERECHNUNG DER SCRATCHKLASSEN. 550
C-----CROP UPPER END 560
C 100 IF(X(N2).LE.XH)GOTO200 570
C DC110N=1,N2 580
C NP=N 590
C IF(X(N).GE.XH)GOTO120 600

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110	CONTINUE	610	GCT0900	1160
120	IF(NP.LE.1)GOTO300	620	C-----ELIMINATE UNNECESSARY POINTS	1170
	DC130M=1,N1	630	C-----CHANGED JULY 1974 BY E.STEIN.	1180
	MP=M	640	C-----POINTS BETWEEN TWO POINTS ARE ONLY ELIMINATED IF ALL POINTS	1190
	IF(NP.LE.NBT(M))GOTO140	650	C-----BETWEEN CAN BE OBTAINED BY INTERPOLATION WITH AN ERROR LESS EPS.	1200
130	CONTINUE	660	400 IF(EPSMIN.LE.0.0.OR.N2.LE.2) GOTO 500	1210
	CALL ERROR(312)	670	K1=1	1220
140	I=JNT(MP)	680	401 K2=K1+2	1230
	CALL TERP1(X(NP-1),Y(NP-1),X(NP),Y(NP),XH,YP,I)	690	C LETZTES INTERVALL ERREICHT?	1240
	X(NP)=XH	700	IF(K2.GT.N2) GOTO 480	1250
	Y(NP)=YP	710	C INTERVALLANFANG GLEICH INTERVALLENDE?	1260
	N2=NP	720	IF(X(K1).EQ.X(K2)) GOTO 420	1270
	N1=MP	730	C INTERPOLATIONSART UND BEREICH FUER INTERVALLENDPUNKT.	1280
C-----	CROP LOWER END	740	402 N=K2	1290
200	IF(X(1).GE.XL)GOTO400	750	ASSIGN 403 TO LAB490	1300
	DC210N=1,N2	760	GCT0 490	1310
	NP=N	770	403 IK2=I	1320
	IF(X(N).GT.XL)GOTO220	780	MK2=M	1330
210	CONTINUE	790	KI=KI+1	1340
	GOTO300	800	C LAUFENDER PUNKT GLEICH INTERVALLANFANG ODER INTERVALLENDE?	1350
220	DC230M=1,N1	810	C ( UND Y-WERT DES LAUFENDEN PUNKTES UNGLEICH DEM Y-WERT	1360
	MP=M	820	C VON INTERVALLANFANG ODER INTERVALLENDE )	1370
	IF(NP.LE.NBT(M))GOTO240	830	IF((X(KI).EQ.X(K1).AND.Y(KI).NE.Y(K1)).OR.(X(KI).EQ.X(K2).AND.	1380
230	CONTINUE	840	*Y(KI).NE.Y(K2))) GOTO 420	1390
	CALL ERROR(312)	850	C INTERPOLATIONSART UND BEREICH FUER LAUFENDEN PUNKT.	1400
240	I=JNT(MP)	860	404 N=KI	1410
	CALL TERP1(X(NP-1),Y(NP-1),X(NP),Y(NP),XL,YP,I)	870	ASSIGN 405 TO LAB490	1420
	X(NP-1)=XL	880	GCT0 490	1430
	Y(NP-1)=YP	890	405 IKI=I	1440
	ND=NP-2	900	MKI=M	1450
	IF(ND.LE.0)GOTO400	910	C BEI AENDERN VON INTERPOLATIONSART KANN KEIN PUNKT ELIMINIERT WERD.	1460
	N2=N2-ND	920	IF(IKI.NE.IK2.OR.Y(KI).LE.1.EE-10) GOTO 420	1470
	DC250N=1,N2	930	IF(MKI.GT.MK2) GCT0 416	1480
	K=N+ND	940	CALL TERP1(X(K1),Y(K1),X(K2),Y(K2),X(KI),YP,IKI)	1490
	X(N)=X(K)	950	EPS=EPSMIN	1500
250	Y(N)=Y(K)	960	IF(LOSSTE.AND.(	1510
	DC260M=1,N1	970	*(((X(KI)-X(KI-1))/X(KI)).LT.5.0E-6).OR.	1520
260	NBT(M)=NBT(M)-ND	980	*(((X(KI+1)-X(KI))/X(KI+1)).LT.5.0E-6))) EPS=EPSMAX	1530
	MP=MP-1	990	IF(ABS((Y(KI)-YP)/Y(KI)).GT.EPS) GCT0 420	1540
	IF(ND.LE.0)GOTC280	1000	C NAECHSTER ZU ELIMINIERENDER PUNKT IM INTERVALL KI...K2.	1550
	N1=N1-ND	1010	KI=KI+1	1560
	DC270M=1,N1	1020	IF(KI.LT.K2) GOTO 404	1570
	K=M+ND	1030	C VERGRUESSERN DES INTERVALLS.	1580
	NBT(M)=NBT(K)	1040	K2=K2+1	1590
270	JNT(M)=JNT(K)	1050	IF(K2.LE.N2) GOTO 402	1600
280	GOTO400	1060	C INTERVALL WAR ZU GROSS. LETZTER PUNKT ERREICHT.	1610
C-----	ENTIRE RANGE ZERO	1070	C GENAUIGKEIT NICHT ERFUELLT. INTERVALL VERKLEINERN UND	1620
300	X(1)=XL	1080	C EVENTUELL PUNKTE ELIMINIEREN.	1630
	Y(1)=0.0	1090	420 K2=K2-1	1640
	X(2)=XH	1100	IF(K2.GT.(K1+1)) GOTO 406	1650
	Y(2)=0.0	1110	C INTERVALL NACH OBEN VERSCHIEBEN.	1660
	N2=2	1120	K1=K1+1	1670
	NBT(1)=2	1130	GOTO 401	1680
	JNT(1)=2	1140	C ELIMINIEREN DER PUNKTE K1+1 BIS K2-1	1690
	N1=1	1150	406 K1P1=K1+1	1700

	K2M1=K2-1	1710	M=1	2260
	KANZ=K2-K1-1	1720	491 IF(N.LE.NBT(M)) GOTO 492	2270
	IF(.NOT.LOSSTE) GOTO 309	1730	M=M+1	2280
C	EINORDNEN DER PUNKTE IN DIE ZUGEHÖRIGE SCRATCH-CLASSE.	1740	IF(M.LE.N1) GOTO 491	2290
	KI=K1P1	1750	CALL ERROR(312)	2300
301	IF(KI.GT.K2M1) GOTO 309	1760	492 I=JNT(M)	2310
	N=KI	1770	GOTO LAB490, (403,405,409,421,302)	2320
	ASSIGN 302 TO LAB490	1780	416 CALL ERROR(777)	2330
	GOTO 490	1790	480 CONTINUE	2340
302	IKI=I	1800	C-----ADD ZEROS TO UPPER END	2350
	CALL TERPI(X(K1),Y(K1),X(K2),Y(K2),X(KI),YP,IKI)	1810	500 IF(X(N2).GE.XH)GOTO600	2360
	EPS=ABS((Y(KI)-YP)/Y(KI))	1820	IF(N2+2.GT.N2X)GOTO700	2370
	IF(EPS.GT.SCRCLA(10)) CALL ERROR(777)	1830	X(N2+1)=X(N2)	2380
	DO 303 K=1,10	1840	Y(N2+1)=0.0	2390
	IF(K.EQ.10) GOTO 304	1850	X(N2+2)=XH	2400
	IF(EPS.LE.SCRCLA(10-K)) GOTO 303	1860	Y(N2+2)=0.0	2410
304	NUMSCR(11-K)=NUMSCR(11-K)+1	1870	N2=N2+2	2420
	GOTO 305	1880	IF(JNT(N1).NE.2)GOTO510	2430
303	CONTINUE	1890	NBT(N1)=N2	2440
305	KI=KI+1	1900	GOTO520	2450
	GOTO 301	1910	510 IF(N1+1.GT.N1X)CALL ERROR(312)	2460
309	DO 407 K=K2,N2	1920	N1=N1+1	2470
	X(K-KANZ)=X(K)	1930	NBT(N1)=N2	2480
407	Y(K-KANZ)=Y(K)	1940	JNT(N1)=2	2490
	N2=N2-KANZ	1950	520 CONTINUE	2500
	K2=K2-KANZ	1960	C-----ADD ZEROS TO LOWER END	2510
C	VERÄENDERN DER INTERPOLATIONSTABELLE.	1970	600 IF(X(1).LE.XL)GOTO800	2520
	N=K1P1	1980	IF(N2+2.GT.N2X)GOTO700	2530
	ASSIGN 409 TO LAB490	1990	DO610N=1,N2	2540
	GOTO 490	2000	K=N2-N+1	2550
409	MK1P1=M	2010	X(K+2)=X(K)	2560
	N=K2M1	2020	610 Y(K+2)=Y(K)	2570
	ASSIGN 421 TO LAB490	2030	X(1)=XL	2580
	GOTO 490	2040	Y(1)=0.0	2590
421	MK2M1=M	2050	X(2)=X(3)	2600
	IF(MK1P1.NE.MK2M1.OR.MK1P1.GT.N1) GOTO 416	2060	Y(2)=0.0	2610
	DO 410 MM=MK1P1,N1	2070	N2=N2+2	2620
410	NBT(MM)=NBT(MM)-KANZ	2080	DO620M=1,N1	2630
C	LCFESCHEN VON IDENTISCHEN EINTRÄGEN IN DER INTERPOLATIONSTABELLE.	2090	NBT(M)=NBT(M)+2	2640
	IF(N1.LE.1) GOTO 412	2100	IF(JNT(1).EQ.2)GOTO640	2650
	K=2	2110	IF(N1+1.GT.N1X)CALL ERROR(313)	2660
414	IF(NBT(K-1).NE.NBT(K)) GOTO 411	2120	DO630M=1,N1	2670
	IF(JNT(K-1).NE.JNT(K)) GOTO 416	2130	K=N1-M+1	2680
	DO 413 KK=K,N1	2140	NBT(K+1)=NBT(K)	2690
	NBT(KK-1)=NBT(KK)	2150	630 JNT(K+1)=JNT(K)	2700
413	JNT(KK-1)=JNT(KK)	2160	NBT(1)=3	2710
	N1=N1-1	2170	JNT(1)=2	2720
	GOTO 415	2180	N1=N1+1	2730
411	K=K+1	2190	640 GOTO800	2740
415	IF(K.LE.N1) GOTO 414	2200	C-----CVERFLOW	2750
412	K1=K2	2210	700 LCF=1	2760
	GOTO 401	2220	C-----FINISHED	2770
490	CONTINUE	2230	800 RETURN	2780
C	DER ZU N GEHÖRENDE INTERPOLATIONSEBEREICH M UNC DIE ZU N	2240	END	2790
C	GEHÖRENDE INTERPOLATIONSART I WIRD GESUCHT.	2250		

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SUBROUTINE LUCY(I,BK)
COMMON/INOUT/NO,NR1(127),N
DIMENSION BK(1)
IF(T.EQ.2) GOTO 2
IF(I.EQ.3) GOTO 3
CALL ERROR(203)
C RESOLVED RESONANCE FORMAT
2 WRITE(NO,5) (BK(J),J=1,N)
5 FORMAT(6E15.6/5E15.6/)
GOTO 7
C STATISTICAL DATA FORMAT
3 WRITE(NO,6) (BK(J),J=1,N)
6 FORMAT(8E15.6)
7 RETURN
END

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IF(N2.GT.N2X) CALL ERROR(105)
READ(NF*IR) (NBT(J),JNT(J),J=1,N1),(X(N),Y(N),N=1,N2)
RETURN
7 IF(N1.GT.NBX) CALL ERROR(104)
READ(NF*IR) (B(N),N=1,N1)
RETURN
2 LNT=1
RETURN
ENTRY IRRECS(NBT,JNT,X,Y,B)
RETURN
END

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SUBROUTINE RRECS(MIS,NAME1,NAME2,EL,EH,LNT)
INTEGER BIB12
COMMON/INOUT/NR1(78),BIB12,NR2(46),NBX,N1X,N2X
COMMON/RECS/MAT,MF,MT,C1,C2,L1,L2,N1,N2
DIMENSION NBT(1),JNT(1),X(1),Y(1),E(1)
DIMENSION ISATZ(400),XSATZ(400)
EQUIVALENCE (ISATZ(1),XSATZ(1))
IT=0
GOTO 6
C
C ENTRY RRECS4(MIS,NAME1,NAME2,LNT)
C
IT=1
6 NF=12
LNT=0
IBIB12=1
8 IW=1
READ(NF*IBIB12) (ISATZ(I),I=1,400)
3 IF(ISATZ(IW).NE.MIS) GOTO 1
IF(ISATZ(IW+1).NE.NAME1) GOTO 1
IF(ISATZ(IW+2).NE.NAME2) GOTO 1
GOTO 5
1 IW=IW+8
IF(IW.LT.400) GOTO 3
IBIB12=IBIB12+1
IF(IBIB12.LE.BIB12) GOTO 8
GOTO 2
5 EL=XSATZ(IW+3)
EF=XSATZ(IW+4)
IR=ISATZ(IW+5)
N1=ISATZ(IW+6)
N2=ISATZ(IW+7)
IF(IT.EQ.1) GOTO 7
IF(N1.GT.N1X) CALL ERROR(103)

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SUBROUTINE PRTPID(NT,B)
C====PRINT TPID RECORD =====
COMMON/RECS/MAT,NR1(8),NS
DIMENSION B(1)
WRITE (NT,20) (B(N),N=1,17), MAT
20 FORMAT(5HOTP ID,82X,5HLABEL/10X,16A4,A2,9X,I7)
NS=NS+1
RETURN
END

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SUBROUTINE PUTPID(NT,B)
C====PUNCH TPID RECORD =====
COMMON/RECS/MAT,NR1(8),NS
DIMENSION B(1)
10 NZ=0
WRITE (NT,20) (B(N),N=1,17), MAT,NZ,NZ,NZ
20 FORMAT(16A4,A2,I4,I2,I3,I5)
NS=NS+1
RETURN
END

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SUBROUTINE RRFC(JT,NT,MCDE,T)
C READ ONE RECORD FROM ENDF/B TAPE
C JT = RECORD TYPE, 1-CONT, 2-LIST, 3-TAB1, 4-TAB2, 5-HOL LIST
C 6-TPID
C NT = INPUT TAPE NUMBER.
C MCDE=1 BINARY TAPE, STANDARD ARRANGEMENT
C =2 BINARY TAPE, ALTERNATE ARRANGEMENT
C =3 BCD CARD IMAGE TAPE
C T = TEMPERATURE(KELVIN). IF T.GE.0 AND RECORD HAS A TEMPERATURE
C DEPENDENCE, DATA WILL BE EVALUATED AT T. IF T.LT.0, ONLY
C FIRST RECORD WILL BE READ.
C-----ERROR STOP 99, NT NOT DEFINED
C ERROR STOP 100, JT OUT OF RANGE 1-6

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C ERROR STOP 101, MODE OUT OF RANGE 1-3 140
C ERROR STOP 102, T NOT IN RANGE GIVEN IN DATA 150
C ERROR STOP 103, INTERPOLATION TABLE TOO LONG OR 0 160
C ERROR STOP 104, LIST TOO LONG OR 0 170
C ERROR STOP 105, TABULATION TOO LONG OR .LT.2 180
C ERROR STOP 106, IMPROPER TEMPERATURE DEPENDENCE 190
C-----ERROR STOP 107, MAT, MF, MT INCORRECT FOR JT=2, 3, 4, 5 200
COMMON/RECS/MAT,MF,MT,C1,C2,L1,L2,N1,N2 210
DIMENSION NBT(1),JNT(1),X(1),Y(1),B(1) 220
COMMON/INOUT/NRI(125),NBX,NIX,N2X 230
C-----PRELIMINARY TESTS 240
100 IF(INT.LE.0)CALL ERROR(99) 250
IF(JT.LT.1.OR.JT.GT.6) CALL ERROR(100) 260
IF(MODE.LT.1.OR.MODE.GT.3) CALL ERROR(101) 270
140 GOTO(200,300,400,500,570,590),JT 280
C-----JT=1, CONT RECORD 290
200 GOTO(210,220,230),MODE 300
210 READ(NT)MAT,MF,MT,C1,C2,L1,L2,N1,N2 310
GOTO240 320
220 READ(NT)MF,MAT,MT,C1,C2,L1,L2,N1,N2 330
GOTO240 340
230 READ(NT,235)C1,C2,L1,L2,N1,N2,MAT,MF,MT 350
235 FORMAT(2E11.0,4I11,I4,I2,I3) 360
240 GOTO900 370
C-----JT=2, LIST RECORD 380
300 GOTO(310,320,330),MODE 390
310 READ(NT)MAT,MF,MT,C1,C2,L1,L2,N1,N2,(B(N),N=1,N1) 400
GOTO340 410
320 READ(NT)MF,MAT,MT,C1,C2,L1,L2,N1,N2,(B(N),N=1,N1) 420
GOTO340 430
330 READ(NT,335)C1,C2,L1,L2,N1,N2,MAT,MF,MT,(B(N),N=1,N1) 440
335 FORMAT(2E11.0,4I11,I4,I2,I3/(6E11.0)) 450
340 IF(N1.GT.NBX.OR.N1.LT.1) CALL ERRCR(104) 460
IF(L1.LE.0.OR.T.LT.0.0.OR.MF.EQ.2) GOTO 800 470
NP=N1 480
DC370N=1,NP 490
370 Y(N)=B(N) 500
GOTO600 510
C-----JT=3, TAB1 RECORD 520
400 GOTO(410,420,430),MODE 530
410 READ(NT)MAT,MF,MT,C1,C2,L1,L2,N1,N2,(NBT(N),JNT(N),N=1,N1),(X(N),Y 540
1(N),N=1,N2) 550
GOTO450 560
420 READ(NT)MF,MAT,MT,C1,C2,L1,L2,N1,N2,(NBT(N),JNT(N),N=1,N1),(X(N),Y 570
1(N),N=1,N2) 580
GOTO450 590
430 READ(NT,435)C1,C2,L1,L2,N1,N2,MAT,MF,MT,(NBT(N),JNT(N),N=1,N1) 600
435 FORMAT(2E11.0,4I11,I4,I2,I3/(6I11)) 610
READ(NT,445)(X(N),Y(N),N=1,N2) 620
445 FORMAT(6E11.0) 630
450 IF(N1.GT.NIX.OR.N1.LT.1) CALL ERROR(103) 640
IF(N2.GT.N2X.OR.N2.LT.2) CALL ERRCR(105) 650
IF(L1.EQ.0.OR.T.LT.0.0.OR.MF.EQ.2) GOTO 800 660
NP=N2 670
GOTO600 680

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C-----JT=4, TAB2 RECORD 690
500 GOTO(510,520,530),MODE 700
510 READ(NT)MAT,MF,MT,C1,C2,L1,L2,N1,N2,(NBT(N),JNT(N),N=1,N1) 710
GOTO540 720
520 READ(NT)MF,MAT,MT,C1,C2,L1,L2,N1,N2,(NBT(N),JNT(N),N=1,N1) 730
GOTO540 740
530 READ(NT,535)C1,C2,L1,L2,N1,N2,MAT,MF,MT,(NBT(N),JNT(N),N=1,N1) 750
535 FORMAT(2E11.0,4I11,I4,I2,I3/(6I11)) 760
540 IF(N1.GT.NIX.OR.N1.LT.1) CALL ERRCR(103) 770
560 IF(N2.GT.N2X.OR.N2.LT.2) CALL ERROR(105) 780
GOTO 800 790
C-----JT=5, HOL LIST 800
570 GOTO(300,300,580),MODE 810
580 READ(NT,585)C1,C2,L1,L2,NCD,N2,MAT,MF,MT 820
585 FORMAT(2E11.0,4I11,I4,I2,I3) 830
N1=17*NCD 840
IF(N1.LE.0.OR.N1.GT.NBX) CALL ERROR(104) 850
READ(NT,588)(B(N),N=1,N1) 860
588 FORMAT(16A4,A2) 870
GOTO800 880
C-----JT=6, TPI0 RECORD 890
590 GOTO(591,592,593),MODE 900
591 READ(NT)MAT,MF,MT,(B(N),N=1,17) 910
GOTO 595 920
592 READ(NT)MF,MAT,MT,(B(N),N=1,17) 930
GOTO 595 940
593 READ(NT,594)(B(N),N=1,17),MAT,MF,MT 950
594 FORMAT(16A4,A2,I4,I2,I3) 960
595 GOTO 900 970
C-----TEMPERATURE DEPENDENCE 980
600 IF(L1.LT.1) CALL ERRCR(106) 990
LT=L1 1000
TA=C1 1010
IF(T.LT.TA) CALL ERROR(102) 1020
IF(T.EQ.TA)GOTO720 1030
650 GOTO(660,660,670),MODE 1040
660 READ(NT)Z,Z,Z,TB,Z,IC,Z,NP1,Z,(B(N),N=1,NP1) 1050
GOTO680 1060
670 READ(NT,675)TB,IC,NP1,(B(N),N=1,NP1) 1070
675 FORMAT(E11.0,11X,I11,11X,I11/(6E11.0)) 1080
680 LT=LT-1 1090
IF(NP1.NE.NP.OR.TB.LT.TA.OR.IC.LT.1.OR.IC.GT.5) CALL ERRCR(106) 1100
IF(T.LE.TB)GOTO700 1110
TA=TB 1120
DO690N=1,NP 1130
690 Y(N)=B(N) 1140
IF(LT.LE.0) CALL ERROR(106) 1150
GOTO650 1160
700 DO71CN=1,NP 1170
CALL TERP1(TA,Y(N),TB,B(N),T,Z,IC) 1180
710 Y(N)=Z 1190
720 GOTO(750,730,750,750),JT 1200
730 DC740N=1,NP 1210
740 B(N)=Y(N) 1220
750 IF(LT.LE.0)GOTO790 1230

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DD780L=1,LT 1240
GOTO(760,760,770),MODE 1250
760 READ(NT)Z,Z,Z,TB,Z,IC,Z,NP1,Z,(B(N),N=1,NP1) 1260
GOTO780 1270
770 READ(NT,675)TB,IC,NP1,(B(N),N=1,NP1) 1280
780 CONTINUE 1290
790 GOTO800 1300
C-----CHECK MAT, MF, AND MT FOR JT=2, 3, 4, 5 1310
800 IF(MAT.LE.0.OR.MF.LE.0.OR.MT.LE.0) CALL ERROR(1C7) 1320
C-----FINISHED 1330
900 RETURN 1340
ENTRY IRREC(NRT,JNT,X,Y,B) 1350
GOTO 900 1360
END 1370

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SUBROUTINE PRHOL(NT,B) 10
C====PRINT HOLLERITH LIST RECORD===== 20
COMMON/RECS/MAT,MF,MT,C1,C2,L1,L2,N1,N2,NS 30
DIMENSION B(1) 40
NCD=N1/17 50
10 WRITE(NT,20) C1,C2,L1,L2,NCD,N2,MAT,MF,MT,NS 60
20 FORMAT('DCOMM',9X,'C1 ',9X,'C2',11X,'L1',11X,'L2 ',9X,'NCD',11X, 70
1*N2',8X,'MAT MF MT SEQ'/E20.4,E13.4,I9,3I13,I11,I4,I5,I7) 80
NS=NS+1 90
NI=1 100
DO 40 NC=1,NCD 110
NF=NI+16 120
WRITE(NT,30) (B(N),N=NI,NF),MAT,MF,MT,NS 130
30 FORMAT(1H / (10X,16A4,A2,9X,I7,I4,I5,I7)) 140
NS=NS+1 150
40 NI=NI+17 160
RETURN 170
END 180

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SUBROUTINE PUHOL(NT,B) 10
C====PUNCH HOLLERITH LIST RECORD===== 20
COMMON/RECS/MAT,MF,MT,C1,C2,L1,L2,N1,N2,NS 30
DIMENSION F(6),S(6),J(6),B(1) 40
10 NCD=N1/17 50
CALL CXFP(C1,F(1),S(1),J(1)) 60
CALL CXFP(C2,F(2),S(2),J(2)) 70
WRITE(NT,20) F(1),S(1),J(1),F(2),S(2),J(2),L1,L2,NCD,N2,MAT,MF,MT, 80
INS 90
20 FORMAT(2(F8.5,A1,I2),4I11,I4,I2,I3,I5) 100
NS=NS+1 110
NI=1 120
DO 40 NC=1,NCD 130
NF=NI+16 140
WRITE(NT,30) (B(N),N=NI,NF),MAT,MF,MT,NS 150

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30 FORMAT(16A4,A2,I4,I2,I3,I5) 160
NS=NS+1 170
40 NI=NI+17 180
RETURN 190
END 200

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SUBROUTINE PRCONT(NT) 10
C====PRINT CONT TYPE RECORD ===== 20
COMMON/RECS/MAT,MF,MT,C1,C2,L1,L2,N1,N2,NS 30
10 IF(MAT.LT.0) GO TO 40 40
IF(MAT.EQ.0) GO TO 100 50
IF(MF.EQ.0) GO TO 80 60
IF(MT.EQ.0) GO TO 60 70
20 WRITE(NT,30)C1,C2,L1,L2,N1,N2,MAT,MF,MT,NS 80
30 FORMAT('DCONT',9X,'C1',11X,'C2',11X,'L1',11X,'L2',11X,'N1',11X, 90
1*N2',9X,'MAT MF MT SEQ'/E20.4,E13.4,I9,3I13,I11,I4,I5,I7) 100
GO TO 120 110
40 WRITE(NT,50) MAT,MF,MT,NS 120
50 FORMAT(5HOFEND,19X,11HEND OF TAPE,53X,I4,I4,I5,I7) 130
GO TO 120 140
60 WRITE(NT,70) MAT,MF,MT,NS 150
70 FORMAT(5HOFEND,4X,14HEND OF SECTION,65X,I4,I4,I5,I7) 160
GO TO 120 170
80 WRITE(NT,90) MAT,MF,MT,NS 180
90 FORMAT(5HOFEND,9X,11HEND OF FILE,63X,I4,I4,I5,I7) 190
GO TO 120 200
100 WRITE(NT,110) MAT,MF,MT,NS 210
110 FORMAT(5HOFEND,14X,15HEND OF MATERIAL,54X,I4,I4,I5,I7) 220
120 NS=NS+1 230
RETURN 240
END 250

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SUBROUTINE PUCONT(NT) 10
C====PUNCH CONT TYPE RECORD ===== 20
COMMON/RECS/MAT,MF,MT,C1,C2,L1,L2,N1,N2,NS 30
10 IF(MT.LE.0) GO TO 40 40
20 CALL CXFP(C1,F1,S1,J1) 50
CALL CXFP(C2,F2,S2,J2) 60
WRITE(NT,30)F1,S1,J1,F2,S2,J2,L1,L2,N1,N2,MAT,MF,MT,NS 70
30 FORMAT(2(F8.5,A1,I2),4I11,I4,I2,I3,I5) 80
GO TO 80 90
40 WRITE(NT,50) MAT,MF,MT,NS 100
50 FORMAT(66X,I4,I2,I3,I5) 110
80 NS=NS+1 120
RETURN 130
END 140

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```

SUBROUTINE PRLIST(NT,B)
C====PRINT LIST TYPE RECORD=====
COMMON/RECS/MAT,MF,MT,C1,C2,L1,L2,N1,N2,NS
DIMENSION B(1)
10 WRITE(NT,20) C1,C2,L1,L2,N1,N2,MAT,MF,MT,NS
20 FORMAT('OLIST',9X,'C1',11X,'C2',11X,'L1',11X,'L2',11X,'N1',11X,
1'N2',8X,'MAT MF MT SEQ'/E20.4,E13.4,I9,3I13,I11,I4,I5,I7)
NS=NS+1
WRITE(NT,30)
30 FCRMAT(' (N)',7X,'B(N)',9X,'B(N+1)',7X,'B(N+2)',7X,'B(N+3)',7X,
1'B(N+4)',7X,'B(N+5)')
NI=1
NF=6
40 IF(NF-N1) 50,50,80
50 WRITE(NT,60)NI,(B(N),N=NI,NF),MAT,MF,MT,NS
60 FORMAT(2H (,I3,2H) ,6E13.4,I7,I4,I5,I7)
NS=NS+1
IF(NF-N1) 70,200,200
70 NI=NI+6
NF=NI+5
GO TO 40
80 NE=NI-NI+1
GO TO(90,110,130,150,170),NE
90 WRITE(NT,100)NI,(B(N),N=NI,N1),MAT,MF,MT,NS
100 FORMAT(2H (,I3,2H) ,6E13.4,65X,I7,I4,I5,I7)
GO TO 190
110 WRITE(NT,120)NI,(B(N),N=NI,N1),MAT,MF,MT,NS
120 FORMAT(2H (,I3,2H) ,2E13.4,52X,I7,I4,I5,I7)
GC TO 190
130 WRITE(NT,140)NI,(B(N),N=NI,N1),MAT,MF,MT,NS
140 FCRMAT(2H (,I3,2H) ,3E13.4,39X,I7,I4,I5,I7)
GO TO 190
150 WRITE(NT,160)NI,(B(N),N=NI,N1),MAT,MF,MT,NS
160 FCRMAT(2H (,I3,2H) ,4E13.4,26X,I7,I4,I5,I7)
GO TO 190
170 WRITE(NT,180)NI,(B(N),N=NI,N1),MAT,MF,MT,NS
180 FCRMAT(2H (,I3,2H) ,5E13.4,13X,I7,I4,I5,I7)
190 NS=NS+1
200 RETURN
END

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SUBROUTINE PULIST(NT,B)
C====PUNCH LIST TYPE RECORD=====
COMMON/RECS/MAT,MF,MT,C1,C2,L1,L2,N1,N2,NS
DIMENSION F(6),S(6),J(6),B(1)
10 CALL CXFP(C1,F(1),S(1),J(1))
CALL CXFP(C2,F(2),S(2),J(2))
WRITE(NT,20)F(1),S(1),J(1),F(2),S(2),J(2),L1,L2,N1,N2,MAT,MF,MT,NS
20 FCRMAT(2(F8.5,A1,I2),4I11,I4,I2,I3,I5)
NS=NS+1
NI=1
30 NE=NI-NI+1

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KX=6
IF(NE-6) 40,50,50
40 KX=NE
50 DO 60 K=1,KX
N=NI+K-1
60 CALL CXFP(B(N),F(K),S(K),J(K))
IF(NE-6) 100,70,70
70 WRITE(NT,80)(F(K),S(K),J(K),K=1,6),MAT,MF,MT,NS
80 FCRMAT(6(F8.5,A1,I2),I4,I2,I3,I5)
NS=NS+1
IF(NE-6) 220,220,90
90 NT=NI+6
GC TO 30
100 GO TO (110,130,150,170,190),NF
110 WRITE(NT,120)(F(K),S(K),J(K),K=1,1),MAT,MF,MT,NS
120 FCRMAT(1(F8.5,A1,I2),55X,I4,I2,I3,I5)
GO TO 210
130 WRITE(NT,140)(F(K),S(K),J(K),K=1,2),MAT,MF,MT,NS
140 FCRMAT(2(F8.5,A1,I2),44X,I4,I2,I3,I5)
GC TO 210
150 WRITE(NT,160)(F(K),S(K),J(K),K=1,3),MAT,MF,MT,NS
160 FCRMAT(3(F8.5,A1,I2),33X,I4,I2,I3,I5)
GC TO 210
170 WRITE(NT,180)(F(K),S(K),J(K),K=1,4),MAT,MF,MT,NS
180 FCRMAT(4(F8.5,A1,I2),22X,I4,I2,I3,I5)
GO TO 210
190 WRITE(NT,200)(F(K),S(K),J(K),K=1,5),MAT,MF,MT,NS
200 FCRMAT(5(F8.5,A1,I2),11X,I4,I2,I3,I5)
210 NS=NS+1
220 RETURN
END

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SUBROUTINE PRTABL(NT,NBT,JNT,X,Y)
C====PRINT TABL TYPE RECORD=====
COMMON/RECS/MAT,MF,MT,C1,C2,L1,L2,N1,N2,NS
DIMENSION NBT(1),JNT(1),X(1),Y(1)
NF=N1
NP=N2
10 WRITE(NT,20)C1,C2,L1,L2,NR,NP,MAT,MF,MT,NS
20 FCRMAT('TABL',9X,'C1',11X,'C2',11X,'L1',11X,'L2',11X,'NR',11X,
1'NP',9X,'MAT MF MT SEQ'/E20.4,E13.4,I9,3I13,I12,I4,I5,I7)
NS=NS+1
WRITE(NT,30)
30 FCRMAT(' (N)',8X,'NBT(N)',7X,'JNT(N)',7X,'NBT(N+1)',5X,
1'JNT(N+1)',5X,'NBT(N+2)',5X,'JNT(N+2)')
NI=1
NF=3
40 IF(NF-NR)50,50,80
50 WRITE(NT,60)NI,(NBT(N),JNT(N),N=NI,NF),MAT,MF,MT,NS
60 FCRMAT(' (,I4,')',110,5I13,I11,I4,I5,I7)
NS=NS+1
IF(NF-NR)70,140,140

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70 NI=NI+3
NF=NI+2
GOTO 40
80 IF(NF-NR-1) 140,110,50
90 WRITE(NT,100) NI, (NBT(N), JNT(N), N=NI, NR), MAT, MF, MT, NS
100 FORMAT(' (', I4, ')', I10, I13, 52X, I11, I4, I5, I7)
GOTO 130
110 WRITE(NT, 120) NI, (NBT(N), JNT(N), N=NI, NR), MAT, MF, MT, NS
120 FCRMAT(' (', I4, ')', I10, 3I13, 26X, I11, I4, I5, I7)
130 NS=NS+1
140 WRITE(NT, 150)
150 FCRMAT('0 (N)', 7X, 'X(N)', 9X, 'Y(N)', 9X, 'X(N+1)', 7X, 'Y(N+1)',
17X, 'X(N+2)', 7X, 'Y(N+2)')
NI=1
NF=3
160 IF(NF-NP) 170, 170, 200
170 WRITE(NT, 180) NI, (X(N), Y(N), N=NI, NF), MAT, MF, MT, NS
180 FORMAT(2H (, I5, I4), 1P6E13.5, I7, I4, I5, I7)
NS=NS+1
IF(NF-NP) 190, 260, 260
190 NI=NI+3
NF=NI+2
GOTO 160
200 IF(NF-NP-1) 260, 230, 210
210 WRITE(NT, 220) NI, (X(N), Y(N), N=NI, NP), MAT, MF, MT, NS
220 FCRMAT(2H (, I5, I4), 1P2E13.5, 52X, I7, I4, I5, I7)
GOTO 250
230 WRITE(NT, 240) NI, (X(N), Y(N), N=NI, NP), MAT, MF, MT, NS
240 FORMAT(2H (, I5, I4), 1P4E13.5, 26X, I7, I4, I5, I7)
250 NS=NS+1
260 RETURN
END

SUBROUTINE PUTAB1(NT, NBT, JNT, X, Y)
C===== PUNCH TAB1 TYPE RECORD =====
COMMON/RECS/ MAT, MF, MT, C1, C2, L1, L2, N1, N2, NS
DIMENSION F(6), S(6), J(6), NBT(1), JNT(1), X(1), Y(1)
NR=NI
NP=N2
10 CALL CXFP(C1, F(1), S(1), J(1))
CALL CXFP(C2, F(2), S(2), J(2))
WRITE(NT, 20) F(1), S(1), J(1), F(2), S(2), J(2), L1, L2, NR, NP, MAT, MF, MT, NS
20 FORMAT (F8.5, A1, I2, F8.5, A1, I2, 4I11, I4, I2, I3, I5)
NS=NS+1
NI=1
NF=3
30 IF(NF-NR) 40, 40, 70
40 WRITE(NT, 50) (NBT(N), JNT(N), N=NI, NF), MAT, MF, MT, NS
50 FORMAT(6I11, I4, I2, I3, I5)
NS=NS+1
IF(NF-NR) 60, 130, 130
60 NI=NI+3

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NF=NI+2
GO TO 30
70 IF(NF-NR-1) 130, 100, 80
80 WRITE(NT, 90) (NBT(N), JNT(N), N=NI, NR), MAT, MF, MT, NS
90 FORMAT(2I11, 44X, I4, I2, I3, I5)
GO TO 120
100 WRITE(NT, 110) (NBT(N), JNT(N), N=NI, NR), MAT, MF, MT, NS
110 FCRMAT(4I11, 22X, I4, I2, I3, I5)
120 NS=NS+1
130 NI=1
NF=3
140 IF(NF-NP) 150, 150, 190
150 DO 160 K=1, 6, 2
N=NI+K/2
CALL CXFP(X(N), F(K), S(K), J(K))
160 CALL CXFP(Y(N), F(K+1), S(K+1), J(K+1))
WRITE(NT, 170) (F(K), S(K), J(K), K=1, 6), MAT, MF, MT, NS
170 FORMAT(6(F8.5, A1, I2), I4, I2, I3, I5)
NS=NS+1
IF(NF-NP) 180, 260, 260
180 NI=NI+3
NF=NI+2
GO TO 140
190 IF(NF-NP-1) 260, 220, 200
200 CALL CXFP(X(NI), F(1), S(1), J(1))
CALL CXFP(Y(NI), F(2), S(2), J(2))
WRITE(NT, 210) (F(K), S(K), J(K), K=1, 2), MAT, MF, MT, NS
210 FCRMAT(2(F8.5, A1, I2), 44X, I4, I2, I3, I5)
GO TO 250
220 DO 230 K=1, 4, 2
N=NI+K/2
CALL CXFP(X(N), F(K), S(K), J(K))
230 CALL CXFP(Y(N), F(K+1), S(K+1), J(K+1))
WRITE(NT, 240) (F(K), S(K), J(K), K=1, 4), MAT, MF, MT, NS
240 FORMAT(4(F8.5, A1, I2), 22X, I4, I2, I3, I5)
250 NS=NS+1
260 CONTINUE
RETURN
END

SUBROUTINE PRITAB2(NT, NBT, JNT)
C===== PRINT TAB2 TYPE RECORD =====
COMMON/RECS/ MAT, MF, MT, C1, C2, L1, L2, N1, N2, NS
DIMENSION NBT(1), JNT(1)
NR=NI
NP=N2
10 WRITE(NT, 20) C1, C2, L1, L2, NR, NP, MAT, MF, MT, NS
20 FORMAT('OTAB2', 9X, 'C1', 11X, 'C2', 11X, 'L1', 11X, 'L2', 11X, 'NR', 11X,
1'NP', 9X, 'MAT MF MT SEQ'/E20.4, E13.4, I9, 3I13, I12, I4, I5, I7)
NS=NS+1
WRITE(NT, 30)
30 FORMAT(' (N)', 8X, 'NBT(N)', 7X, 'JNT(N)', 7X, 'NBT(N+1)',

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15X,'INT(N+1)',5X,'NBT(N+2)',5X,'INT(N+2)')
NI=1
NF=3
40 IF(NF-NR)50,50,80
50 WRITE(NT,60)NI,(NBT(N),JNT(N),N=NI,NF),MAT,MF,MT,NS
60 FORMAT(' (',I4,')',I10,5I13,I11,I4,I5,I7)
NS=NS+1
IF(NF-NR)70,140,140
70 NT=NT+3
NF=NI+2
GOTO40
80 IF(NF-NR-1)140,110,90
90 WRITE(NT,100)NI,(NBT(N),JNT(N),N=NI,NR),MAT,MF,MT,NS
100 FORMAT(' (',I4,')',I10,I13,52X,I11,I4,I5,I7)
GOTO130
110 WRITE(NT,120)NI,(NBT(N),JNT(N),N=NI,NR),MAT,MF,MT,NS
120 FORMAT(' (',I4,')',I10,3I13,26X,I11,I4,I5,I7)
130 NS=NS+1
140 RETURN
END

SUBROUTINE PUTAB2(NT,NBT,JNT)
C=====PUNCH TAB2 TYPE RECORD =====
COMMON/RECS/MAT,MF,MT,C1,C2,L1,L2,N1,N2,NS
DIMENSION F(6),S(6),J(6),NBT(1),JNT(1)
NR=N1
NP=N2
10 CALL CXFP(C1,F(1),S(1),J(1))
CALL CXFP(C2,F(2),S(2),J(2))
WRITE(NT,20)F(1),S(1),J(1),F(2),S(2),J(2),L1,L2,NR,NP,MAT,MF,MT,NS
20 FORMAT(F8.5,A1,I2,F8.5,A1,I2,4I11,I4,I2,I3,I5)
NS=NS+1
NI=1
NF=3
30 IF(NF-NR)40,40,70
40 WRITE(NT,50)(NBT(N),JNT(N),N=NI,NF),MAT,MF,MT,NS
50 FORMAT(6I11,I4,I2,I3,I5)
NS=NS+1
IF(NF-NR)60,130,130
60 NI=NI+3
NF=NI+2
GOTO30
70 IF(NF-NR-1)130,100,80
80 WRITE(NT,90)(NBT(N),JNT(N),N=NI,NR),MAT,MF,MT,NS
90 FORMAT(2I11,44X,I4,I2,I3,I5)
GOTO120
100 WRITE(NT,110)(NBT(N),JNT(N),N=NI,NR),MAT,MF,MT,NS
110 FORMAT(4I11,22X,I4,I2,I3,I5)
120 NS=NS+1
130 RETURN
END

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SUBROUTINE CXFP(X,F,S,N)
C=====CONVERT X FOR PUNCHING =====
C X - FLOATING POINT NUMBER = F*10.0**N
C F - 0.999995 LE F LT 9.999995
C S - SIGN (HOLLERITH + OR -) OF EXPONENT
C N - EXPONENT
DATA SP/1H+/,SM/1H-/
IF(X,NE,J.0) GC TC 10
F=0.0
S=SP
N=C
RETURN
10 N=ALOG10(ABS(X))
IF(ABS(X)-1.0)40,20,20
20 F=X/10.0**N
S=SP
IF(ABS(F)-9.999995)70,30,30
30 F=F/10.0
N=N+1
GOTO70
40 N=1-N
F=X*10.0**N
S=SM
IF(ABS(F)-9.999995)70,50,50
50 F=F/10.0
N=N-1
IF(N)60,60,70
60 S=SP
70 CONTINUE
RETURN
END

SUBROUTINE XTEXT(I,J)
J=I
RETURN
END

SUBROUTINE EPT(FMIN,EMAX,NPR,NTP,NR,FR,GTR,ENS,N2X)
COMMON/RECS/NR1(9),NS
DIMENSION XP(200),XPT(100),ER(1),GTR(1),ENS(1)
CALCULATES ENERGY MESH BASED ON NPR POINTS PER RESONANCE
NTP IS CALCULATED TOTAL POINTS TO DESCRIBE NR RESONANCES FROM
FMIN TO EMAX - ENS IS CALCULATED ENERGY MATRIX
ER(J) IS RESONANCE ENERGY FOR JTH RESONANCE WHICH HAS TOTAL WIDTH
EQUAL GTR(J) AND XI(J) = GTR(J)/DELTA(J) WHERE DELTA(J) IS THE
DOPPLER WIDTH
METHOD BASED ON UNICORN CODE (REF. NAA-SR-1198C,VOL.VI)
OPTIMIZED POINTWISE REPRESENTATION OF (1+X**2)**-1 USING LOG

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C      INTERPOLATION. THE CRITERION USED IS THAT BETWEEN ANY TWO POINTS 130
C      THE DIFFERENCE BETWEEN THE EXACT INTEGRAL AND THE DIFFERENCE 140
C      ASSUMING LOG Y, LINEAR X INTERPOLATION BE WITHIN 1 PERCENT OF A 150
C      CONSTANT,  $Q = (\pi/2)(3/APR)**3$  160
C      IN THIS ROUTINE,  $X = 2(E-ER)/WIDTH$  WHERE WIDTH IS THE MAXIMUM 170
C      OF GTR OR DELTA 180
C      SET UP ENERGY POINTS XP IN VARIABLE X TO DESCRIBE A RESONANCE 190
C      PNR=NPR 200
C      CCNS=42.4115/PNR**3 210
C      XPT(1)=0.0 220
C      XPT(2)=1.817121*CCNS**(1.0/3.0) 230
C      XT=0.0 240
C      DO 25 N=1,98 250
C      CPX=1.0+XPT(N)**2 260
C      Q=CPX/(1.0+XPT(N+1)**2) 270
C      CLN=ALOG(Q) 280
C      XIN=XPT(N+1)-XPT(N) 290
C      ULO=XIN*(Q-1.0)/CLN/CPX 300
C      VXT=ATAN(XPT(N+1)) 310
C      UXT=VXT-XT 320
C      CRI=ABS(ULO-UXT)-CCNS 330
C      IF(ABS(CRI).LE.0.01*CCNS) GO TO 20 340
C      15 DELX= 350
C      1(Q-1.0)-1.0/CLN)+ABS(CRI)/XPT(N+1)) 360
C      CHECK FOR OSCILLATION ABOUT SOLUTION 370
C      DEL=DFLX+DELX1 380
C      IF(ABS(DEL)/XPT(N+1)-5.0E-6) 17,17,16 390
C      16 IF((XPT(N+1)+DELX)-XPT(N)) 17,17,19 400
C      17 DELX=DELX/2.0 410
C      GO TO 16 420
C      19 XPT(N+1)=XPT(N+1)+DELX 430
C      DFLX1=DELX 440
C      NS=N 450
C      IF(XPT(N+1)-10000.0) 10,10,30 460
C      20 XT=VXT 470
C      IF(N-1) 22,22,24 480
C      22 XPT(N+2)=XPT(N+1)**2.0 490
C      GO TO 25 500
C      24 XPT(N+2)=XPT(N+1)**2/XPT(N) 510
C      25 CONTINUE 520
C      30 NPT=2*NS-1 530
C      DO 35 I=1,NS 540
C      J=NS-I+1 550
C      XP(I)=-XP(J) 560
C      K=NPT-I+1 570
C      35 XP(K)=XP(J) 580
C      NPT1=NPT+1 590
C      DO 40 I=NPT1,200 600
C      40 XP(I)=0.0 610
C      FIND RESONANCE NO. MM NEAREST TO EMIN TO START ENERGY CALCULATION 620
C      ENS(1)=EMIN 630
C      NRM=MAX0(NR-1,1) 640
C      DO 45 MM=1,NRM 650
C      IF(2.0*ENS(1).LT.(ER(MM)+ER(MM+1))) GO TO 46 660
C      45 CONTINUE 670

```

```

MM=NRM 680
C      FIND POINT N OF XP ARRAY FOR THE MMTH RESONANCE WHICH LIES JUST 690
C      BELOW EMIN IN ABSOLUTE ENERGY 700
C      46 WIDTH=GTR(MM) 710
C      XI=2.*(ENS(1)-ER(MM))/WIDTH 720
C      DO 47 K=1,NPT 730
C      N=MAX0(K-1,1) 740
C      IF(XI-XP(K)) 48,47,47 750
C      47 CONTINUE 760
C      CALCULATE ENERGY POINTS 770
C      49 IF(MM.LT.NR)WIDTH1=GTR(MM+1) 780
C      I=1 790
C      GO TO 52 800
C      50 IF(ENS(I).LE.ENS(1))GO TO 51 810
C      52 I=I+1 820
C      51 IF(ENS(I-1).GE.FMAX) GO TO 96 830
C      N=N+1 840
C      IF(I.GT.N2X) CALL RESERR(105,MM,NR,2) 850
C      ENS(I)=XP(N)*WIDTH/2.0+ER(MM) 860
C      IF(MM=NR) 93,87,96 870
C      87 IF(N=NPT) 50,50,96 880
C      93 XMD=2.0*(ER(MM+1)-ER(MM))/(WIDTH+WIDTH1) 890
C      IF(N.GT.NPT) GO TO 89 900
C      IF(XP(N).LT.XMD) GO TO 50 910
C      89 MM=MM+1 920
C      WIDTH=WIDTH1 930
C      IF(MM.LT.NR)WIDTH1=GTR(MM+1) 940
C      DO 91 K=1,NPT 950
C      N=MAX0(K-1,1) 960
C      IF(-XMD-XP(K)) 92,91,91 970
C      91 CONTINUE 980
C      92 IF(I.GT.N2X) CALL RESERR(105,MM,NR,2) 990
C      ENS(I)=-XMD*WIDTH/2.0+ER(MM) 1000
C      GO TO 50 1010
C      96 NTP=I-1 1020
C      IF(NTP.GT.N2X) CALL RESERR(105,MM,NR,2) 1030
C      ENS(NTP)=FMAX 1040
C      RETURN 1050
C      END 1060

SUBROUTINE SCALE 10
C      SCALE GENERATES COMPLETE SCALE IF IT IS SHORTER THAN NCMAX IT 20
C      IS KEPT IN MAIN STORAGE, OTHERWISE DATASETS NIN,NOUT ARE USED. 30
C      40
C      50
C      DIMENSION E(1),TEMP(1),EC(1) 60
C      DATA NIN,NOUT/3,4/ 70
C      ENTRY INITSC(NCMAX) 80
C      NC=0 90
C      ND=0 100
C      NDS=0 110
C      RETURN 120

```

C	ENTRY GENER(MAX,E,TEMP,EC,ND,NCU)	130	IF(KC.LE.NCMAX) GOTO 36	680
	IF(NC.NE.0) GOTO 10	140	WRITE(NOUT) NCMAX,(TEMP(I),I=1,NCMAX)	690
C		150	KC=1	700
C	FIRST ENTRY TO GENER.	160	NEWNC=NEWNC+NCMAX	710
	IF(MAX.LE.NCMAX) GOTO 4	170	36 TEMP(KC)=EC(IC)	720
	REWIND NOUT	180	IC=IC+1	730
	DC 2 M=1,MAX,NCMAX	190	IF(IC.LE.LDAT) GOTO 26	740
	ME=MINJ(M+NCMAX-1,MAX)	200	ICDONE=ICDONE+LDAT	750
	LCAT=ME-M+1	210	IF(ICDONE.GE.NC) GOTO 60	760
2	WRITE(NOUT) LDAT,(E(I),I=M,ME)	220	READ(NIN) LDAT,(EC(I),I=1,LDAT)	770
	REWIND NOUT	230	IC=1	780
	NCS=NOUT	240	GOTO 26	790
	NOUT=NIN	250		800
	NIN=NDS	260	C EC>E	810
	GOTO 6	270	C CHECK IF E(M) IS NEAR BY TEMP(KC).	820
4	DO 5 M=1,MAX	280	40 IF(KC.GT.0.AND.E(M).LE.TEMP(KC)*1.CCCC10) GOTO 43	830
5	EC(M)=E(M)	290	KC=KC+1	840
6	NC=MAX	300	IF(KC.LE.NCMAX) GOTO 42	850
	ND=NDS	310	WRITE(NOUT) NCMAX,(TEMP(I),I=1,NCMAX)	860
	NCU=NC	320	KC=1	870
	RETURN	330	NEWNC=NEWNC+NCMAX	880
C		340	42 TEMP(KC)=E(M)	890
C	FURTHER ENTRIES.	350	43 M=M+1	900
10	ICDONE=0	360	IF(M.LE.MAX) GOTO 26	910
	NEWNC=0	370		920
	KC=0	380	C	930
	IF(NC.LE.NCMAX) GOTO 19	390	C	940
	REWIND NIN	400	END OF DATA FOR E.	950
	REWIND NOUT	410	50 DO 52 J=IC,LDAT	960
C		420	KC=KC+1	970
C	ENERGIES ) E(I) ARE SIMPLY COPIED.	430	IF(KC.LE.NCMAX) GOTO 52	980
12	READ(NIN) LDAT,(E(I),I=1,LDAT)	440	WRITE(NOUT) NCMAX,(TEMP(I),I=1,NCMAX)	990
	IF(E(I).LT.EC(LDAT)*1.000001) GOTO 20	450	KC=1	1000
	IF(LDAT.LT.NCMAX) GOTO 20	460	NEWNC=NEWNC+NCMAX	1010
	WRITE(NOUT) LDAT,(E(I),I=1,LDAT)	470	52 TEMP(KC)=EC(J)	1020
	ICDONE=ICDONE+LDAT	480	ICDONE=ICDONE+LDAT	1030
	NEWNC=ICDONE	490	IF(ICDONE.GE.NC) GOTO 70	1040
	IF(ICDONE.GE.NC) GOTO 60	500	READ(NIN) LDAT,(E(I),I=1,LDAT)	1050
	GOTO 12	510	IC=1	1060
C		520	GOTO 50	1070
19	LDAT=NC	530		1080
20	IC=1	540	C	1090
	M=1	550	C	1100
C		560	END OF DATA FOR EC.	1110
C	INSERT NEW ENERGIES IF NECESSARY.	570	60 DO 62 J=M,MAX	1120
26	IF(EC(IC).GT.E(M)*1.000010) GOTO 40	580	KC=KC+1	1130
	IF(EC(IC).LT.E(M)*0.99999) GOTO 30	590	IF(KC.LE.NCMAX) GOTO 62	1140
C		600	WRITE(NOUT) NCMAX,(TEMP(I),I=1,NCMAX)	1150
C	EC=E	610	KC=1	1160
C	M=M+1	620	NEWNC=NEWNC+NCMAX	1170
	IF(M.GT.MAX) GOTO 50	630	62 TEMP(KC)=E(J)	1180
C		640		1190
C	EC<E	650	C	1200
30	KC=KC+1	660	C	1210
		670	70	1220
			WRITE LAST BLOCK AND UPDATE COUNTERS.	
			NC=NEWNC+KC	
			IF(NC.LE.NCMAX) GOTO 72	
			WRITE(NOUT) KC,(TEMP(I),I=1,KC)	
			NDS=NOUT	
			NOUT=NIN	
			NIN=NDS	

```

REWIND NIN
REWIND NOUT
GOTO 100
72 DO 73 I=1,KC
73 EC(I)=TEMP(I)
100 ND=NDS
ACU=NC
RETURN
END

SUBROUTINE COMBP(X1P,YA1P,YB1P,X2P,YA2P,YB2P,IAF,IBP,CFER,CON,EPSP
1,LCF,NBT,JNT,X,Y)
C====COMBINE TAB1 FUNCTIONS A AND B IN A SINGLE PANEL=====
C (DELETE THE TERMINAL P ON ITEMS IN LIST ABOVE. I.E.,READ X1P AS X1
C I)
C X1 AND X2 ARE THE X LIMITS OF THE PANEL.
C YA1 AND YA2 ARE THE VALUES OF FUNCTION A AT X1 AND X2.
C YB1 AND YB2 ARE THE VALUES OF FUNCTION B AT X1 AND X2.
C IA AND IB ARE THE INTERPOLATION CODES FOR FUNCTIONS A AND B.
C OPER IS THE NAME OF THE OPERATION TO BE PERFORMED.
C YC=OPER(YA,YB,CON)
C CCA(N) IS AN ARRAY OF CONSTANTS REQUIRED BY FUNCTION OPER.
C EPS IS THE RELATIVE ERROR CRITERION.
C THE ANSWER FUNCTION IS ADDED TO THE ANSWER FUNCTION FROM
C PREVIOUS PANELS ALREADY STORED IN COMMON/RECS/ AND
C ARRAYS NBT,JNT,X, AND Y.
C LCF=0, NORMAL RETURN.
C =1, ANSWER OVERFLOWS /RECS/.
C-----NC ERROR STOPS
COMMON/RECS/MAT,MF,MT,C1,C2,L1,L2,N1,N2
DIMENSION NBT(1),JNT(1),X(1),Y(1)
COMMON/INOUT/NR1(124),ICOMB,NBX,NIX,N2X,NR2(22),FAICOD
C====ICOMB=0 THE INTERPOLATION SCHEME IS CHOSEN BY THIS ROUTINE
C =1 THE INTERPOLATION SCHEME IS ALWAYS LINEAR-LINEAR
EXTERNAL OPER
C====INITIALIZE
100 LCF=0
X1=X1P
YA1=YA1P
YB1=YB1P
X2=X2P
YA2=YA2P
YB2=YB2P
IA=IAP
IB=IBP
EPS=EPSP
IFX=0
EPPS=1.E-6
IF(IA.EQ.1)YA2=YA1
YA2E=YA2
IF(IB.EQ.1)YB2=YB1
YB2F=YB2

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1240
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1260
127C
1280
1290
1300
1310
10
20
30
40
5C
60
7C
80
9C
100
110
12C
130
140
15C
160
17C
180
19C
200
210
22C
230
24C
250
260
270
280
29C
300
310
32C
330
340
350
360
37C
380
39C
400
410
42C

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C-----CALCULATE VALUES AT END POINTS
YC1=OPER(YA1,YB1,CON)
110 YC2=OPER(YA2,YB2,CON)
C-----TEST FOR DISCONTINUITY
IF(X1.NE.X2)GOTO201
IC=2
IFX=1
GOTO500
C-----TEST FOR EQUALITY OF X1 AND X2 UP TO 6 SIGNIFICANT DIGITS
201 IF(ABS((X1-X2)/X1).GT.EPPS)GOTO 200
TEX=1
IC=2
GOTO 500
C====CALCULATE INTERMEDIATE VALUE, (X3,YC3)
200 X3=0.5*(X1+X2)
IF(ABS((X3-X1)/X2).GT.EPPS)GOTO 202
GOTO 900
202 CALL TERP1(X1,YA1,X2,YA2,X3,YA3,IA)
CALL TERP1(X1,YB1,X2,YB2,X3,YB3,IB)
YC3=OPER(YA3,YB3,CON)
C====TEST TO SEE IF YC3 CAN BE OBTAINED BY INTERPOLATION
IF(YC1.LE.0.0)GOTO310
IF(YC2.LE.0.0)GOTO310
IF(X1.LE.0.0)GOTO320
IF(X2.LE.0.0)GOTO320
GOTO340
C-----LINEAR INTERPOLATION ONLY
310 YC4=0.5*(YC1+YC2)
IF(YC3.EQ.YC4) GO TO 335
EPS4=ABS(YC4-YC3)*3./ (ABS(YC1)+ABS(YC2)+ABS(YC3))
GOTO330
320 YC4=0.5*(YC1+YC2)
EPS4=ABS(YC4-YC3)/YC3
330 IF(EPS4.GT.EPS)GOTO400
335 IC=2
GOTO500
C-----LAST INTERPOLATION CODE EPS4 AND BEST INTERPOLATION CODE EPS5
C-----IS PRODUCED. IF EPS4.LE.FAICOD*EPS5 THE LAST INTERPOLATION CODE
C-----IS USED. IN THE OTHER CASE THE BEST INTERPOLATION CODE IS USED.
340 EPS5=1.0E+10
EPS4=1.0E+10
NFIN=5
IF(ICOMB.EQ.1) NFIN=2
IF(N1.GE.1) IC4=JNT(N1)
DO 360 N=2,NFIN
CALL TERP1(X1,YC1,X2,YC2,X3,YC4,N)
EPS6=ABS(YC4-YC3)/YC3
IF(N1.GE.1.AND.IC4.EQ.N) EPS4=EPS6
IF(EPS6.GE.EPS5)GOTO360
EPS5=EPS6
IC5=N
360 CONTINUE
IF(EPS5.GT.EPS) GOTO 400
IF(N1.LT.1.OR.EPS4.GT.EPS) GOTO 341
IF(EPS4.LE.FAICOD*EPS5) GOTO 342

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341 IC=IC5
    GOTO 500
342 IC=IC4
    GOTO 500
C====REJECT VALUE, HALVE INTERVAL, AND TRY AGAIN
4CC X2=X3
    YA2=YA3
    YB2=YB3
    YC2=YC3
    GOTO200
C====OK, ACCEPT VALUE YC2
500 IF(N2.GT.0)GOTO510
    X(1)=X1
    Y(1)=YC1
    X(2)=X2
    Y(2)=YC2
    N2=2
    NBT(1)=2
    JNT(1)=IC
    N1=1
    GOTO700
510 N2=N2+1
    IF(N2.GT.N2X)GOTC800
    X(N2)=X2
    Y(N2)=YC2
    IF(IC.EQ.JNT(N1)) GC TO 520
    N1=N1+1
    IF(N1.GT.N1X)GOTC800
    JNT(N1)=IC
    NBT(N1)=N2
    GOTO700
520 NBT(N1)=N2
    IF(1EX.EQ.1)GOTC900
C====TEST TO SEE IF PREVIOUS POINT CAN BE ELIMINATED
600 IF(X1.NE.X2)GOTO610
    IF(N1.LE.1)GOTO900
    IF(NBT(N1-1).NE.N2-1)GOTO900
    IF(JNT(N1-1).NE.1)GOTO900
    X(N2-1)=X(N2)
    Y(N2-1)=Y(N2)
    N2=N2-1
    N1=N1-1
    NBT(N1)=N2
    GOTO900
610 CCNTINUE
C FOLLOWING STATEMENTS DELETED JULY 1974 BY E.STEIN.
C THE STATEMENTS WILL BE REPLACED BY A CROP-CALL IN COMB.
C 610 CALL TERP1(X(N2-2),Y(N2-2),X(N2),Y(N2),X(N2-1),YC6,IC)
C EPS6=ABS(YC6-Y(N2-1))
C IF(EPS6.EQ.0)GOTO620
C IF(Y(N2-1).EQ.0)GOTO700
C EPS6=EPS6/ABS(Y(N2-1))
C IF(EPS6.GT.EPS)GOTO700
C 620 X(N2-1)=X(N2)
C Y(N2-1)=Y(N2)

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1010
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1080
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1100
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1120
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1190
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1500
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1520

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C N2=N2-1
C NBT(N1)=N2
C====SEE IF MORE POINTS ARE NEEDED
700 IF(1EX.EQ.1)GOTO900
    IF(X2.LT.X2P)GOTC720
    IF(YA2E.NE.YA2P)GOTC710
    IF(YB2E.EQ.YB2P)GOTO900
710 YA2F=YA2P
    YB2F=YB2P
720 X1=X2
    YA1=YA2
    YB1=YB2
    YC1=YC2
    X2=X2P
    YA2=YA2E
    YB2=YB2E
    GOTO110
C====OVERFLOW
800 LCF=1
C====FINISHED
900 RETURN
END

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1600
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1700
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1730
1740

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SUBROUTINE DENISE(MATP,LRP,LISRES,LISUNR,NIS,LFW,LNU,BK,NBT,JNT,
*X,Y,B,A)
C CMGN/DENS/NR3(402),MX
C CMGN/INDUT/NO,NR1(5),LEDITS,NR2(122),NBKX,NR4(10),FPSR
DIMENSION BK(1),NBT(1),JNT(1),X(1),Y(1),B(1),A(1)
EPS=FPSR
NCMAX=NBKX/2
CALL TICUMB(NBT,JNT,X,Y,A,A)
CALL IFFTCH(NBT,JNT,X,Y,B,A,A)
CALL ISTORE(NBT,JNT,X,Y,B,A,A)
CALL IFPDS(A,A)
CALL IIPDS(A,A)
CALL IRRFCS(NBT,JNT,X,Y,B)
MX=100
CALL DELETE(,A)
IF(LISRES.EQ.1.AND.LISUNR.EQ.1)GOTC1
IF(LISRES.EQ.1)GOTO2
CALL MYRIAM(NIS,LISRES,LISUNR,4HSCAT,EL,EH,EPS,C,Y)
CALL WRECS2(,4HSGN,4HPRES,EL,EH)
20 IF(LISUNR.EQ.1)GOTO21
CALL MYRIAM(NIS,LISRES,LISUNR,4HSCAT,EL,EH,EPS,1,Y)
CALL WRECS2(,4HSGN,4HSTAT,EL,EH)
21 IF(LISRES.EQ.1)GOTO22
CALL MYRIAM(NIS,LISRES,LISUNR,4HCAPT,EL,EH,EPS,C,Y)
CALL WRECS2(,4HSGG,4HRES,EL,EH)
22 IF(LISUNR.EQ.1)GOTO23
CALL MYRIAM(NIS,LISRES,LISUNR,4HCAPT,EL,EH,EPS,1,Y)
CALL WRECS2(,4HSGG,4HSTAT,EL,EH)
23 IF(LFW.NE.1)GOTC2

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IF(LISRES.EQ.1)GOTO 24
CALL MYRTAM(NIS,LISRES,LISUNR,4HFISS,EL,EH,EPS,C,Y)
CALL WRECS2(0,4HSGF,4HRES,EL,EH)
24 IF(LISUNR.EQ.1)GOTO 2
CALL MYRTAM(NIS,LISPES,LISUNR,4HFISS,EL,EH,EPS,1,Y)
CALL WRECS2(0,4HSGF,4HSTAT,EL,EH)
2 CALL XTEXT(4HRES,NAME2)
IF(LISRES.EQ.1)GOTO 10
CALL MARION(LFW,EL,EH,EPS,NAME2)
IF(LEBITS.NE.1)GOTO 947
WRITE(ND,3)
3 FORMAT(1H1,'POINT WISE TOTAL CROSS SECTION CALCULATED FROM',
* ' RESOLVED RESONANCE PARAMETERS'//)
CALL WREC(3,NC,4)
947 CONTINUE
CALL WRECS2(0,4HSGT,4HRES,EL,EH)
10 IF(LISUNR.EQ.1)GOTO 1
CALL XTEXT(4HSTAT,NAME2)
EPS=EPSR
CALL MARION(LFW,EL,EH,EPS,NAME2)
IF(LEBITS.NE.1)GOTO 948
WRITE(ND,4)
4 FORMAT(1H1,'POINT WISE TOTAL CROSS SECTION CALCULATED FROM',
* ' STATISTICAL RESONANCE PARAMETERS'//)
CALL WREC(3,ND,4)
948 CONTINUE
CALL WRECS2(0,4HSGT,4HSTAT,EL,EH)
C-----FILE 3
1 CALL SOPHIE(MATP,LISRES,LISUNR,LFW,LANU,EPSR,BK,PK,BK(1+NCMAX),
*NCMAX,NBT,JNT,X,Y,B,B)
RETURN
END

SUBROUTINE FETCH(MA,LNT)
C=====FETCH RECORD FROM DENSE STORAGE=====
C MA - RECORD IDENT TO BE FETCHED
C LNT=0, RECORD FOUND AND MOVED TO COMMON/RECS/ AND
C ARRAYS NBT,JNT,X, AND Y.
C =1, RECORD NOT IN DENSE STORAGE
C-----NO ERROR STOPS
COMMON/RECS/MAT,MF,MT,C1,C2,L1,L2,N1,N2
COMMON/DENS/JMT(100),JAT(100),JTT(100),JLT(100),
1 JNS,MNS,MX
COMMON/INDUT/ NR1(125),NBX,N1X,N2X
DIMENSION LA(1),ABT(1),JNT(1),X(1),Y(1),B(1),A(1)
C-----SEARCH FOR RECORD
100 LNT=0
IF(MNS.LE.1)GOTO120
MNSP=MNS-1
DO110M=1,MNSP
MP=M
IF(JMT(M).EQ.MA)GOTO130

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600
610

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110 CONTINUE
120 LNT=1
GOTO300
130 JA=JAT(MP)
JT=JTT(MP)
IF(JT.EQ.6)GOTO210
C-----MOVE FIRST 9 WORDS
MAT=LA(JA)
MF=LA(JA+1)
MT=LA(JA+2)
C1=A(JA+3)
C2=A(JA+4)
L1=LA(JA+5)
L2=LA(JA+6)
N1=LA(JA+7)
N2=LA(JA+8)
JA=JA+9
C-----TEST IF RECORD IS TOO LONG.
IF((JT.EQ.2.OR.JT.EQ.5).AND.N1.GT.NBX) CALL ERRCR(104)
IF((JT.EQ.3.OR.JT.EQ.4).AND.N1.GT.N1X) CALL ERRCR(103)
IF(JT.EQ.3.AND.N2.GT.N2X) CALL ERRCR(105)
C-----MOVE REMAINDER OF RECORD
GOTO(300,140,160,190,140),JT
140 DC150N=1,N1
J=JA+N-1
150 B(N)=A(J)
GOTO300
160 DO170N=1,N1
J=JA+2*(N-1)
NRT(N)=LA(J)
170 JNT(N)=LA(J+1)
JA=JA+2*N1
DO180N=1,N2
J=JA+2*(N-1)
X(N)=A(J)
180 Y(N)=A(J+1)
GOTO300
190 DC200N=1,N1
J=JA+2*(N-1)
NRT(N)=LA(J)
200 JNT(N)=LA(J+1)
GOTO300
C-----TIPID RECORD
210 MAT=LA(JA)
MF=LA(JA+1)
MT=LA(JA+2)
IF(NBX.LT.17) CALL ERROR(104)
DO220N=1,17
J=JA+2*N
220 B(N)=A(J)
C-----FINISHED
300 RETURN
ENTRY IFETCH(NRT,JNT,X,Y,B,A,LA)
GOTO 300
END

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SUBROUTINE STORE(JT,MA,LOF)
C====STORE RECORD IN DENSE STORAGE=====
C JT = RECORD TYPE, 1-CONT, 2-LIST, 3-TAB1, 4-TAB2, 5-HOL, 6-TPID
C MA = IDENT OF RECORD TO BE STORED. IF A RECORD WITH
C THE SAME IDENT IS ALREADY STORED, IT IS DELETED.
C LOF= OVERFLOW INDICATOR NORMALLY ZERO. IF LOF=1, RECORD
C WILL NOT FIT.
C-----ERROR STOP 300, JT NOT IN RANGE 1-6
C ERROR STOP 301, MA=C NOT ALLOWED
COMMON/DENS/JMT(100),JAT(100),JTT(100),JLT(100),
1 JNS,MNS,MX
COMMON/INOUT/NR1(130),JX
DIMENSION LA(1),NBT(1),JNT(1),X(1),Y(1),B(1),A(1)
COMMON/RECS/MAT,MF,MT,C1,C2,L1,L2,N1,N2
C-----INITIALIZE ANT TEST JT AND MA
100 LOF=0
IF(MA.EQ.0)CALL ERROR(301)
IF((JT.LT.1).OR.(JT.GT.6))CALL ERROR(300)
C-----TEST FOR RECORD WITH SAME IDENT
110 IF(MNS.LE.1)GOTO130
MNSP=MNS-1
DC120M=1,MNSP
IF(JMT(M).EQ.MA) CALL DELETE(MA,A)
120 CONTINUE
IF(MNS.GT.MX)GOTO320
C-----TEST FOR TOO MANY VALUES
130 GOTO(140,150,160,170,150,180),JT
140 JL=9
GOTO190
150 JL=9+N1
GOTO190
160 JL=9+2*(N1+N2)
GOTO190
170 JL=9+2*N1
GOTO190
180 JL=20
190 IF((JNS+JL-1).GT.JX)GCTC320
C-----MOVE FIRST 9 WORDS
200 LA(JNS)=MAT
LA(JNS+1)=MF
LA(JNS+2)=MT
A(JNS+3)=C1
A(JNS+4)=C2
LA(JNS+5)=L1
LA(JNS+6)=L2
LA(JNS+7)=N1
LA(JNS+8)=N2
C-----MOVE REST OF RECORD
210 GCTC(310,220,240,270,220,290),JT
220 DC230N=1,N1
J=JNS+8+N

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230 A(J)=B(N)
GOTO310
240 DC250N=1,N1
J=JNS+7+2*N
LA(J)=NBT(N)
250 LA(J+1)=JNT(N)
JF=JNS+7+2*N1
DC260N=1,N2
J=JP+2*N
A(J)=X(N)
260 A(J+1)=Y(N)
GCTO310
270 DC280N=1,N1
J=JNS+7+2*N
LA(J)=NBT(N)
280 LA(J+1)=JNT(N)
GCTO310
290 DC300N=1,17
J=JNS+2+N
300 A(J)=B(N)
C-----UPDATE TABLES AND COUNTERS
310 JMT(MNS)=MA
JTT(MNS)=JT
JAT(MNS)=JNS
JLT(MNS)=JL
MNS=MNS+1
JNS=JNS+JL
GCTO400
C-----OVERFLOW
320 LCF=1
C-----FINISHED
400 RETURN
ENTRY ISTORE(NBT,JNT,X,Y,B,A,LA)
GOTO 400
END

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SUBROUTINE DELETE(MA,A)
C====DELETE RECORD MA FROM DENSE STORAGE=====
C IF MA=0, CLEAR DENSE STORAGE
C-----NO ERROR STOPS
COMMON/DENS/JMT(100),JAT(100),JTT(100),JLT(100),
1 JNS,MNS,MX
COMMON/INOUT/NR1(130),JX
DIMENSION A(1)
C=====TEST MA
100 IF(MA.NE.0)GOTO200
C=====CLEAR DENSE STORAGE
MNS=1
JNS=1
DC110M=1,MX
JMT(M)=0
JAT(M)=0

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      JLT(M)=0
110 JTT(M)=0
      DC120J=1,JX
120 A(J)=0.0
      GCT0300
C====DELTE RECORD MA, SEARCH FOR LOCATION
200 IF(MNS.LE.1)GOTO300
      MNSP=MNS-1
      DC210M=1,MNSP
      MP=M
      IF(JMT(M).EQ.MA)GOTO220
210 CCNTINUE
      GOTO300
220 JA=JAT(MP)
      JL=JLT(MP)
C-----CLOSE UP ARRAY A(J)
      JP=JNS-JA-JL
      IF(JP.LE.0)GOTO240
      DO230N=1,JP
      J1=JA+N-1
      J2=J1+JL
230 A(J1)=A(J2)
240 JNS=JNS-JL
      DO250N=1,JL
      J=JNS+N-1
250 A(J)=0.0
C-----CLOSE UP J TABLES
      MXP=MNS-MP-1
      IF(MXP.LE.0)GOTO270
      DC260M=1,MXP
      M1=MP+M-1
      JMT(M1)=JMT(M1+1)
      JTT(M1)=JTT(M1+1)
260 JLT(M1)=JLT(M1+1)
270 MNS=MNS-1
      JMT(MNS)=0
      JAT(MNS)=0
      JTT(MNS)=0
      JLT(MNS)=0
      MXP=MXP-1
      IF(MXP.LE.0) GO TO 300
      DC 280 M=1,MXP
      M1=MP+M
280 JAT(M1)=JAT(M1-1)+JLT(M1-1)
C====FINISHED
300 RETURN
      END

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      SUBROUTINE FPDS(JA,NP,XP,YP,IP)
C====FETCH POINT FROM /DENS/ STORAGE=====
10
C      JA - STARTING INDEX IN /DENS/ ARRAY A OF THE TAB1 RECORD
30
C      CONTAINING ARRAYS X(N),Y(N),N=1,N2
40
C      NP - DESIRED POINT NUMBER
50
C      XP - X VALUE CORRESPONDING TO NP
60
C      YP - Y VALUE CORRESPONDING TO NP
70
C      IP - INTERPOLATION CODE USED BETWEEN X(NP) AND X(NP+1)
80
C      THE FOLLOWING TABLE SUMMARIZES THE OUTPUT VALUES OF XP,YP,IP
90
C      NP      XP      YP      IP
100
C      -----
110
C      NP.LE.0      -1.0E+20      0.0      1
120
C      NP.GT.0.AND.LT.N2      X(NP)      Y(NP)      I
130
C      NP.EQ.N2      X(N2)      Y(N2)      2
140
C      NP.EQ.N2+1      X(N2)      0.0      2
150
C      NP.GT.N2+1      1.0E+20      0.0      2
160
C-----NO ERROR STOPS
170
C-----DIMENSION LA(1),A(1)
180
C-----INITIALIZE
190
10 N1=LA(JA+7)
200
      N2=LA(JA+8)
210
      J1=JA+9
220
      J2=J1+2*N1
230
C-----TEST NP
240
20 IF(NP.LE.0)GOTO30
250
      IF(NP.LE.N2)GOTO40
260
      IF(NP.EQ.N2+1)GOTO70
270
      GOTO80
280
C-----NP.LE.0
290
30 XP=-1.0E+20
300
      YP=0.0
310
      IP=1
320
      GOTO100
330
C-----NP.GT.0 AND NP.LE.N2
340
40 K=J2+2*(NP-1)
350
      XP=A(K)
360
      YP=A(K+1)
370
      DO50M=1,N1
380
      K=J1+2*(M-1)
390
      IF(NP+1.LE.LA(K))GOTO60
400
50 CCNTINUE
410
      IP=2
420
      GOTO100
430
60 IP=LA(K+1)
440
      GOTO100
450
C-----NP.EQ.N2+1
460
70 K=J2+2*(N2-1)
470
      XP=A(K)
480
      YP=0.0
490
      IP=2
500
      GCT0100
510
C-----NP.GT.N2+1
520
80 XP=1.0E+20
530
      YP=0.0
540
      IP=2
550

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C-----FINISHED
100 RETURN
ENTRY IFPDS(A,LA)
GOTO 100
END
560
570
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600

SUBROUTINE IPDS(JA,NP,XP,YP,IP)
C===== INTERPOLATE POINT IN /DENS/ STORAGE=====
C JA - STARTING INDEX IN /DENS/ ARRAY A OF THE TAB1 RECORD
C XP - GIVEN VALUE OF X
C YP - CORRESPONDING(DESIRE) VALUE OF Y
C IP - INTERPOLATION CODE USED TO COMPUTE YP
C NP - INDEX OF X AND Y ARRAYS SUCH THAT XP LIES
C BETWEEN X(NP) AND X(NP+1). IF NP.GT.0 ON INPUT, IT
C IS USED TO START SEARCH.
C JA, XP, AND NP(GUESSED VALUE OR 0) ARE GIVEN AS INPUT
C YP, IP, AND NP(CORRECT VALUE) ARE OUTPUT FROM S.F.
C IF XP.LT.X(1), YP=0.0, NP=C, IP=1
C IF XP.GT.X(N2), YP=0.0, NP=N2+1, IP=1, WHERE N2 IS THE
C NUMBER OF X,Y VALUES GIVEN IN THE TAB1 RECORD.
C-----ERROR STOP 314, IMPROPER INTERPOLATION TABLE
DIMENSION LA(1),A(1)
C-----INITIALIZE
10 N1=LA(JA+7)
N2=LA(JA+8)
J1=JA+9
J2=J1+2*N1
C-----TEST FOR XP OUT OF RANGE
20 IF(XP.GE.A(J2))GOTO30
YP=0.0
NF=0
IP=1
GOTO100
30 K=J2+2*(N2-1)
IF(XP.LE.A(K))GOTO40
YP=0.0
NP=N2+1
IP=1
GOTO100
C-----SEARCH TABLE FOR PROPER PANEL
40 IF(NP.LE.0)NP=1
IF(NP.GE.N2)NP=N2-1
K=J2+2*(NP-1)
IF(XP.LT.A(K))NP=1
N2P=N2-1
DO50N=NP,N2P
NPP=N+1
K=J2+2*N
IF(XP.LT.A(K))GOTO60
50 CCNTINUE
YP=A(K+1)
IP=2
460

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NP=N2+1
GOTO100
C-----FIND INTERPOLATION CODE
60 DO70M=1,N1
MP=M
K=J1+2*(M-1)
IF(NPP.LE.LA(K))GOTO80
70 CCNTINUE
CALL ERROR(314)
80 IP=LA(K+1)
C-----INTERPOLATE
90 K=J2+2*(NPP-2)
CALL TERP1(A(K),A(K+1),A(K+2),A(K+3),XP,YP,IP)
NP=NPP-1
C-----FINISHED
100 RETURN
ENTRY IIPDS(A,LA)
GOTO 100
END
470
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490
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520
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540
550
560
570
580
590
600
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620
630
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SUBROUTINE LRIDS(MA,JA,LNT)
C===== LOCATE RECORD IN DENSE STORAGE=====
C MA - RECORD INDENT
C JA - STARTING INDEX IN ARRAY A FOR RECORD MA
C LNT =0, NORMAL RETURN
C =1, RECORD MA NOT IN DENSE STORAGE
C-----NO ERROR STOPS
COMMON/DENS/JMT(100),JAT(100),JTT(100),JLT(100),
1 JNS,MNS,MX
10 LNT=0
MXP=MNS-1
IF(MXP.LE.0)GOTO30
IF(MA.EQ.0)GOTO30
DO20M=1,MXP
JA=JAT(M)
IF(JMT(M).EQ.MA)GOTO40
20 CONTINUE
30 LNT=1
40 RETURN
END
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100
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120
130
140
150
160
170
180
190
200

SUBROUTINE SICKY(NIS,NISO,IF4,IF5,BK,TAB,ITAB)
COMMON/TAB/ENDTAB(25),KEDTAB(50),KEDR(240),NCNAME(39),NRIF3,NRIF4
DIMENSION KDR1(123),KDR2(1)
EQUIVALENCE (KEDR(1),KDR1(1)),(KEDR(124),KDR2(1))
INTEGER ENDTAB
COMMON/INDUT/NRE1(80),NAMKDK(2),MATKDK,NTYP,NAMISO(20),MATISO(10),
*NRE2(14),NKED,NRE3(3),ITD1M
NAMELIST/LIB/IFIRST,IMOD,IDAT
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REAL*8 MTNAM,MATTAB(20),DATE,TIME
EQUIVALENCE(MTNAM,NAMKDK(1))
DIMENSION IF1(100),IF2(100),IF3(100),KEDA(3),IN(20,2),
*IT1(100),IT2(100),IT3(100),IT4(100),
*NRI(20),NTAB(7),BK(1),TAB(1),ITAB(1)
EQUIVALENCE(NINP,NI),(NCUTP,NO)
INTEGER*2 S,LSED
EQUIVALENCE(LSED,NAME1)
DATA S/'SE'/
DATA KEDA/'KEDA','BIBL','IOTH'/
DATA NRI/ 0, 4,
* 0, 3, 0, 3,
1 1, 1, 0, 4,
2 1, 3, 3, 8,
3 2, 6, 0, 3,
4 3, 8/
DIMENSION ISATZ(880),XSATZ(880)
EQUIVALENCE(ISATZ(1),XSATZ(1))
201 FORMAT('1 INSERTION OF TRANSLATED DATA INTO KEC*K/'2X,39('**')//')
NINP=5
NCUTP=6
IFILE=1
IFIL=11
L=1
NREAC2=69
NK=41
IF(NISO.EQ.0)NIS=1
CALL RDA1(BK)
DC 200 I=1,880
200 XSATZ(I)=0.0
CALL DATUM(DATE,TIME)
CALL SPLIT(DATE,IDAT1,IDAT2,IDAT3)
IDAT=IDAT1*10000+IDAT2*100+IDAT3
WRITE(NOUTP,201)
CALL PRIZE1
C-----IFIRST=0 LIBRARY STILL INITIALIZED
C 1 LIBRARY TO BE INITIALIZED
C-----IMOD=0 ISOTOPE NOT YET PRESENT IN LIBRARY
C 1 ISOTOPE YET PRESENT IN LIBRARY AND TO BE REPLACED
C
C IDAT=DATE DE COMPILATION
C
IFIRST=0
IMOD=0
READ(NINP,LIB)
WRITE(NOUTP,LIB)
MTK=MATKDK
NMTTAB=0
IF(NIS.GT.1)NIS=NIS+1
DO 10 MIS=1,NIS
ICVFL=0
IF(NIS.EQ.1)GOTO 11
IF(MIS.EQ.1)GOTO 11
IFIRST=0
NAMKDK(1)=NAMISO((MIS-1)*2-1)

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9C
100
110
120
130
14C
150
160
170
180
19C
200
210
220
230
24C
250
260
27C
280
29C
300
310
32C
330
340
350
36C
37C
380
39C
400
41C
420
430
44C
450
46C
470
480
490
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51C
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530
54C
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56C
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580
59C
600
610
620
630
NAMKDK(2)=NAMISO((MIS-1)*2)
64C
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660
670
680
69C
700
710
720
730
74C
750
760
770
780
79C
80C
810
82C
830
840
85C
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87C
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92C
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101C
1020
103C
104C
1050
106C
1070
1080
1090
1100
111C
1120
1130
114C
1150
116C
1170
1180
CALL RDA3(MTK)
NTP3=NTYP
IF(MIS.EQ.1)NTYP=NTP3+NRIF4
IF(IFIRST)210,210,220
220 DO 251 I=1,3
ISATZ(I)=KEDA(I)
251 CCNTINUE
NISOT=0
NISOT2=0
NXTR=2
GOTO 260
210 IF(IMOD.EQ.0)GOTO 327
CALL LISA(NAMKDK(1),NAMKDK(2),KFIRST)
IF(KFIRST.EQ.0)GOTO 327
IFIRST=1
GOTO 220
327 READ(IFILE'L')(ISATZ(I),I=1,880)
NISOT=ISATZ(5)
NISOT2=ISATZ(11)
IF(NISOT.EQ.0.OR.NISOT2.EQ.0)GOTO 220
IF(NISOT.EQ.NISOT2) GOTO 326
WRITE(NOUTP,1001) NISOT,NISOT2
1001 FORMAT(' WARNING: NISOT=',I3,' IS NOT EQUAL NISOT2=',I3,'. THIS MA
1Y PRODUCE ERRORS. ')
326 NXTR=ISATZ(880)
IF(NXTR.GT.1.AND.NXTR.LT.10000) GOTO 329
WRITE(NOUTP,1002)
1002 FORMAT(' LIBRARY UNCHANGED, BECAUSE WORD 880 IN RECCRD 1 IS WRONG.
1 ')
GOTO 328
329 K=ISATZ(7)
IFEND=3*NISOT
C
C-----FIRST TABLE
C
DC 290 J=1,NISOT
IF1(J)=ISATZ(K)
IF2(J)=ISATZ(K+1)
IF3(J)=ISATZ(K+2)
K=K+3
290 CCNTINUE
KC=ISATZ(13)
JENC=4*NISOT2
C
C-----THIRD TABLE
C
DO 300 J=1,NISOT2
IT1(J)=ISATZ(KC)
IT2(J)=ISATZ(KC+1)
IT3(J)=ISATZ(KC+2)
IT4(J)=ISATZ(KC+3)
KC=KC+4
300 CCNTINUE

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260	WRITE(NOUTP,270)NISOT	1190	ISATZ(KC+2)=IT3(J)	1740	
270	FORMAT(1H,'NUMBER OF ISOTOPES STILL INCLUDED INTO THE LIBRARY=',	1200	ISATZ(KC+3)=IT4(J)	1750	
	X I6)	1210	KC=KC+4	1760	
	N=1	1220	IF(INW.EQ.0.AND.J.EQ.NISOT2)GOTO 382	1770	
	ISATZ(4)=IDAT	1230	380	CCNTINUE	1780
	ISATZ(5)=NISOT+N	1240	GOTO 360	1790	
	ISATZ(6)=1	1250	330	K=14	1800
	ISATZ(7)=14	1260	KC=ISATZ(13)	1810	
	ISATZ(8)=NREAC2	1270	ISATZ(K)=NAMKDK(1)	1820	
	ISATZ(9)=1	1280	ISATZ(K+1)=NAMKDK(2)	1830	
	ISATZ(10)=3*ISATZ(5)+14	1290	ISATZ(K+2)=MTK	1840	
	ISATZ(11)=NISOT2+N	1300	K=K+3	1850	
	ISATZ(12)=1	1310	ISATZ(KC)=MTK	1860	
	ISATZ(13)=ISATZ(10)+NREAC2*3	1320	ISATZ(KC+1)=NTYP	1870	
	IF(IFIRST)320,320,330	1330	ISATZ(KC+2)=2	1880	
320	KC=KC+3*N	1340	ISATZ(KC+3)=1	1890	
	K1=14	1350	KC=KC+4	1900	
	INW=0	1360	360	K2=ISATZ(10)	1910
	DO 370 J=1,NISOT	1370	J2=NREAC2*3	1920	
	IF(INW.EQ.1)GOTO 371	1380	DO 390 J=1,J2	1930	
	IF(IF3(J).LT.MTK)GOTO 371	1390	ISATZ(K2)=KFDR(J)	1940	
	IF(IF3(J).NE.MTK)GOTO 372	1400	K2=K2+1	1950	
	IF(IMOD.EQ.1)GOTO 372	1410	390	CCNTINUE	1960
	WRITE(NO,325)NAMKDK(1),NAMKDK(2),MTK	1420	C	ADRESSE DE LA FIN DE LA TRCISIEME TABLE	1970
325	FORMAT(1H0,'THE ISOTOPE',2X,2A4,2X,I6,' IS YET IN THE LIBRARY.',	1430		KC=ISATZ(13)+4*ISATZ(11)-1	1980
	*' IMOD IS NOT SET TO 1. LIBRARY UNCHANGED')	1440	C	IMPRESSION DES TABLES	1990
	RETURN	1450		WRITE(NOUTP,400)(ISATZ(I),I=1,3)	2000
372	ISATZ(K1)=NAMKDK(1)	1460	400	FORMAT(1X,3A4)	2010
	ISATZ(K1+1)=NAMKDK(2)	1470		WRITE(NOUTP,410)(ISATZ(I),I=4,13)	2020
	ISATZ(K1+2)=MTK	1480	410	FORMAT(1X,10I6)	2030
	K1=K1+3	1490		ITAB1=ISATZ(7)	2040
	J1=J	1500		ITAB2=ISATZ(10)-1	2050
	INW=1	1510		DO 430 K=1,2	2060
	IF(J.GT.NISOT)GOTO 370	1520		WRITE(NOUTP,420)(ISATZ(I),I=ITAB1,ITAB2)	2070
371	ISATZ(K1)=IF1(J)	1530	420	FORMAT(1X,2A4,6X,I6)	2080
	ISATZ(K1+1)=IF2(J)	1540		ITAB1=ISATZ(10)	2090
	ISATZ(K1+2)=IF3(J)	1550		ITAB2=ISATZ(13)-1	2100
	K1=K1+3	1560		WRITE(NO,412)	2110
	IF(INW.EQ.0.AND.J.EQ.NISOT)GOTO 373	1570	412	FORMAT(1H0)	2120
	GOTO 370	1580	430	CONTINUE	2130
373	J=J+1	1590		ITAB1=ISATZ(13)	2140
	GOTO 372	1600		WRITE(NOUTP,411)(ISATZ(I),I=ITAB1,KC)	2150
370	CONTINUE	1610	411	FORMAT(4I10)	2160
	KC=ISATZ(13)	1620		WRITE(NO,412)	2170
	INW=0	1630		NT=0	2180
	DO 380 J=1,NISOT2	1640		NTAB(6)=NXTR	2190
	IF(J1.NE.J)GOTO 381	1650		NTAB(7)=1+7*NTYP	2200
382	ISATZ(KC)=MTK	1660	C	INDEX D'ECRIURE DANS ITAB	2210
	ISATZ(KC+1)=NTYP	1670		JINIT=1	2220
	ISATZ(KC+2)=NXTR	1680		NTD=NTAB(7)	2230
	ISATZ(KC+3)=1	1690		K=1	2240
	KC=KC+4	1700		IF(IF4.EQ.1.OR.IF5.EQ.1)GOTO 805	2250
	INW=1	1710	C		2260
381	ISATZ(KC)=IT1(J)	1720	C	DATA FROM FILES 1,2,3	2270
	ISATZ(KC+1)=IT2(J)	1730	C		2280

DC 530 I=1,NK	2290	IF(NE=880)650,650,660	2840
NAME1=KEDR(K)	2300	650 NTAB(7)=NE	2850
NAME2=KEDR(K+1)	2310	GOTO 670	2860
CALL RDA2(MTK ,MIS,NAME1,NAME2,LFT)	2320	660 NTAB(6)=NTAB(6)+(NE-1)/880	2870
IF(LFT.EQ.1)GOTO 531	2330	NTAB(7)=MOD(NE-1,880)+1	2880
NT=NT+1	2340	670 JINIT=JINIT+7	2890
NTDS=NTD+NKED-1	2350	NTC=NTDS+1	2900
NTAB(1)=KEDR(K+2)	2360	531 K=K+3	2910
IF(NTAB(1).LE.21550) GCTO 570	2370	530 CONTINUE	2920
NTAB(3)=1	2380	IF(IOVFL.EQ.0)GOTO 914	2930
NTAB(4)=1	2390	IRECR=NXTR+IOVFL*ITDIM/880	2940
GOTO 580	2400	IF(IOVFL.GT.1)IRECR=IRECR-(IOVFL-1)	2950
570 KTI=(K+2)*2/3	2410	WRITE(IFILE'IRECR)(TAB(I),I=881,IREEL)	2960
NTAB(3)=NR1(KTI-1)	2420	WRITE(IFILE'NXTR)(TAB(I),I=1,880)	2970
NTAB(4)=NR1(KTI)	2430	GOTO 805	2980
580 IF(NTAB(1).EQ.30050) ) GCTO 590	2440	914 WRITE(IFILE'NXTR)(TAB(I),I=1,NTDS)	2990
NTAB(2)=0	2450	C	3000
GOTO 610	2460	C DATA FROM FILE 4	3010
590 NTAB(2)=1	2470	C	3020
NTAB(5)=NRIF3	2480	805 IF(NRIF4.EQ.0)GOTO 902	3030
GOTO 830	2490	IF(MIS.GT.1)GOTO 901	3040
610 NTAB(5)=NKED/(NTAB(3)+NTAB(4))	2500	IF(IF4.EQ.1.OR.IF5.EQ.1)GOTO 806	3050
830 JP=1	2510	LASTW=MOD(NTDS-1,880)+1	3060
DO 760 JK1=NTD,NTDS	2520	LASTR=NXTR+(NTDS-1)/880	3070
IF(IOVFL.EQ.0)GOTO 913	2530	NTD=LASTW+1	3080
IF(JK1-(ITDIM*(IOVFL+1)-IOVFL*880))910,910,911	2540	IF(LASTR.EQ.NXTR) GOTO 807	3090
913 IF(JK1-ITDIM)910,910,911	2550	READ(IFILE'LASTR)(TAB(I),I=881,1760)	3100
911 IOVFL=IOVFL+1	2560	NTD=LASTW+891	3110
IECR=1	2570	GOTO 807	3120
IF(IOVFL.GT.1)IECR=881	2580	806 LASTW=7*NTP	3130
IRECR=NXTR+(IOVFL-1)*ITDIM/880	2590	LASTR=NXTR	3140
IF(IOVFL.GT.1)IRECR=IRECR-(IOVFL-2)	2600	JINIT=1	3150
WRITE(IFILE'IRECR)(TAB(IJKL),IJKL=IECR,ITDIM)	2610	NTD=1+7*NTP	3160
910 IREEL=JK1-IOVFL*(ITDIM-880)	2620	807 NTAB(6)=LASTR	3170
TAB(IREEL)=BK(JP)	2630	NTAB(7)=LASTW+1	3180
JP=JP+1	2640	IF(NTAB(7).LE.880)GOTO 812	3190
760 CONTINUE	2650	NTAB(7)=1	3200
IF(NTAB(1).NE.30050)GOTO 841	2660	NTAB(6)=NTAB(6)+1	3210
IS=NTD	2670	812 K=1	3220
DO 821 IRL=1,NRIF3	2680	DC 801 IK=1,39	3230
ITAB(IS+2)=ITAB(IS+2)+NTAB(6)-1	2690	IF(NCNAME(IK).EQ.0)GOTO 817	3240
ITAB(IS+3)=ITAB(IS+3)+NTAB(7)-1	2700	NTAB(5)=NCNAME(IK)	3250
IF(ITAB(IS+3).LE.880)GOTO 822	2710	NAME1=KDR2(K)	3260
ITAB(IS+2)=ITAB(IS+2)+1	2720	NAME2=KDR2(K+1)	3270
ITAB(IS+3)=ITAB(IS+3)-880	2730	NT=NT+1	3280
822 IS=IS+4	2740	CALL RDA4(MTK,1,NAME1,NAME2,LNT,IR)	3290
821 CONTINUE	2750	NTDS=NTD+NKED-1	3300
841 JLS=JINIT+6	2760	IF(NTDS.GE.ITCIM)GOTO 641	3310
KP=1	2770	READ(IFIL'IR)(TAB(I),I=NTD,NTDS)	3320
DC 640 J=JINIT,JLS	2780	C-----SET NUMERICAL NAME	3330
ITAB(J)=NTAB(KP)	2790	NTAB(1)=KDR2(K+2)	3340
KP=KP+1	2800	C-----SET NUMBER OF ARGUMENTS	3350
IF(J-ITDIM)640,641,641	2810	NTAB(3)=1	3360
640 CONTINUE	2820	C-----SET NUMBER OF FUNCTIONAL VALUES	3370
NE=NKED+NTAB(7)	2830	NTAB(4)=1	3380

```

      IF(LSEQ.EQ.S.AND.NTAB(1).GE.50000)NTAB(4)=3
C-----SET NUMBER OF ADD. NAMES
      IF(NTAB(1).GE.40051)GOTO 808
      NTAB(2)=1
      GOTO 809
808 IF(NTAB(1).GE.44651)GOTO 810
      NTAB(2)=2
      GOTO 809
810 IF(NTAB(1).GE.50000)GOTO 811
      NTAB(2)=3
      GOTO 809
811 NTAB(2)=1
809 JLS=JINIT+6
      KP=1
      DO 813 J=JINIT,JLS
      ITAB(J)=NTAB(KP)
      KP=KP+1
      IF(J-ITDIM)813,641,641
813 CCNTINUE
C RELOCATION OF THE REACTION TYPE TABLE
      NTRL=NCNAME(IK)
      IS=NTD
      IS=IS+NTAB(2)-1
      DC 818 IRL=1,NIRL
      ITAB(IS+2)=ITAB(IS+2)+NTAB(6)-1
      ITAB(IS+3)=ITAB(IS+3)+NTAB(7)-1
      IF(ITAB(IS+3).LE.880)GOTO 820
      ITAB(IS+2)=ITAB(IS+2)+1
      ITAB(IS+3)=ITAB(IS+3)-880
820 IS=IS+NTAB(2)+3
818 CCNTINUE
      IS=IS-1
C
      NKED=NKED+NTAB(7)
      IF(NKED-880)814,814,815
814 NTAB(7)=NKED
      GOTO 816
815 NTAB(6)=NTAB(6)+(NKED-1)/880
      NTAB(7)=MOD(NKED-1,880)+1
816 JINIT=JINIT+7
      NTC=NTDS+1
817 K=K+3
801 CCNTINUE
901 JINIT=JINIT-1
      WRITE(NO,900)(ITAB(I),I=1,JINIT)
900 FORMAT(7I12)
      IF(NT.NE.NTYP)WRITE(NO,740)NT,NTYP
740 FORMAT(1H1,'NT=',I6,' NTYP=',I6//)
      JP=JINIT+1
C
C WRITING OF DATA IN KEDAK
C
      IF(MIS.GT.1)GOTO 902
      IF(IF4.EQ.1.OR.IF5.EQ.1.OR.LASTR.EQ.NXTR) GOTO 819
      WRITE(IFILE'NXTR')(TAB(I),I=1,880)

```

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3390
3400
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3880
3890
3900
3910
3920
3930

```

```

      WRITE(IFILE'LASTR')(TAB(I),I=881,NTDS)
      NXTR=(NTDS-881)/880+1+LASTR
      GOTO 900
819 WRITE(IFILE'NXTR')(TAB(I),I=1,NTDS)
902 NXTR=(NTDS-1)/880+1+NXTR
900 ISATZ(880)=NXTR
      WRITE(IFILE'I')(ISATZ(I),I=1,880)
      NMTTAB=NMTTAB+1
      MATTAB(NMTTAB)=MTNAM
      WRITE(NO,750)NXTR
750 FORMAT(1H1,'NEXT AVAILABLE RECORD IN THE KEDAK LIBRARY',I6//)
      10 CCNTINUE
328 CALL TABPRT(MATTAB,NMTTAB,IFILE)
      RETURN
641 WRITE(NO,642)
642 FORMAT(1H1,'OVERFLOW IN TAB VECTOR')
      RETURN
      END
3940
3950
3960
3970
3980
3990
4000
4010
4020
4030
4040
4050
4060
4070
4080
4090
4100
4110
C
      SUBROUTINE LISAI(INAM1,INAM2,IFIRST)
      J.CANIELS AUGUST 1971
C NTEL(1)=LENGTH OF MATERIAL CONVERSION TABLE/NTEL(2-3)=ADDRESS
C NTEL(4)=LENGTH OF TYPE CONVERSION TABLE/NTEL(5-6)=ADDRESS
C NTEL(7)=LENGTH OF MATERIAL ADDRESS TABLE / NTEL(8-9)=ADDRESS
C ISATZ(880)=TOTREC=FIRST AVAILABLE RECCRD
      INTFGER OLDRREC,TOTREC,TEMP
      DIMENSION ISATZ(880),NRECO(62),NREAC(62),NTEL(9),NTEST(62)
      DIMENSION IS2(880)
      NC=6
      NT=1
      NFWREC=1
      OLDRREC=1
      ILAST=0
      IFIRST=0
      DC 303 J=1,62
      303 NTEST(J)=0
C READ FIRST RECORD *****
      READ(NT'OLDREC')(ISATZ(I),I=1,880)
      DC 301 J=1,9
      301 NTEL(J)=ISATZ(J+4)
      TOTREC=ISATZ(880)
      M=NTEL(1)
      N=NTEL(3)
C TEST ON FIRST TABLE *****
      DC 2 J=1,M
      IF(INAM1.EQ.ISATZ(N+3*J-3).AND.INAM2.EQ.ISATZ(N+3*J-2))GOTO 6
      2 CCNTINUE
      WRITE(NO,10)INAM1,INAM2
      10 FORMAT(' THIS NAME ',2A4,' NOT FOUND IN TABLE 1 OF THIS LIBRARY
      1'/' LIBRARY UNCHANGED')
      RETURN
      6 NUM=ISATZ(N+3*J-1)

```



```

INDEX1=N+3*J-3
INDEX3=N+3*J
M=NTEL(7)
N=NTEL(9)
DC 300 J=1,M
NRECO(J)=ISATZ(N+4*J-2)
300 NREAC(J)=ISATZ(N+4*J-3)
C SECOND TABLE UNCHANGED *****
C TEST ON THIRD TABLE *****
DO 3 J=1,M
IF(NUM.EQ.ISATZ(N+4*J-4))GO TO 8
3 CONTINUE
WRITE(NO,11)NUM
11 FORMAT(' THE NUMBER CORRESPONDING THE NAME OF THE ASKED ELEMENT ',
110,' IS NOT FOUND IN TABLE 3'/' LIBRARY IS UNCHANGED')
RETURN
8 INDEX2=N+4*J-4
INDEX4=N+4*J
KN=J
IF(M.GT.1)GOTO 200
IFIRST=1
RETURN
200 IF(KN.EQ.M)ILAST=1
NREC=NRECO(KN)
C CHANGING FIRST RECORD(BEGIN + TABLE 1 + TABLE 2 TILL THE WRIGHT
C PLACE IN TABLE 3 ) *****
ISATZ(5)=ISATZ(5)-1
ISATZ(10)=ISATZ(10)-3
ISATZ(11)=ISATZ(11)-1
ISATZ(13)=ISATZ(13)-3
NTOT=INDEX2-4
DO 20 J=INDEX1,NTOT
20 ISATZ(J)=ISATZ(J+3)
IF(ILAST.EQ.1)GOTO 101
NINIT=NTOT+1
NFIN=ISATZ(13)+4*ISATZ(11)-1
DC 100 J=NINIT,NFIN
100 ISATZ(J)=ISATZ(J+7)
C-----ARRANGEMENT DES NUMEROS DE RECORDS EN CRDRE CROISSANT
101 M1=M-1
DO 102 J=1,M1
J1=J+1
DC 102 J1=J1,M
IF(NRECO(J).LT.NRECO(J1))GOTO 102
TEMP=NRECO(J)
NRECO(J)=NRECO(J1)
NRECO(J1)=TEMP
TEMP=NREAC(J)
NREAC(J)=NREAC(J1)
NREAC(J1)=TEMP
102 CONTINUE
WRITE(NO,1500)(NRECO(J),NREAC(J),J=1,M)
1500 FORMAT(1X,2I6)
C-----L'ISOTOPE A S-T-IL LA PLUS HAUTE ADRESSE
IF(NRECO(M).GT.NREC)GOTO 103

```

```

34C
350
36C
370
380
39C
400
41C
420
43C
44C
450
460
470
48C
49C
500
510
520
530
54C
550
560
57C
580
590
600
610
62C
630
64C
650
660
67C
680
690
70C
71C
720
73C
740
750
760
770
78C
790
80C
810
820
83C
840
850
86C
87C
880

```

```

ISATZ(880)=NREC
WRITE(NT'1)(ISATZ(I),I=1,880)
RETURN
103 DC 104 J=1,M
IF(NRECO(J).EQ.NREC)GOTO 105
104 CCNT INUE
STOP
105 KN=J+1
NDIFF=NRECO(KN)-NREC
C-----MODIFICATION DE LA 3E TABLE DU RECCRD 1
DC 108 J=1,M1
KLM=ISATZ(13)+J*4-2
IF(ISATZ(KLM).LE.NREC)GOTO 108
ISATZ(KLM)=ISATZ(KLM)-NDIFF
108 CONTINUE
ISATZ(880)=ISATZ(880)-NDIFF
WRITE(NT'1)(ISATZ(I),I=1,880)
C-----DEPLACEMENT DES ISOTOPE DONT L'ADRESSE EST SUP. A NREC
DO 107 J=KN,M
N1=NRECO(J)
IF(J.LT.M)GOTO 109
N2=TOTREC-1
GOTO 110
109 N2=NRECO(J+1)-1
110 DC 107 K=N1,N2
READ(NT'K)(ISATZ(I),I=1,880)
IF(K.GT.N1)GOTO 111
JK=NREAC(J)
DO 112 JR=1,JK
J6=JR*7-1
IF(ISATZ(JR*7-5).EQ.0)GOTO 120
NFNAM=ISATZ(JR*7-5)
NRFN=ISATZ(JR*7-1)
NWFN=ISATZ(JR*7)
NLONG=ISATZ(JR*7-2)
READ(NT'NRFN)(IS2(I),I=1,880)
NRFNN=NRFN
DC 121 NLO=1,NLONG
INLO=NWFN+NLO*(3+NFNAM)-2
LESDIF=(INLO-1)/880
INLOLO=INLO-LESDIF*880
IF((NRFN+LESDIF).EQ.NRFNN)GOTO 415
WRITE(NT'NRFNN)(IS2(I),I=1,880)
NRFNN=NRFN+LESDIF
READ(NT'NRFNN)(IS2(I),I=1,880)
415 IS2(INLOLO)=IS2(INLOLO)-NDIFF
IF(K.EQ.NRFNN.AND.INLO.LE.880)ISATZ(INLO)=ISATZ(INLO)-NDIFF
121 CONTINUE
WRITE(NT'NRFNN)(IS2(I),I=1,880)
120 ISATZ(J6)=ISATZ(J6)-NDIFF
112 CONTINUE
111 KL=K-NDIFF
WRITE(NT'KL)(ISATZ(I),I=1,880)
WRITE(NO,1051)K,KL
1051 FORMAT(1X,'K=',I6,' KL=',I6)

```

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890
900
910
920
930
940
95C
960
970
980
99C
1000
101C
1020
1030
104C
1050
1060
1070
1080
109C
1100
1110
1120
1130
114C
1150
116C
1170
1180
119C
1200
121C
1220
1230
124C
1250
126C
1270
128C
129C
1300
131C
1320
133C
1340
1350
136C
1370
1380
1390
1400
141C
1420
1430

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```

IF(ATYCON(1).EQ.LX) GOTC 110
LX=ATYCON(1)
READ(IFIL'LX) RW
110 LS=ATYCON(2)
LE=LTYCON*3+LS-1
CALL PAGE(TABLES,IL)
WRITE(KOUT,606)
606 FCRMAT(/5X,10('*'),2X,'TYPECONVERSIONTABLE:',2X,10('*'))//11X,
1'ALPHA',T21,'NUMERIC')
J=1
K=1
M=LS-1
DC 120 I=LS,LE,3
M=M+1
IF(M.LE.LRECD) GOTO 112
LX=LX+1
READ(IFIL'LX) RW
M=1
112 MS=M
LXS=LX
WTYP(J)=W(M)
M=M+1
IF(M.LE.LRECD) GOTO 114
LX=LX+1
READ(IFIL'LX) RW
M=1
114 WTYP(J+1)=W(M)
M=M+1
IF(M.LE.LRECD) GOTO 116
LX=LX+1
READ(IFIL'LX) RW
M=1
116 ITYP(K)=IW(M)
IF(IL.GT.ILM) CALL PAGE(TABLES,IL)
IL=IL+1
WRITE(KOUT,616) LXS,MS,WTYP(J),WTYP(J+1),ITYP(K)
616 FCRMAT(1X,'(',I3,',',I3,')',1X,2A4,T20,I7)
K=K+1
J=J+2
120 CONTINUE
C
C     FETCH MATERIAL ADRESSTABLE.
C
IF(AMATAB(1).EQ.LX) GOTO 210
LX=AMATAB(1)
READ(IFIL'LX)RW
210 LS=AMATAB(2)
LE=LMATAB*4+LS-1
J=0
CALL PAGE(TABLES,IL)
WRITE(KOUT,617)
617 FCRMAT(/5X,10('*'),2X,'MATERIALADRESSTABLE:',2X,10('*'))//12X,
1'NAME',T30,'LENGTH',T40
1,'ADDRESS',T20,'ALPHA')
M=LS-1

```

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104C
1050
106C
107C
1080
109C
1100
1110
1120
1130
114C
1150
1160
117C
1180
1190
1200
1210
122C
1230
1240
1250
1260
127C
1280
1290
130C
1310
132C
1330
1340
135C
1360
137C
1380
1390
140C
1410
142C
1430
1440
1450
1460
147C
1480
1490
150C
1510
1520
1530
1540
1550
1560
157C
1580

```

```

DC 220 I=LS,LE,4
J=J+1
M=M+1
IF(M.LE.LRECD) GCTO 208
LX=LX+1
READ(IFIL'LX) RW
M=1
208 MS=M
LXS=LX
ISNAM(J)=IW(M)
M=M+1
IF(M.LE.LRECD) GOTO 211
LX=LX+1
READ(IFIL'LX) RW
M=1
211 LISTYP(J)=IW(M)
M=M+1
IF(M.LE.LRECD) GOTO 212
LX=LX+1
READ(IFIL'LX) RW
M=1
212 AISTYP(1,J)=IW(M)
M=M+1
IF(M.LE.LRECD) GOTO 214
READ(IFIL'LX) RW
M=1
LX=LX+1
214 AISTYP(2,J)=IW(M)
DC 215 L=1,LMACON
IF(IMAT(L).NE.ISNAM(J)) GOTO 215
ALPHAN=MATNAM(L)
GOTO 216
215 CCNTINUE
216 IF(IL.GT.ILM) CALL PAGE(TABLES,IL)
IL=IL+1
WRITE(KOUT,607) LXS,MS,ISNAM(J),ALPHAN,LISTYP(J),(AISTYP(K,J),K=1,
12)
607 FCRMAT(1X,'(',I3,',',I3,')',1X,I7,T20,A8,T30,I5,T38,'(',I3,',',I3,
1')')
220 CONTINUE
C
C     FOR EACH MATERIAL PRINT ADRESSTABLE OF TYPES.
C
DC 1000 L=1,LMATAB
DO 225 I=1,LMACON
IF(IMAT(I).NE.ISNAM(L)) GOTO 225
ALPHAN=MATNAM(I)
GCTO 226
225 CONTINUE
ALPHAN=NF
226 IF(MAT(I).EQ.ALL) GCTO 230
DO 240 I=1,NMAT
IF(MAT(I).NE.ALPHAN ) GOTO 240
MAT(I)=BLANK
GOTO 230

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159C
1600
1610
162C
1630
164C
1650
1660
167C
1680
169C
1700
1710
172C
1730
1740
1750
1760
177C
1780
1790
1800
1810
182C
1830
1840
1850
1860
1870
1880
189C
1900
1910
192C
1930
194C
1950
1960
1970
1980
199C
2000
2010
2020
2030
204C
2050
2060
2070
2080
209C
2100
2110
2120
2130

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240	CCONTINUE	2140	ADDNAM(5,KW)=XALPH(1)	2690
	GOTO 1000	2150	ACCNAM(7,KW)=XALPH(2)	2700
230	TITLE=ALPHAN	2160	310 CCONTINUE	2710
	CALL PAGE(TITLE,IL)	2170	IF(KW.EQ.0) GOTO 1000	2720
	WRITE(KOUT,618) ISNAM(L),ALPHAN	2180	IF(IL.GT.45) CALL PAGE(TITLE,IL)	2730
618	FORMAT(/5X,10('*')) ,2X,'TYPETABLE FOR:',I7,' ALPHA NAME: ',A8,	2190	WRITE(KOUT,610)	2740
	12X,10('*'))//	2200	IL=IL+4	2750
	1T15,'TYPE',T33,'ADD.NAMES',T45,'ARG.NUM',T55,'FW.NUM',T65,'DP.NUM'	2210	DC 400 K=1,KW	2760
	2,T75,'ADRESS',T21,'ALPHA'//	2220	IF(ADDNAM(4,K).EQ.LX) GOTO 406	2770
	IL=IL+6	2230	LX=ADDNAM(4,K)	2780
	IF(AISTYP(1,L).EQ.LX) GOTO 300	2240	READ(IFIL'LX) RW	2790
	LX=AISTYP(1,L)	2250	406 NW=ADDNAM(2,K)	2800
	READ(IFIL'LX) RW	2260	LS=ADDNAM(5,K)	2810
300	LS=AISTYP(2,L)	2270	LC=NW+3	2820
	LE=LISTYP(L)*7+LS-1	2280	LE=ADDNAM(3,K)*LC+LS-1	2830
	KW=0	2290	IF(IL.GT.50) CALL PAGE(TITLE,IL)	2840
	M=LS-1	2300	WRITE(KOUT,611) ADDNAM(1,K),ADDNAM(6,K),ADDNAM(7,K)	2850
	DC 310 I=LS,LE,7	2310	IL=IL+4	2860
	M=M+1	2320	M=LS-1	2870
	IF(M.LE.LRECD) GOTO 302	2330	DO 420 I=LS,LE,LC	2880
	LX=LX+1	2340	LEI=I+NW-1	2890
	READ(IFIL'LX) RW	2350	MS=M+1	2900
	M=1	2360	LXS=LX	2910
302	MS=M	2370	IF(MS.GT.LRECD) LXS=LXS+1	2920
	LXS=LX	2380	IF(MS.GT.LRECD) MS=1	2930
		2390	KJ=0	2940
C	ALPHA NAME OF THIS TYPE.	2400	DC 410 J=I,LEI	2950
C	DC 305 K=1,LYCON	2410	KJ=KJ+1	2960
	IF(ITYP(K).NE.IW(M)) GOTO 305	2420	M=M+1	2970
	ALPHAN=TYP(K)	2430	IF(M.LE.LRECD) GOTO 408	2980
	GOTO 306	2440	LX=LX+1	2990
305	CONTINUE	2450	READ(IFIL'LX) RW	3000
	ALPHAN=NF	2460	M=1	3010
306	ITW(1)=IW(M)	2470	408 TW(KJ)=W(M)	3020
	DC 308 J=2,7	2480	410 CCONTINUE	3030
	M=M+1	2490	IF(IL.GT.ILM) CALL PAGE(TITLE,IL)	3040
	IF(M.LE.LRECD) GOTO 307	2500	WRITE(KOUT,612) LXS,MS,(ITW(J),J=1,KJ)	3050
	LX=LX+1	2510	JS=I+NW	3060
	READ(IFIL'LX) RW	2520	LEI=JS+2	3070
	M=1	2530	MS=M+1	3080
307	ITW(J)=IW(M)	2540	LXS=LX	3090
308	CONTINUE	2550	IF(MS.GT.LRECD) LXS=LXS+1	3100
	IF(IL.GT.ILM) CALL PAGE(TITLE,IL)	2560	IF(MS.GT.LRECD) MS=1	3110
	WRITE(KOUT,608) LXS,MS,ITW(1),ALPHAN,(ITW(J),J=2,7)	2570	KJ=0	3120
608	FORMAT(1X,'(',I3,',',I3,')',1X,I7,2X,A8,3I10,I12,2X,'(',I3,',',I3,	2580	DC 416 J=JS,LEI	3130
	1,')')	2590	KJ=KJ+1	3140
	IL=IL+1	2600	M=M+1	3150
	NW=ITW(2)	2610	IF(M.LE.LRECD) GOTO 414	3160
	IF(NW.EQ.0) GOTO 310	2620	LX=LX+1	3170
	KW=KW+1	2630	READ(IFIL'LX) RW	3180
	ADDNAM(1,KW)=ITW(1)	2640	M=1	3190
	ADDNAM(2,KW)=NW	2650	414 ITW(KJ)=IW(M)	3200
	ACCNAM(3,KW)=ITW(5)	2660	416 CCONTINUE	3210
	ADDNAM(4,KW)=ITW(6)	2670	WRITE(KOUT,613) LXS,MS,(ITW(J),J=1,KJ)	3220
	ACCNAM(5,KW)=ITW(7)	2680	IL=IL+3	3230

```

420 CONTINUE
400 CCNTINUE
610 FCRMAT(///5X,5('*'),2X,'TABLES FOR TYPES WITH MORE THAN ONE NAME',
12X,5('*'))
611 FCRMAT(///5X,'** TYPE',I7,' ALPHANAME:',2A4,' **/')
612 FORMAT(10X,'(,I3,',',I3,')',4X,' ADD. NAMES:',5(1PE13.5))
613 FORMAT(10X,'(,I3,',',I3,')',4X,'DP.NUM,ADR.:', ' I5,2X,'(,I3,',',
1,I3,')'/)
1000 CONTINUE
IF(MAT(I).EQ.ALL) GOTO 1600
K=C
DO 1100 I=1,NMAT
IF(MAT(I).EQ.BLANK) GOTO 1100
IF(K.EQ.0) CALL PAGE(END,IL)
K=1
IF(IL.GT.ILM) CALL PAGE(END,IL)
WRITE(KOUT,622) MAT(I)
IL=IL+5
622 FORMAT(///10X,'***** ',A8,' NOT FOUND *****//')
1100 CONTINUE
IF(K.EQ.1) GOTO 2000
1600 CALL PAGE(END,IL)
WRITE(KOUT,620)
620 FORMAT(///// ' ***** NORMAL END OF TABPRT *****' )
2000 RETURN
END

```

```

3240
3250
326C
3270
328C
3290
3300
3310
3320
333C
3340
3350
336C
3370
3380
3390
3400
341C
3420
3430
3440
3450
346C
3470
348C
3490

```

```

SUBROUTINE PAGE(TITLE,IL)
REAL*8 TITLE
DATA KOUT/6/
NPG=NPG+1
WRITE(KOUT,600) IFIL,NPG,TITLE
600 FORMAT('I',9I1X,' FT-NUM-',I2,' PAGE-',I3,' **',A8,'**')
IL=3
RETURN
ENTRY INITPG(IFIL)
NPG=0
RETURN
END

```

```

10
20
30
40
50
60
70
80
90
100
11C
120

```

```

SUBROUTINE SPLIT(A,I1,I2,I3)
REAL*8 A,AA
INTEGER*2 I(4),FPZ(10),FZP(10)
EQUIVALENCE (AA,I(1))
DATA FZP/'0.', '1.', '2.', '3.', '4.', '5.', '6.', '7.', '8.', '9.'/
DATA FPZ/'0.', '1.', '2.', '3.', '4.', '5.', '6.', '7.', '8.', '9.'/
AA=A
CALL CON(I(1),I1)
CALL CON(I(4),I3)

```

```

10
2C
30
40
50
60
7C
80
90

```

```

I2E=0
DO 1 J=1,10
IF(I(3).EQ.FZP(J))GOTO 2
I2E=I2E+1
1 CONTINUE
2 CONTINUE
I2Z=0
DO 3 J=1,10
IF(I(2).EQ.FPZ(J))GOTO 4
I2Z=I2Z+1
3 CONTINUE
4 CONTINUE
I2=I2E+I2*I2Z
IF(I2E.GE.10.OR.I2Z.GE.10)WRITE(6,100)I2E,I2Z
100 FCRMAT(///' FEHLER IN SPLIT . I2E=',I4,' I2Z=',I4)
RETURN
END

```

```

100
110
12C
130
140
150
160
17C
180
19C
200
210
22C
230
240
250
260

```

```

SUBROUTINE CON(IA,IN)
INTEGER*2 IA,IN2(2),II(10)
DATA IFOR/'I2 ' /
DATA II/'00', '01', '02', '03', '04', '05', '06', '07', '08', '09' /
IM=0
IN=0
DO 1 J=1,100
IF(J.GT.10)GOTO 3
IN2(1)=II(J)
GOTO 4
3 CONTINUE
CALL CONVX(IM,IN2,IFOR)
4 CONTINUE
IF(IN2(1).EQ.IA)GOTO 2
IM=IM+1
IN=IN+1
1 CONTINUE
2 CONTINUE
IF(IN.GE.100)WRITE(6,100)IA,IN
100 FORMAT(///' FEHLER IN CCN . IA=',A4,' IN=',I4)
RETURN
END

```

```

10
20
30
40
5C
60
7C
80
9C
10C
110
12C
130
140
15C
160
170
180
190
20C
210
220

```

```

SUBROUTINE PRIEIN(IEIN, IOUT, ZEILMA)
C
C AUSDRUCK DER VON DER EINHEIT IEIN KOMMENDEN EINGABE AUF IOUT.
C DIE EINGABE-EINHEIT MUSS DIE DCB-SPEZIFIKATIONEN RECFM=F BZW.
C RECFM=FB SOWIE LRECL=80 HABEN.
C WEITERHIN SOLLTE DIE DATEI AN IHREM LOGISCHEN ANFANG STEHEN.
C DIE AUSGABEEINHEIT MUSS EINE LOGISCHE SATZLAENGE VON MINDESTENS
C 91 BYTE HABEN.

```

```

10
2C
30
4C
50
60
70
80

```

C	INTEGER ZEIL,ZEILMA	9C	BUFREG	START	1C
	DIMENSION A(20)	100	*	AUFRUF: CALL BUFREG(INTREG,&LABEL)	20
	LOGICAL*4 SECONO	110	*	INTREG ENTHAELT NACH DEM AUFRUF IN K-BYTE DIE REGION, DIE FUER	30
	SECONO=.FALSE.	12C	*	PUFFER BENOETIGT WIRD.	40
	IF(ZEILMA.LT.10.OR.ZEILMA.GT.60) ZEILMA=60	130	*	IN DER TIOT WERDEN ALLE VORHANDENEN DD-NAMEN DER FORM FTXXFOOI	50
1	WRITE(IOUT,100)	140	*	MIT AUSNAHME DES DATASETS FUER DIE DRUCKAUSGABE	60
100	FORMAT(1H1)	150	*	( DIE FILENUMMER STEHEN IM COMMON INOUT )	70
	WRITE(IOUT,101)	160	*	DURCHGEGANGEN, DIE BLKSIZE GEHOLT, DIE PUFFERANZAHL GLEICH 2	8C
101	FORMAT(' ',90('*'))	17C	*	GESETZT, DER FUER DIE PUFFER BENOETIGTE PLATZ BERECHNET UND	90
	WRITE(IOUT,102)	180	*	DIESER ZU INTREG ADDIERT.	100
102	FORMAT(' ',34('*'),' AUSDRUCK DER EINGABE '	19C	*	BEI VORKOMMEN DER DDNAMEN SYSUDUMP BZW. SYSAEENC WERDEN JE 2K	110
	IF(.NOT.SECONO) WRITE(IOUT,103)	200	*	ZU INTREG ADDIERT.	120
103	FORMAT('+',56X,34('*'))	210	*	TRITT EIN FEHLER AUF, SO WIRD RETURN 1 GEMACHT, WOBEI DIE IN	13C
	IF(SECONO) WRITE(IOUT,104)	220	*	INTREG STEHENDE NEGATIVE ZAHL DEN FEHLERGRUND ANGIBT.	140
104	FORMAT('+',56X,'( FORTSETZUNG ) ',18('*'))	230	*	2 '' '' AUSGABE '' '' ''	15C
	WRITE(IOUT,101)	24C	*	5 '' '' AUSGABE '' '' .	160
	WRITE(IOUT,105)	250	*	7 PROGRAMM- ODER MASCHINENFEHLER.	170
105	FORMAT(' ')	26C	*	9 FEHLERMELDUNG STEHT IN SYMSMG.	18C
	ZEIL=4	270	*		190
	IF(SECONO) GOTC 2	280	Q	EQU 3 HILFSSPEICHER. ( KURZZEITSPEICHER ) .	200
	N=0	29C	AUS	EQU 5 FILENUMMER VON AUSGABE.	210
	SECONO=.TRUE.	300	REGION	EQU 11 SUMME DER PUFFERREGION.	220
3	READ(IEIN,106,END=10) A	310	DDLAUF	EQU 5	23C
106	FORMAT(20A4)	320	DDINCR	EQU 6	240
	N=N+1	330	I	EQU 8	250
	2 ZEIL=ZEIL+1	340	ACHT	EQU 12	260
	IF(ZEIL.GT.ZEILMA) GOTC 1	350	J	EQU 9	270
	WRITE(IOUT,107) A,N	360	NFWOLD	EQU 4	28C
107	FORMAT(' ',20A4,' #',I5)	370	BRANCH1	EQU 14	290
	GOTO 3	380	BRANCH2	EQU 2	300
10	WRITE(IOUT,105)	39C	PBLKS	EQU 7	310
	WRITE(IOUT,101)	400	*		320
	WRITE(IOUT,102)	41C	*		33C
	WRITE(IOUT,108)	420	*		340
108	FORMAT('+',56X,' ( ENDE ) ',24('*'))	430		SAVE (14,12),,*	350
	WRITE(IOUT,101)	440		USING BUFREG,10	360
	REWIND IEIN	450		EXTRN INJUT	370
	RETURN	46C		LR 10,15	38C
	END	470	*	SAVE-AREA-VERKETTUNG.	390
		480		LR 3,13	400
				LA 13,SAVE	41C
				ST 13,8(0,3)	420
				ST 3,4(0,13)	430
				L Q,0(1) ADRESSE DER ZU UEBERGEHENDEN REGION.	440
				ST Q,PARMACR	450
				L Q,=A(INOUT) ADRESSE VON COMMON INOUT.	460
				L AUS,0(Q) FILE-NUMMER VON AUSGABE.	47C
				C AUS,=F'1'	480
				BL FEHLER2	490
				C AUS,=F'99'	50C
				BH FEHLER5	510
			*	KONVERTIERUNG DER FILENUMMER IN DEZIMALDARSTELLUNG UND	520
			*	SPEICHERN NACH FTAUS.	530
				CVD AUS,DHILF	540
				UNPK DHILF+6(2),DHILF+6(2)	550

	MVZ	DHILF+7(1),DHILF+6	560		LA	J,7	1110
	MVC	FTAUS+2(2),DHILF+6	570		B	LABWEI	1120
*		FREIE REGION VOR GETMAIN.	580	DDWRIT	EQU	*	1130
	LA	1,PARM1	590	*	DDNAME	IN DCB BRINGEN.	1140
	L	15,=V(FREESP)	600		MVC	DCBALL+40(8),DDNAM	1150
	BALR	14,15	610	*	LFSFN	VON JFCB.	1160
*		ADRESSE DER TIOT HOLEN. ADRESSE DES 1. DDNAMEN SETZEN.	620		RDJFCB	DCBALL	1170
	SR	REGION,REGION PUFFERPLATZ NULL SETZEN.	630	*	RDJFCB	ERFOLGREICH?	1180
*		MISREG NULL SETZEN.	640		LTR	15,15	1190
	ST	REGION,MISREG	650	*	BNZ	FEHLER7	1200
*		ERROR NULL SETZEN.	660	*		DATASET NEW ODER OLD? ( NEWOLD=1 DATASET IST OLD.	1210
	ST	REGION,ERROR	670	*		NEWOLD=C DATASET IST NEW. )	1220
	EXTRACT	TIOTADR,FIELDS=(TIOT)	680		LA	BRANCH1,M7	1230
	L	Q,TIOTADR	690		LA	BRANCH2,M6	1240
	LA	DDLAUF,24(Q) ADRESSE DES 1. DD-ENTRY.	700		SR	NEWOLD,NEWOLD	1250
*		PRUEFEN OB LETZTER DD-ENTRY.	710		TM	JFCB+87,X'CO'	1260
	LA	ACHT,8	720	*		REI OLD UND MOD WIRD NEWOLD=1 GESETZT. BEI CC SPRUNG NACH	1270
WEITER	SR	Q,Q	730	*		FEHLER7.	1280
	IC	Q,Q(DDLAUF)	740		BD	M1	1290
	C	Q,=F*0'	750		BZ	FEHLER7	1300
	BE	RETURN	760		LA	NEWOLD,1	1310
	LR	DDINCR,Q LAENGE DES DD-ENTRIES.	770	M1	EQU	*	1320
*		DDNAMEN HOLEN.	780	*		HOLEN VON BLKSIZE AUS JFCB.	1330
	MVC	DDNAM(8),4(DDLAUF)	790		L	Q,JFCB+100	1340
	LD	0,DDNAM	800		SLL	Q,16	1350
*		VERGLEICH MIT FTAUS.	810		SRL	Q,16	1360
	CD	0,FTAUS	820		ST	Q,BLKSZ	1370
	BE	ERHOEHE	830		LTR	Q,Q	1380
*		HAT NAMEN DIE FORM FTXXFY, SYSABEND BZW. SYSUDUMP?	840		BNP	NDBLKS	1390
	CD	0,SYSA8	850	M2	EQU	*	1400
	BE	ADD2K ES WERDEN 2K-BYTE ADDIERT.	860	*		AUF VIELFACHE VON 4 AUFRUNDEN.	1410
	CD	0,SYSD	870		LR	I,Q	1420
	BE	ADD2K ES WERDEN 2K-BYTE ADDIERT.	880		SRL	Q,2	1430
	LH	Q,DDNAM	890		SLL	Q,2	1440
	C	Q,=XL4'FFFFC6E2'	900		CR	I,Q	1450
	BNE	ERHOEHE	910		BE	M3	1460
	SR	Q,Q	920		LA	Q,4(0)	1470
	IC	Q,DDNAM+4	930	M3	EQU	*	1480
	C	Q,=XL4'000000C6'	940	*		BUFFER ALIGNMENT PRUEFEN. GEGEBENENFALLS Q ALF VIELFACHE	1490
	BNE	ERHOEHE	950	*	VON 8 AUFRUNDEN.		1500
	L	I,=F*-8'	960		TM	JFCB+92,X'01'	1510
	LA	J,2	970		BD	M4	1520
LABWEI	IC	Q,DDNAM(J)	980		LR	I,Q	1530
	C	Q,=XL4'000000F0'	990		SRL	Q,3	1540
	BL	ERHOEHE	1000		SLL	Q,3	1550
	C	Q,=XL4'000000F9'	1010		CR	I,Q	1560
	BH	ERHOEHE	1020		BE	M4	1570
	AR	I,ACHT	1030		LA	Q,8(Q)	1580
	B	LABEL(I)	1040	M4	EQU	*	1590
LABEL	LA	J,3	1050		LR	RBLKS,Q	1600
	B	LABWEI	1060		B	J(BRANCH2)	1610
	LA	J,5	1070	*		WENN DATASET NEW, DANN SPRUNG NACH GTMN. SCNST	1620
	B	LABWEI	1080	*		OLD DATASET LABEL LESEN UND PRUEFEN OB BLKSIZE UEBER=	1630
	LA	J,6	1090	*		EINSTIMMT.	1640
	B	LABWEI	1100	M6	LTR	NEWOLD,NEWOLD	1650

	BZ	GTMN	1660
*	LABEL	VORHANDEN?	1670
	TM	JFCB+66,X'11'	168C
	BZ	M5	1690
	B	GTMN	1700
M5	EQU	*	1710
*	DIRECT	ACCESS DEVICE OR MAGNETIC TAPE?	1720
	L	Q,16(DDLAUF) ADRESSE VON UCB.	173C
	TM	18(Q),X'20' DIRECT ACCESS DEVICE?	1740
	BO	M8	175C
	TM	18(Q),X'80' MAGNETIC TAPE?	176C
	BO	M9	1770
	B	FEHLER7	178C
	EQU	*	1790
M9	*	MAGNETIC TAPE.	180C
	OPEN	DCBALL	1810
	RDJFCB	DCBALL	1820
*	RDJFCB	ERFOLGREICH?	183C
	LTR	15,15	1840
	BNZ	FEHLER7	1850
	CLOSE	DCBALL	186C
	L	Q,JFCB+100 BLKSIZE	1870
	B	M13	188C
	EQU	*	1890
M8	*	LESEN VON FORMAT 1 DSCB. ZUERST VOL-SER-NUMBER AUS	1900
*	UCB	NACH VOLSER SPEICHERN. Q ENTHAELT ACRESSE VON UCB.	1910
	MVC	VOLSER(6),28(Q) VOL-SER-NUMBER MOVEN.	1920
	OBTAIN	DSCB	193C
*	OBTAIN	ERFOLGREICH?	1940
	LTR	15,15	195C
	BNZ	FEHLER7	196C
	L	Q,WKA+40	1970
M13	SLL	Q,16	198C
	SRL	Q,16	1990
	B	D(BRANCH1)	200C
M7	C	Q,BLKSZ	2010
	BNE	MSG1	2020
	B	GTMN	2030
NOBLKS	EQU	*	2040
*	WENN	DATASET OLD UND LABEL VORHANDEN BRANCH1 MIT M2 LADEN	205C
*	BRANCH2	MIT GTMN LADEN UND SPRUNG NACH M5.	2060
*	WENN	DATASET OLD UND KEIN LABEL VORHANDEN SPRUNG NACH	2070
*	MSG4.		2080
*	WENN	DATASET NEW PRUEFEN OB DDNAME GLEICH FT07F001.	209C
*	WENN	JA BLKSIZE GLEICH 80 SETZEN. SPRUNG NACH GTMN.	210C
*	WENN	DATASET NEW UND DDNAME GLEICH FT05F001 ELKSIZE GLEICH 80	2110
*	SETZEN,	WENN DDNAME GLEICH FT07F001 BLKSIZE GLEICH 936 SETZEN,	212C
*	SONST	BLKSIZE GLEICH 800 SETZEN. DANACH SPRUNG NACH GTMN.	2130
*	BENUTZER	WIRD BENACHRICHTIGT VOM EINSETZEN DER DEFAULT-WERTE.	2140
	LA	BRANCH1,M2	2150
	LA	BRANCH2,GTMN	2160
*	LABEL	VORHANDEN?	2170
	TM	JFCB+66,X'11'	218C
	BZ	M11	2190
	LTR	NEWOLD,NEWOLD	2200

	BP	MSG4	2210
	B	M12	2220
M11	LTR	NEWOLD,NEWOLD	223C
	BP	M5	2240
M12	CD	O,FT07F001	225C
	BE	SET80	2260
	CD	O,FT05F001	2270
	BE	SET80	2280
	CD	O,FT06F001	2290
	BE	SET936	230C
	B	SET800	2310
SET80	LA	RBLKS,80	2320
	B	NAC	2330
SFT936	LA	RBLKS,936	2340
	B	NAC	235C
SET800	LA	RBLKS,800	2360
NAC	EQU	*	2370
*	DDNAME	IN NACHRICHT MOVEN. BLKSIZE IN NACHRICHT MOVEN.	2380
	MVC	WT05+51(8),DDNAM	2390
	LTR	RBLKS,RBLKS	240C
	BM	FEHLER7	2410
	C	RBLKS,=F'9999'	2420
	RH	FEHLER7	2430
	CVD	RBLKS,DHILF	2440
	LA	15,HIER	245C
STEZA	SR	I,I	2460
	L	J,DHILF+4	247C
	SRL	J,4	2480
STEZA1	LA	I,1(1)	2490
	SRL	J,4	2500
	LTR	J,J	2510
	RP	STEZA1	2520
	UNPK	DHILF(5),DHILF+5(3)	2530
	MVZ	DHILF+4(1),DHILF+3	254C
	BR	15	255C
HIER	C	I,=F'4'	256C
	BE	MOV4	2570
	C	I,=F'3'	2580
	BE	MOV3	2590
	C	I,=F'2'	2600
	BE	MOV2	261C
	C	I,=F'1'	2620
	BNE	FEHLER7	2630
MOV1	MVC	WT05+72(1),DHILF+4	264C
	B	WT05	2650
MOV2	MVC	WT05+71(2),DHILF+3	266C
	B	WT05	2670
MOV3	MVC	WT05+70(3),DHILF+2	2680
	B	WT05	269C
MOV4	MVC	WT05+69(4),DHILF+1	2700
WT05	WTO	' MISSING BLKSIZE IN CDCARD WITH DDNAME=	. BL\$ 2710
	KSIZE	ASSUMED.',ROUTCDE=11	2720
GTMN	EQU	*	273C
*	RBLKS	ENTHAELT BLKSIZE. ES WIRD ANGENOMMEN, DASS	2740
*	2	PUFFER UND EIN KONTROLLBLOCK MIT 8 BYTE ANGELEGT	2750



*	WERDEN SOLLEN.	2760	LTR	Q,Q	3310
*	WENN BUFL GROESSER ALS BLKSIZE WIRD BUFL GENOMMEN.	2770	BNZ	MSG3	3320
	L Q,JFCB+88	2780	TM	ERROR,X'01'	3330
	SLL Q,16	2790	BD	FEHLER8	3340
	SRL Q,16	2800	L	Q,PARMAER	3350
	CR RBLKS,Q	2810	ST	REGION,0(Q)	3360
	BNL TAKEBL	2820	L	13,4(13)	3370
	LR I,Q	2830	RETURN	(14,12),T,RC=0	3380
	TM JFCB+92,X'01'	2840	RFTURN1	L Q,PARMAER	3390
	BD ALIGN4	2850	ST	Q,Q(Q)	3400
	SRL Q,3	2860	L	13,4(13)	3410
	SLL Q,3	2870	RETURN	(14,12),T,RC=4	3420
	CR I,Q	2880	FEHLER2	L Q,=F'-2'	3430
	BE SETBLK	2890	B	RETURN1	3440
	LA Q,8(Q)	2900	FEHLER5	L Q,=F'-5'	3450
	B SETBLK	2910	B	RETURN1	3460
ALIGN4	SRL Q,2	2920	FEHLER7	L Q,=F'-7'	3470
	SLL Q,2	2930	B	RETURN1	3480
	CR I,Q	2940	FEHLER8	L Q,=F'-8'	3490
	BE SETBLK	2950	R	RETURN1	3500
SETBLK	LA Q,4(Q)	2960	MSG1	EQU *	3510
TAKEBL	LR RBLKS,Q	2970	LA	I,MSG10	3520
	AR RBLKS,RBLKS	2980	B	MSGW	3530
	LA RBLKS,8(RBLKS)	2990	MSG3	EQU *	3540
	A RBLKS,=XL4'80000000'	3000	*	REGION MUSS UM MINDESTENS Q BYTE ERFOEDT WERDEN.	3550
	ST RBLKS,PUFFER	3010	B	FEHLER8	3560
	GETMAIN LC,LA=PUFFER,A=ADRESS	3020	MSG4	EQU *	3570
	LTR 15,15	3030	LA	I,MSG40	3580
	BZ M10	3040	B	MSGW	3590
*	PUFFER=FRVGM ZU MISREG ADDIEREN.	3050	MSGW	EQU *	3600
	L Q,PUFFER	3060	*	DDNAMEN IN MESSAGE MOVEN.	3610
	LA Q,0(Q)	3070	MVC	WT01+41(8),DDNAM	3620
	S Q,FRVGM	3080	WT01	' ERROR IN DD CARD WITH CCNAME=	3630
	BM FEHLER7	3090	B	0(I)	3640
	A Q,MISREG	3100	MSG10	EQU *	3650
	ST Q,MISREG	3110	*	BLKSIZE AUS DD-KARTE UNGLEICH BLKSIZE AUS	3660
	B ERHOEHE	3120	*	OLD-DATASET-LABEL.	3670
M10	EQU *	3130	*	Q ENTHAELT BLKSIZE AUS OLD DATASET LABEL	3680
*	FREIE REGION NACH GETMAIN FESTSTELLEN.	3140	LTR	Q,Q	3690
	LA 1,PARM2	3150	BM	FEHLER7	3700
	L 15,=V(FREE SP)	3160	C	Q,=XL4'0000FFFF'	3710
	BALR 14,15	3170	RH	FEHLER7	3720
*	DIFFERENZ ZWISCHEN VORHER UND NACHHER IN FREIER	3180	CVD	Q,DHILF	3730
*	REGION BESTIMMEN.	3190	LA	15,DA	3740
	L Q,FRVGM	3200	B	STEZA	3750
	S Q,FRNGM	3210	DA	EQU *	3760
	BM FEHLER7	3220	C	I,=F'5'	3770
	AR REGION,Q	3230	BE	MOVDA5	3780
	FREEMAIN L,LA=PUFFER,A=ADRESS	3240	C	I,=F'4'	3790
ERHOEHE	AR DD LAUF,DD INCR	3250	RE	MOVDA4	3800
	B WEITER	3260	C	I,=F'3'	3810
ADD2K	LA REGION,2(REGION)	3270	BE	MOVDA3	3820
	B ERHOEHE	3280	C	I,=F'2'	3830
RETURN	EQU *	3290	BE	MOVDA2	3840
	L Q,MISREG	3300	C	I,=F'1'	3850

	BE	MOVDA1	3860
	B	FEHLER7	3870
MOVDA1	MVC	WTO4+55(1),DHILF+4	3880
	B	WTO4	389C
MOVDA2	MVC	WTO4+54(2),DHILF+3	3900
	B	WTO4	391C
MOVDA3	MVC	WTO4+53(3),DHILF+2	3920
	B	WTO4	3930
MOVDA4	MVC	WTO4+52(4),DHILF+1	394C
	B	WTO4	3950
MOVDA5	MVC	WTO4+51(5),DHILF	396C
WTO4	WTO	' BLKSIZE OF THE ( CLC ) DATASET EQUALS	BYTE.'\$ 3970
		,ROUTCDE=11	398C
	B	MSG	399C
MSG40	EQU	*	4000
*		DATASET IST OLD. KEIN LABEL VORFANDEN. KEINE BLKSIZE	4C1C
*		IST AUF DER DDKARTE ANGEGBEN.	4020
WTO2	WTO	' MISSING BLKSIZE ON DD CARD FOR OLD DATASET WITH NO \$	4030
		LABEL.',ROUTCDE=11	404C
	B	MSG	4050
MSG	MVI	ERROR,X'01'	4060
	B	ERHOEHE	4070
DHILF	DC	D'0'	4080
FTAUS	DC	C'FT00F001'	409C
SYSAB	DC	C'SYSABEND'	410C
SYSUD	DC	C'SYSUDUMP'	411C
FT05F001	DC	C'FT05F001'	4120
FT06F001	DC	C'FT06F001'	4130
FT07F001	DC	C'FT07F001'	414C
DDNAM	DC	D'0'	4150
TICTACR	DS	F	4160
PUFFER	DC	F'0'	4170
ERROR	DC	F'0'	4180
BLKSZ	DC	F'0'	419C
EXL	DC	XL1'87'	4200
	DC	AL3(JFCB)	4210
ADRESS	DC	F'0'	4220
MISREG	DC	F'0'	4230
FRVGM	DC	F'0'	424C
FRNGM	DC	F'0'	4250
PARMAER	DC	F'0'	4260
PARM1	DC	XL1'80'	4270
	DC	AL3(FRVGM)	4280
PARM2	DC	XL1'80'	429C
	DC	AL3(FRNGM)	4300
DCBALL	DCB	DSORG=PS,MACRF=R,EXLST=EXL	4310
JFCB	DC	44F'0'	4320
DSCB	CAMLST	SEARCH,JFCB,VOLSER,WKA	4330
VOLSER	DC	CL8'00000000'	4340
WKA	DS	DD	435C
	DS	148C	4360
SAVE	DS	18F	4370
	END		438C

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Section 3

Production of KEMA-input  
from BRIGITTE-output by  
the program COPEND

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## Survey

To insert the translated (see BRIGITTE section 2.) ENDF/B-data into the local public KEDAK-library, the program COPEN is used. This program produces for one or several materials KEMA input (see II) out of the translated ENDF/B-data, which are stored in a KEDAK-library created by BRIGITTE (see 2.2.4). To control the program only the material name (resp. names) is (resp. are) necessary. All datatypes available for a specified material are processed. The output of the program COPEN can be read by KEMA (see II) without performing any changes.

The following description is in German, because the program has been first described by the author<sup>+</sup> in an internal report, from which the text was taken.

For the user of the program we will give some helps in English. The input and the JCL-cards (Job Control Cards) are described in 3.2.2 and 3.2.3. The JCL-cards are selfexplaining (for an IBM-user). The input consists of the number of materials, which are to be copied, followed by the material names (see 3.2.3 after //G.SYSIN DD \*). The rest of this section describes the subroutines of COPEN, which are also listed as source programs.

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<sup>+</sup> The author has left the KFK.

Inhalt

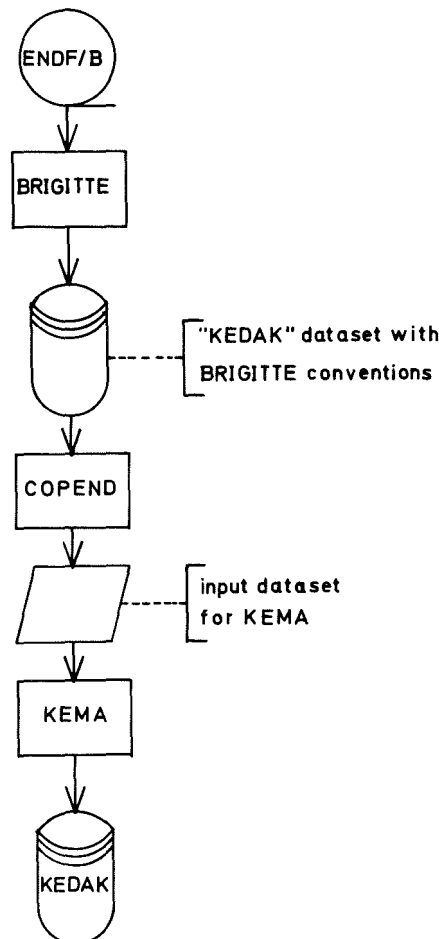
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### 3.1 Einleitung

Das in Abschnitt 2. beschriebene Programm BRIGITTE überträgt Daten aus der ENDF-Darstellung in die KEDAK-Darstellung.

BRIGITTE schreibt diese Daten nicht direkt in die Kerndatenbibliothek KEDAK, sondern im FORMAT<sup>+</sup> der KEDAK Bibliothek auf eine temporäre Datei, um die Korrektur möglicher Fehler in den konvertierten Daten zu erleichtern.

Aus diesem Grund ist ein Anschlußprogramm erforderlich, das es gestattet, diese konvertierten Datensätze in die offizielle KEDAK-Bibliothek zu übernehmen. Das hierfür geschriebene Programm COPENDE ist im folgenden beschrieben. Es erstellt aus den konvertierten Datensätzen eine Datei im Eingabeformat für das Verwaltungsprogramm der Kerndatenbibliothek KEMA (siehe II oder /2/), mit dessen Hilfe die Daten schließlich in KEDAK eingebracht werden. Nachfolgendes Flußdiagramm soll diesen Ablauf veranschaulichen.



---

<sup>+</sup> Dieses Format unterscheidet sich geringfügig von dem "KEMA" KEDAK-Format (siehe 2.2.4).

### 3.2 Das Programm COPEND

#### 3.2.1 Die Funktion des Programmes COPEND

COPEND ist ein Programm, welches ENDF/B-Daten /1/ aus einer von BRIGITTE (siehe 2.) erstellten Datei in eine Datei im Eingabeformat für KEMA (siehe II oder /2/) umwandelt.

Die Datei der konvertierten ENDF/B-Daten wird auf Einheit Nr. 1 ( $\hat{=}$  data set reference number = 1) erwartet, die Datei im KEMA-Eingabeformat wird auf die Einheit Nr. 10 geschrieben. Die Eingabe für COPEND ist formatfrei, es gelten die FREEFO-Konventionen /3/. Für FREEFO muß die Einheit Nr. 8 bereitgestellt werden.

Das Programm COPEND benutzt das Retrieval Paket LDFPAC (siehe III.2) und die Subroutine FREEFO /3/.

#### 3.2.2 Beschreibung der Eingabe

NMAT	≤ 30, Anzahl der von der Datei der ENDF/B-Daten zu kopierenden Isotope (Materialien)
NATNAM(I) (I=1,NMAT)	Doppelworte, die Materialnamen in KEDAK-Konvention (siehe II, III.1, /2/) enthalten.

#### 3.2.3 Die Jobkontrollkarten

Liste der Jobkontrollkarten mit einem Eingabebeispiel:

```
//INRO48PA JOB (0048,101,P6M1A),LANGNER,CLASS=A
// EXEC FGG,LIB=NUSYS,NAME=COPEND
//G.FTO8FOO1 DD UNIT=SYSDA,SPACE=(TRK,1)
//G.FTO1FOO1 DD UNIT=2314,VOL=SER=GFKO50,DSN=KEDAK3,DISP=SHR
//G.FT10FOO1 DD UNIT=SYSDA,SPACE=(TRK,20)
//G.SYSIN DD +
2 'AL 27 ' 'C 12 '
/*
```

Das Programm COPEND befindet sich in der Bibliothek LOAD.NUSYS; es ist 74 k Bytes lang



### 3.3 Die einzelnen Subroutinen von COPEND

#### 3.3.1 Die Subroutine INPDAT

##### 3.3.1.1 Die Organisation der ENDF/B-Daten für WRIREC

Die Subroutine INPDAT wird für jedes gewünschte Isotop einmal aufgerufen. Mit Hilfe der Subroutine LDFITN (siehe III.2) wird festgestellt, welche Reaktionstypen auf dem von BRIGITTE erzeugten KEDAK-File vorhanden sind; die Namen dieser Typen werden in dem COMMON /TYPE/ im Feld TYPES abgelegt.

Die Daten für jeden vorhandenen Reaktionstyp werden von den Subroutinen LOCDAT (s. 3.3.3) und DAT (s. 3.3.4) in das Feld X gebracht, und von der Subroutine WRIREC (s. 3.4.2) auf die Einheit Nr. 10 geschrieben.

##### 3.3.1.2 Die Argumentliste der Subroutine INPDAT

Der Aufruf der Subroutine INPDAT lautet:

```
CALL INPDAT(MATNA,NTYP,MAXNUM,NTAPE)
```

MATNA	- Doppelwort, Isotopname (KEDAK-Konvention)
NTYP	- Anzahl der Reaktionstypen in der Liste ((s. 3.2.2), z.Z. = 40)
MAXNUM	- maximale Anzahl der Daten, die für einen Eingabesatz für KEMA geliefert werden dürfen
NTAPE	- Nummer der Einheit, auf welche die umge- wandelten Daten geschrieben werden sollen

#### 3.3.2 Die Subroutine WRIREC

##### 3.3.2.1 Schreiben eines Satzes im KEMA-Eingabeformat

Die Subroutine WRIREC schreibt jeweils einen Satz im KEMA-Eingabeformat auf die Einheit NTAPE und berücksichtigt dabei, daß die Anzahl der Worte in diesem Satz nicht größer als 2000 sein darf. Außerdem werden diese Daten über die Ausgabeeinheit Nr. 6 ausgedruckt.

Das Herausschreiben auf NTAPE geschieht über die Subroutine WRIII (N,Z,NTAPE). Dabei ist  $N \leq 2000$  die Anzahl der Worte im Satz, Z(N) das Feld der Daten des Satzes, NTAPE die Nummer der Einheit, auf die geschrieben wird.

### 3.3.2.2 Die Argumentliste der Subroutine WRIREC

Der Aufruf für WRIREC lautet:

```
CALL WRIREC(NARG,NAMES,X,NUMX,NTAPE)
```

NARG	Feld von 3 Worten, in NARG(1) steht die Anzahl der Namen für die gewünschte Namenskombination (vgl. 3.3.3.2)
NAMES	Doppelwortfeld für die Namen (KEDAK-Konvention) (vgl. 3.3.3.2)
X	Feld für die Daten (Dimension $\leq 2000$ )
NUMX	Anzahl der gelieferten Daten
NTAPE	die Nummer der Einheit, auf die geschrieben werden soll.

### 3.3.2.3 Der Aufbau des herausgeschriebenen Satzes

Die von WRIREC geschriebenen Sätze sind ADD-Sätze für KEMA (II, /2/). Sie haben den folgenden Aufbau:

N	Anzahl der Worte im Satz
ADD	= 'ADD ', Doppelwort
NNAM	Anzahl der Namen in der Namenskombination
NAMES(I) (I=1,NNAM)	Namen (KEDAK-Konvention)
NARG	Anzahl der Argumente
NWERT	Anzahl der Werte
X(I) (I=1,NUMX)	Feld der Daten

Anmerkung: Für NARG = 0, wird als formales Argument der Wert 'ARG ' gesetzt.

### 3.3.3 Die Subroutine LOCDAT mit Entry NXTDAT

#### 3.3.3.1 Die Funktion der Subroutine LOCDAT

Die Subroutine LOCDAT mit Entry NXTDAT liest Daten für einen beliebigen Reaktionstyp von einer Datei im KEDAK-Format in ein Feld X.

LOCDAT wird für den ersten Aufruf einer Namenskombination benutzt. Für eventuell nötige weitere Aufrufe für die gleiche Namenskombination muß das Entry NXTDAT benutzt werden. LOCDAT verwendet die Subroutinen LDFLOC und LDFNXT (siehe III.2).

Das Einlesen wird abgebrochen, wenn das Feld X soweit gefüllt ist, daß ein weiterer logischer Satz der KEDAK-Bibliothek, bestehend aus (NARG + NWERT) Daten, keinen Platz darin findet (vgl. 3.3.2.3).

#### 3.3.3.2 Argumentliste der Subroutine LOCDAT

Der Aufruf lautet:

```
CALL LOCDAT(NARG,NAMES,NAMNXT,X,NUMX,MAXNUM,NR)
```

NARG	Feld der Länge 3. NARG(1) enthält die Anzahl der Namen des gewünschten Datentypes, NARG(2) und NARG(3) werden mit der Anzahl der Argumente und Werte des Datentyps gefüllt.
NAMES	Feld von Doppelworten. Muß vor dem Aufruf mit den alphanumerischen Namen des gewünschten Datentyps gefüllt werden
NAMNXT	falls NR = 2, enthält NAMNXT die Namen des als nächsten zu lesenden Datentyps
X	Feld, das mit den gelesenen Daten gefüllt wird.
NUMX	Anzahl der gelesenen Datenworte
MAXNUM	Länge des Feldes X

NR	Returncode
	= 0 für die angegebenen Namen gibt es noch weitere Daten, die im Feld X keinen Platz mehr fanden; zum Lesen dieser Daten ist NXTDAT aufzurufen
	= 2 der gewünschte Datentyp hat auch numerische Namen, für die angegebenen numerischen gibt es keine weiteren Daten, aber NAMNXT enthält die numerischen Namen der als nächstes zu lesenden Daten; hierfür ist LOCDAT mit den geänderten Namen zu benutzen
	= 3 Datentyp hat keine numerischen Namen, alle Daten für den gewünschten Datentyp wurden übertragen

Der Returncode wird von der Subroutine DAT (vgl. 3.3.4) interpretiert und verändert.

### 3.3.3.3 Argumentliste des Entry NXTDAT

Der Aufruf für NXTDAT lautet:

```
CALL NXTDAT(NUMX,NR)
```

Die Argumente haben die gleiche Bedeutung wie in 3.3.3.2 beschrieben. Das Entry NXTDAT ist immer dann aufzurufen, wenn der vorher von LOCDAT oder NXTDAT zurückgegebene Returncode NR = 0 war.

### 3.3.4 Die Subroutine DAT und Entry REPDAT

#### 3.3.4.1 Aufgabe der Subroutine DAT

Unter Verwendung von LOCDAT und NXTDAT stellen DAT und REPDAT die Daten für einen im Aufruf spezifizierten Datentyp in einem im Aufruf angegebenen Datenfeld zur Verfügung.

Für jeden Datentyp ist zunächst ein Aufruf von DAT erforderlich. Um die Daten des Datentyps vollständig zu lesen, ist REPDAT solange aufzurufen, wie der zuletzt zurückgegebene Returncode  $\leq 1$  ist (s.u.).

Falls für einen Datentyp numerische Namen existieren, sind diese beim ersten Aufruf (von DAT) mit Null zu füllen. Die tatsächlichen numerischen Namen werden von DAT und REPDAT eingefüllt.

#### 3.3.4.2 Die Argumentliste der Routine DAT

Der Aufruf von DAT lautet:

CALL DAT(NARG,NAMES,X,NUMX,MAXNUM,NR)

NR

Returncode

= 0 für die spezifizierte Kombination von alphanumerischen Namen gibt es noch weitere Daten, die im Feld X keinen Platz mehr fanden; zum Lesen dieser Daten ist REPDAT aufzurufen.

= 1 für die angegebene Kombination von alphanumerischen Namen sind alle Daten in das Feld X übertragen worden; für diesen Datentyp gibt es jedoch noch weitere Kombinationen von numerischen Namen; um diese numerischen Namen und die Daten hierfür zu erhalten, ist REPDAT aufzurufen

= 2,3 alle Daten des gewünschten Datentyps wurden in das Feld X übertragen.

Bemerkung: Die anderen obigen Argumente sind in 3.3.3.2 beschrieben.

#### 3.3.4.3 Die Argumentliste der Routine REPDAT

Der Aufruf lautet:

CALL REPDAT(NUMX,NR)

Bezüglich der Interpretation der Argumente s.o. REPDAT wird immer dann aufgerufen, wenn der Returncode des vorhergehenden Aufrufes von DAT oder REPDAT  $\leq 1$  war.

Literaturhinweise

- /1/ M.K. Drake, Data Formats and Procedures for the ENDF Neutron Cross Section Library, National Neutron Cross Section Center, BNL 50274 (T-601), 1970
- /2/ B. Krieg, Handling and Service Programs for the Karlsruhe Nuclear Data File KEDAK. Part I: Management and Retrieval Programs. Institut für Neutronenphysik und Reaktortechnik, Gesellschaft für Kernforschung mbH., Karlsruhe, KFK 1725, Juni 1973

### 3.5 Quellprogrammlisten

```
C
C
C          PROGRAM COPENO
C  THIS PROGRAM WRITES ENDF/B DATA IN KEMA-INPUT-FORMAT ON A DATA SET
C  WITH THE REFERENCE NUMBER = 10
C
C  COMMON /TYPE/ TYPES
C  REAL*8 NAMES(5),TYPES(70),MATNAM(30),MATNA,ENDE
C  DIMENSION F(100),NF(100),NP(70)
C  EQUIVALENCE (F(1),NF(1))
C  DATA ENDE/8HENDE /
C
C  KNDF=1
C  NIN=5
C  NCUT=6
C  INP=8
C  NTAPE=10
C
C  MAXNUM=1970
C  NT=70
C
C  CALL LDFOPN (KNDF,NADAT,&1000)
C
C  CALL FREEFO(NIN,INP,NOOUT,F,F,F)
C
C  READ (INP) NMAT,(MATNAM(I),I=1,NMAT)
C  DO 10 IMAT=1,NMAT
C  MATNA=MATNAM(IMAT)
C
C  CALL LDFITN(NR,MATNA,TYPES,NP,NTYP,NT,2)
C  IF(NR.EQ.0) GO TO 11
C  CALL SHELL(TYPES,NTYP,1,NTYP)
C
C  CALL INPDAT(MATNA,NTYP,MAXNUM,NTAPE,&1000)
C  GO TO 10
C
C  11 IF(NTYP.GT.70) GO TO 12
C  WRITE (NOOUT,101) MATNA
C  101 FORMAT(1X,'FUER ',A8,' SIND KEINE DATENTYPEN VORHANDEN')
C  GO TO 10
C  12 WRITE (NOOUT,102) MATNA,NT
C  102 FORMAT (1X,'FUER ',A8,' SIND ',I4,' (MAX=70) DATENTYPEN VORHANDEN'
C  1)
C  10 CONTINUE
C
C  N=2
C  WRITE (NTAPE) N,ENDE
C  1000 STOP
C  END
```

```
C
C
SUBROUTINE INPDAT(MATNA,NTYP,MAXNUM,NTAPE,*)
C
COMMON /TYPE/ TYPES
DIMENSION NARG(3),X(1990)
REAL*8 TYPES(70),MATNA,NAMES(10),BLANK
DATA BLANK/8H /
C
NOUT=6
NARG(1)=2
C
DO 10 ITYP=1,NTYP
IN=0
NAMES(1)=MATNA
NAMES(2)=TYPES(ITYP)
C
CALL DAT (NARG,NAMES,X,NUMX,MAXNUM,NR)
C
NZ=NARG(1)
IF(NR-3)11,12,13
12 WRITE (NOUT,101)(NAMES(L),L=1,NZ)
101 FORMAT(1X,'NO DATA FOR THIS NAME CCMBINATION - ',4A8)
GO TO 10
C
15 WRITE (NOUT,102) (NAMES(L),L=1,NZ)
102 FORMAT(1X,'DIMENSION OF THE WORKAREA FOR LDFLCC OR LDFNXT TO SMALL, THE REACTION TYPE IS ',4A8)
GO TO 10
C
16 CALL REPDAT(NUMX,NR)
IF(NR-3)11,12,13
C
11 CALL WRIREC (NARG,NAMES,X,NUMX,MAXNUM,NTAPE)
IF (NR.EQ.1) GO TO 17
IF (NR.EQ.0) GO TO 16
IF (NR.EQ.2) GO TO 20
WRITE (NOUT,103) NR,(NAMES(L),L=1,NZ)
103 FORMAT(1X,'RETURN-CODE = ',I3,'THE REACTION TYPE = ',4A8)
RETURN 1
C
13 IF(NR.EQ.4) GO TO 15
17 IN=IN+1
GO TO 16
C
20 IF(NZ.GT.2) IN=IN+1
WRITE (NOUT,104) (NAMES(L),L=1,2),IN
104 FORMAT (1X,'THE REACTION TYPE ',A8,1X,A8,' HAS ',I5,' DIFFERENT COMBINATIONS OF SECONDARY NAMES')
10 CONTINUE
RETURN
END
```



```
C
C
C      SUBROUTINE WRIREC (NARG,NAMES,X,NUMX,MAXNUM,NTAPE)
C
C      WRITREC WRITES A RECRD IN KEMA-INPUT-FORMAT ON A DATA SET WITH
C      THE REFERENCE NUMBER NTAPE, THE NUMBER OF FULLWORDS IN THE RECORD
C      MAY BE LESS THAN 2000
C
C      REAL *8 NAMES(1),ADD,ZD(2)
C      DIMENSION NARG(1),X(1)
C      DIMENSION Z(2100),ZX(4),IZ(2100)
C      DATA ADD/8HADD      /,ARG/4HARG /
C      EQUIVALENCE (Z(1),ADD),(Z(3),NNAM),(ZX(1),ZD(1)),(Z(4),ZX(1)),
C      1(Z(1),IZ(1))
C
C      NOUT=6
C      NNAM=NARG(1)
C      NARG2=NARG(2)
C      NWERT=NARG(3)
C      NNAM2=NNAM-2
C      ZD  (1)=NAMES(1)
C      ZD  (2)=NAMES(2)
C
C      IF(NUMX-MAXNUM)14,14,15
C      15 WRITE (NOUT,112) NUMX
C      112 FORMAT(1X,'THE NUMBER OF DATA =',I6,' IS GREATER THAN ALLOWED')
C      GO TO 3333
C
C      14 IF(NNAM2)11,12,13
C      11 WRITE (NOUT,111) NAMES
C      111 FORMAT(1X,'THE NUMBER OF NAMES LESS THAN 2',4A6)
C      GO TO 3333
C
C      NUMBER OF NAMES EQUAL 2
C
C      12 N=NUMX+9
C      IZ(8)=NARG2
C      IZ(9)=NWERT
C      I1=9
C      ASSIGN 19 TO I2
C      IF (NARG2.EQ.0) GO TO 16
C
C      18 DC 20 I=1,NUMX
C      20 Z(I1+I)=X(I)
C      GO TO I2,(19,21,22,23,25)
C      16 I1=10
C      N=N+1
C      Z(I1)=ARG
C      ASSIGN 21 TO I2
C      GO TO 18
C
C      THE NUMBER OF NAMES GREATER THAN 2
C
C      13 N=NUMX+9+NNAM2
```

```
DO 10 INAM=1,NNAM2
10 Z(INAM+7)=NAMES(2+INAM)
   I3=8+NNAM2
   IZ(I3)=NARG2
   IZ(I3+1)=NWERT
   I1=I3+1
   IF(NARG2.EQ.0) GO TO 24
   ASSIGN 22 TO I2
   IF(NNAM2.EQ.1) ASSIGN 25 TO I2
   GO TO 18
C
24 N=N+1
   I1=I1+1
   Z(I1)=ARG
   ASSIGN 23 TO I2
   GO TO 18
C
19 WRITE (NOUT,1) N,(Z(I),I=1,N)
   GO TO 17
21 WRITE (NOUT,3) N,(Z(I),I=1,N)
   GO TO 17
22 WRITE (NOUT,2) N,(Z(I),I=1,N)
   GO TO 17
23 WRITE (NOUT,4) N,(Z(I),I=1,N)
   GO TO 17
25 WRITE (NOUT,5) N,(Z(I),I=1,N)
   1 FORMAT(1X,I5,1X,2A4,1X,I5,1X,2A4,1X,2A4,1X ,2I5/(1X,10E13.5))
   2 FCRMAT(1X,I5,1X,2A4,1X,I5,1X,2A4,1X,2A4,1X ,2E15.6,2I5/(1X,10E13.5
   *) )
   3 FORMAT(1X,I5,1X,2A4,1X,I5,1X,2A4,1X,2A4,1X,2I5,1X,A4/(1X,10E13.5))
   4 FORMAT(1X,I5,1X,2A4,1X,I5,1X,2A4,1X,2A4,1X,2E15.6,2I5,1X,A4/(1X,10
   *E13.5))
   5 FCRMAT(1X,I5,1X,2A4,1X,I5,1X,2A4,1X,2A4,1X,E15.6,2I5/(1X,10E13.5))
C
17 CALL WRIT1 (N,Z,NTAPE)
3333 RETURN
   END

C
C
SUBROUTINE WRIT1 (N,X,NTAPE)
DIMENSION X(N)
WRITE (NTAPE) N,X
RETURN
END
```

```
C
C
SUBROUTINE DAT(NARG,NAMES,X,NUMX,MAXNUM,NR)
C
REAL*8 NAMES(1),NAMNXT(10)
DIMENSION X(1),NARG(1)
C
      USE DAT,REPDAT FOR RETRIEVAL OF KEDAK DATA OF ANY TYPE,WHICH
      SHOULD BE FILLED IN SUCCESSIVE ORDER INTO X.
      USE DAT FOR FIRST CALL WITH THIS FIRST TWO NAMES.
      USE REPDAT FOR FURTHER CALLS.
      NR=1 MEANS,THAT THE LAST DATA FOR THIS NAME COMBINATION ARE
      TRANSMITTED.THERE ARE HOWEVER FURTHER DATA WITH SAME
      FIRST TWO NAMES.TO GET THESE AND THE NEW NAMES
      CALL REPDAT.
      REST OF RETURN CODE SETTING SEE IN LOCDAT.
C
      NAMZ=NARG(1)
      IF(NAMZ.LE.2) GOTO 10
      DO 5 I=3,NAMZ
5 NAMES(I)=0.
10 CALL LOCDAT(NARG,NAMES,NAMNXT,X,NUMX,MAXNUM,NR)
      NAMZ=NARG(1)
12 IF(NAMZ.LE.2) GOTO 100
      IF(NR.NE.2) GOTO 100
      DO 20 I=1,NAMZ
      IF(NAMNXT(I).EQ.NAMES(I)) GOTO 20
      GOTO 30
20 CONTINUE
      GOTO 100
30 NR=1
      GOTO 100
C
      ENTRY REPDAT(NUMX,NR)
      IF(NR.NE.1) GOTO 50
      DO 40 I=1,NAMZ
40 NAMES(I)=NAMNXT(I)
      GOTO 10
50 CALL NXTDAT(NUMX,NR)
      GOTO 12
100 RETURN
      END
```

```
C
C
SUBROUTINE LOCDAT(NARG,NAMES,NAMNXT,X,NUMX,MAXNUM,NR)
C
DIMENSION X(1),Z(20),NARG(1)
REAL*8 NAMES(1),NAMNXT(1),NAMSV(10)
C
      LOCDAT,NXTDAT MAY BE USED TO RETRIEVE DATA FOR ANY TYPE
      FROM KEDAK.THE DATA ARE FILLED IN SUCCESSIVE ORDER INTO X.
```

C           LOC DAT MUST BE USED FOR FIRST CALL WITH THIS COMBINATION  
C           OF NAMES, NXT DAT FOR FURTHER CALLS.  
C           NR ..... RETURN CODE.  
C           0 = THERE ARE FURTHER DATA FOR THIS COMBINATION OF NAMES.  
C           2 = NO SUCH DATA  
C           3 = NO DATA FOR THIS NAME COMBINATION.  
C           4 = DIMENSION OF Z EXCEEDED.  
C           READING STOPS WHEN LENGTH OF X(MAXNUM) WOULD BE EXCEEDED.  
C           RETURN CODE IS SET 0. NUMX IS SET LENGTH OF DATA RETRIEVED.  
C           IN CASE OF NR=2 NAMNXT CONTAINS THE NAMES OF THE NEXT TYPE.  
C

          NAMZ=NARG(1)  
          I=0  
          CALL LDFLOC(NERR,NARG,NAMES,Z)  
          NAMZ=NARG(1)  
          IF(NERR.EQ.0) GOTO 30  
          DO 10 L=1,NAMZ  
10       NAMSV(L)=NAMES(L)  
          LL=NARG(2)+NARG(3)  
          IF(LL.GT.20) GOTO 40  
          GOTO 21  
20       CALL LDFNXT(NERR,NARG,NAMES,Z)  
          IF(NERR.EQ.0) GOTO 25  
21       IE=I+LL  
          IF(IE.GT.MAXNUM) GOTO 35  
          DO 22 L=1,LL  
          I=I+1  
22       X(I)=Z(L)  
          GOTO 20

C           ENTRY NXT DAT(NUMX,NR)  
          I=0  
          GOTO 21

C  
25       NR=2  
          DO 26 L=1,NAMZ  
          NAMNXT(L)=NAMES(L)  
26       NAMES(L)=NAMSV(L)  
          GOTO 200  
30       NR=3  
          GOTO 200  
35       NR=0  
          GOTO 200  
40       NR=4  
          GOTO 200  
200       NUMX=I  
          RETURN  
          END

```
C
C
C
C
SUBROUTINE SHELL(LISTE,L,LGA,LGE)
C
C   SORTIEREN NACH DEM SHELL-VERFAHREN.
C
C   DIMENSION LISTE(L)
C   REAL*8 LISTE,HELP
C   INTEGER S
C
C   S=L/2
1  LGZ=LGA
2  LTZ=LGZ
3  IF(LISTE(LTZ).GE.LISTE(LTZ+S)) GOTO 4
   HELP=LISTE(LTZ)
   LISTE(LTZ)=LISTE(LTZ+S)
   LISTE(LTZ+S)=HELP
   LTZ=LTZ-S
   IF(LTZ.GE.LGA) GOTO 3
4  LGZ=LGZ+1
   IF((LGZ+S).LE.LGE) GOTO 2
   S=S/2
   IF(S.NE.0) GOTO 1
   RETURN
   END
```

Section 4

Translating from KEDAK into  
ENDF/B (Version 1) format by the  
program KTOE

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The only program, which is available for the automated translation from KEDAK to ENDF/B, is the program KTOE. A NEA abstract of this code, which is distributed by the NEA Computer Programme Library, will be given.

It is planned, either to extend this code in order to fulfill ENDF/B-IV conventions, and above all to refine the rough translation, or to develop own methods.

1. NAME OR DESIGNATION OF PROGRAMME. KTOE
2. COMPUTER FOR WHICH THE PROGRAMME IS DESIGNED AND OTHERS UPON WHICH IT IS OPERABLE. COMPLETELY MACHINE INDEPENDENT
3. NATURE OF PHYSICAL PROBLEM SOLVED. THIS CODE PERFORMS A FULLY AUTOMATED TRANSLATION FROM KEDAK INTO ENDF/B (VERSION 1) FORMAT. OUTPUT IS ON TAPE IN DECIMAL CARD IMAGE FORMAT.
4. METHOD OF SOLUTION. BEFORE TRANSLATION EACH REACTION IS CROPPED IN ORDER TO REDUCE THE AMOUNT OF DATA POINTS OVER WHICH IT IS SPECIFIED. LINEAR-LINEAR INTERPOLATION RULE IS PRESERVED. ELASTIC ANGULAR DISTRIBUTIONS ARE GIVEN IN TABULATED FORM.  $\mu_L$  AND  $\xi_{e1}$  ARE DIRECTLY CALCULATED. THE RESONANCE DATA BOTH, RESOLVED AND UNRESOLVED, ARE ALSO TRANSLATED; THIS LEADS TO A REDUNDANCY OF INFORMATION SINCE THE CORRESPONDING TABULATED CROSS-SECTIONS ARE GIVEN TOO.
5. RESTRICTIONS ON THE COMPLEXITY OF THE PROBLEM. THE EXISTING RESTRICTIONS BOTH ON KEDAK AND ENDF/B HAVE BEEN APPLIED TO THE ARRAY SIZES USED EVERYWHERE IN THE CODE.
6. TYPICAL RUNNING TIME. 3 TO 15 MINUTES PER NUCLIDE ON THE IBM 7094.
7. UNUSUAL FEATURES OF THE PROGRAMME. FIVE SIGNIFICANT FIGURES ARE ALLOWED FOR A REAL ITEM IN THE FORMAT. SECONDARY ENERGY DISTRIBUTION LAWS FOR THE (N,N) AND (N,2N) PROCESSES ARE GENERATED BY ASSUMING AN EVAPORATION MODEL WITH THE TEMPERATURE EXPRESSED AS A FUNCTION OF THE INCIDENT MASS.
8. RELATED AND AUXILIARY PROGRAMMES. DICTON SHOULD BE USED TO INSERT AN INDEX OF THE CONTENTS IN THE FILE. CHECKER SHOULD BE USED TO TEST THE TRANSLATED FILE.
9. STATUS. IN USE.



10. REFERENCES. ENDF/B AND KEDAK DESCRIPTION REPORTS.
11. MACHINE REQUIREMENTS. 32K OF DIRECTLY ADDRESSABLE CORE LOCATIONS  
2 TAPE UNITS AND 5 TEMPORARY TAPES WHICH CAN BE SIMULATED ON  
DISK OR OTHER DRIVES.
12. PROGRAMMING LANGUAGE USED. FORTRAN IV
13. OPERATING SYSTEM OR MONITOR UNDER WHICH PROGRAMME IS EXECUTED.  
NONE IN PARTICULAR.
14. ANY OTHER PROGRAMMING OR OPERATING INFORMATION OR RESTRICTIONS.  
THE CODE CAN TRANSLATE ONE NUCLIDE EACH TIME.
15. NAME AND ESTABLISHMENT OF AUTHOR. GIAN CARLO PANINI, C.N.E.N.,  
C.D.C., BOLOGNA.
16. MATERIAL AVAILABLE.
17. CATEGORY. M  
KEYWORDS. KEDAK, ENDF/B