

**KERNFORSCHUNGSZENTRUM  
KARLSRUHE**

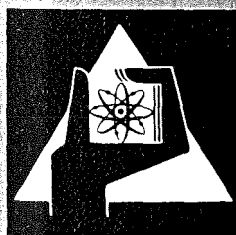
April 1977

KFK 2387/V

Institut für Neutronenphysik und Reaktortechnik  
Projekt Schneller Brüter

**The KEDAK Program Compendium  
Part V  
KEDAK Evaluation Aids**

Compiled by: F. H. Fröhner  
with contributions from  
C. Broeders, I. Broeders, B. Goel, I. Langner,  
R. Meyer, H. W. Wiese



**GESELLSCHAFT  
FÜR  
KERNFORSCHUNG M.B.H.**

**KARLSRUHE**

Als Manuskript vervielfältigt

Für diesen Bericht behalten wir uns alle Rechte vor

GESELLSCHAFT FÜR KERNFORSCHUNG M. B. H.  
KARLSRUHE

KERNFORSCHUNGSZENTRUM KARLSRUHE

KFK 2387/V

Institut für Neutronenphysik und Reaktortechnik  
Projekt Schneller Brüter

The KEDAK Program Compendium

Part V

KEDAK Evaluation Aids

Compiled by

F. H. Fröhner

with Contributions from

C. Broeders, I. Broeders, B. Goel, I. Langner,  
R. Meyer\*, H. W. Wiese

\*Present address: Software AG, Darmstadt

Gesellschaft für Kernforschung mbH., Karlsruhe



## Abstract

Part V of the KEDAK Program Compendium contains descriptions of the more important codes that are used in KEDAK evaluation work at Karlsruhe, in particular: check programs, programs performing arithmetic operations on whole data sets, programs for weighting and smoothing and programs for estimation of nuclear-model and cross section parameters.

Das KEDAK-Programm-Compendium

Teil V

Programmhilfsmittel für die KEDAK-Auswertearbeit

---

## Zusammenfassung

Teil V des KEDAK-Programm-Compendiums enthält Beschreibungen der wichtigeren Programme, die bei der KEDAK-Auswertearbeit in Karlsruhe benutzt werden - speziell Prüfprogramme, Programme für arithmetische Operationen an ganzen Datensätzen, Programme zum Wichten und Glätten und Programme für die statistische Schätzung von Kernmodell- und Wirkungsquerschnitts-Parametern.

## V. KEDAK Evaluation Aids

Contents:	page
1. <u>GENERAL REMARKS</u>	1
2. <u>CHECK PROGRAMS</u>	2
2.1 The general-purpose check program PRTSTKED	3
2.2 SELPLO: Checking and processing program for differential elastic-scattering data	25
3. <u>PROGRAMS PERFORMING ARITHMETIC OPERATIONS ON NUCLEAR DATA</u>	52
3.1 Simple averages (AVERAGE)	53
3.2 Calculation of composite cross sections (CALCUL)	88
4. <u>PROGRAMS FOR WEIGHTING AND SMOOTHING OF CROSS SECTION DATA</u>	94
4.1 Program SIGPLO	95
4.2 Karlsruhe version of SCORE	151
5. <u>PROGRAMS FOR ESTIMATION OF NUCLEAR-MODEL AND CROSS SECTION PARAMETERS</u>	213
5.1 Multi-level shape analysis program for transmission data (FANAL)	213
5.2 Multi-level shape analysis program for capture data (FANAC)	215
5.3 Estimation of level-statistical parameters from individual resonance parameters (STARA)	217
5.4 Estimation of level-statistical parameters from average cross sections (FITACS)	218
5.5 INCH3 : Coupled-channels program for inelastic and charge exchange reactions with automatic parameter optimization	220
5.6 DWBA: Program calculating inelastic-scattering and charge-exchange cross sections	222
5.7 SECDIST: Program for the calculation of energy distributions of secondary neutrons produced by inelastic scattering	223

## 1. GENERAL REMARKS

The KEDAK library of evaluated neutron data must be regularly revised, updated and expanded in response to the users' needs. A continuous evaluation effort is necessary to maintain the usefulness of the file. The following sections deal with computer programs that were developed or adapted in the course of this evaluation work. It should be pointed out that as in previous chapters the program descriptions are little more than a collection from various sources and differ considerably with respect to depth and detail. References are found after each program description.

## 2. CHECK PROGRAMS

The large size and rapid growth of modern neutron data files make it difficult to ensure formal correctness and consistency with the laws of physics in evaluated files. Efficient check programs are required to ease the burden on the compiler. Two such programs are being used in KEDAK evaluation work:

- PRTSTKED for angle-integrated cross sections and related data such as cross section ratios that depend on incident energy only, and
- SELPLO for differential elastic-scattering cross sections.

They are described in Sects. 2.1 and 2.2.



## 2.1 THE GENERAL-PURPOSE CHECK PROGRAM PRTSTKED

### Contents:

#### 2.1.1 Introduction

#### 2.1.2 Consistency checks and physics tests

##### 2.1.2.1 Tests of nuclear properties

##### 2.1.2.2 Tests on partial cross sections

##### 2.1.2.3 Consistency checks on redundant quantities

#### 2.1.3 Input/output specifications

Appendix: Brief user's guide to the KEDAK print and test program PRTSTKED

### 2.1.1 Introduction

The KEDAK print and test program PRTSTKED, developed in 1972 by R. Meyer, is designed to print a large number of KEDAK reaction types and to perform consistency checks and physics test. Among the data that cannot be tested yet are individual and statistical resonance parameters, angular distributions and secondary-particle spectra.

### 2.1.2 Consistency checks and physics tests

As explained in the appended PRTSTKED input/output description the program is sectioned into a number of modules. The actual test operations are performed in the program modules ISOPAC, TSTBL1 and TSTBL2:

- ISOPAC tests nuclear properties (charge, mass, spin, neutron binding energy, nuclear radius etc.);

- TSTBL1 checks individual energy-dependent cross sections;
- TSTBL2 checks consistency between individual partial cross sections and derived quantities such as the total cross section or the capture-to-fission ratio  $\alpha$ .

#### 2.1.2.1 Tests of nuclear properties

The following tests are performed by ISOPAC:

- Is the isotope in question on file?
- Does the charge number  $Z$  satisfy the condition  
 $1 \leq Z \leq 104$  ?
- Is  $Z$  integer?
- In case both isotopic and elemental data are stored, is the isotopic  $Z$  equal to the elemental  $Z$ ?
- Is the nucleon number  $A=N+Z$  compatible with the atomic weight?
- Is  $Z$  consistent with the chemical symbol?
- Is the combination  $Z,A$  inside the range of the nuclide chart?
- Does  $Z,A$  belong to an unstable isotope? If yes, is the isotope listed as component of an element?
- Is the atomic weight consistent with the nuclide chart?
- Is the ground state spin  $I$  integer for even  $A$ , half-integer for odd  $A$ ?  
Is  $I=0$  for even  $A$ ?
- Is the condition  $0 \leq I \leq 10$  fulfilled?
- Is the nuclear radius inside the expected range?
- Is the neutron binding energy inside the expected range?
- Is  $\lambda\sqrt{E}(1+m/M)$  consistent with the atomic weight?  
( $m$ : neutron mass,  $M$ : nuclear mass)
- Do the isotopic abundances  $a_i$  for an element fulfill the conditions  
 $0 < a_i \leq 1, \quad \sum_i a_i = 1$  ?
- Is more than 1 isotopic component given for an isotope?

### 2.1.2.2 Checks on partial cross sections

Partial cross sections must be on file in the order of ascending energy. Tests are performed by TSTBL1 as follows:

- Energy values are tested with respect to sign and order and reordered if necessary.
- The occurrence of double or multiple energy values is tested.
- Completeness of the energy range (1 meV - 15 MeV) is checked. Suspicious gaps or clusters are identified in the sequence of energies.
- It is checked whether cross section values fulfill the condition

$$10^{-10}_b < \sigma < 10^6_b$$

- In case of angular distributions the condition

$$- 1 \leq \mu = \cos \theta \leq + 1$$

is checked.

- In case of inelastic scattering thresholds are checked against excitation energies of the residual nucleus.

### 2.1.2.3 Consistency checks on redundant quantities

The following tests are performed by TSTBL2:

- Is the total cross section equal to the sum of all partial cross sections,

$$\sigma_T = \sum_c \sigma_c \quad ?$$

- Is the nonelastic cross section  $\sigma_{\text{nonel}}$  consistent with

$$\sigma_{\text{nonel}} = \sigma_T - \sigma_n = \sum_{c \neq n} \sigma_c \quad ?$$

- Is the absorption cross section  $\sigma_{\text{abs}}$  consistent with

$$\sigma_{\text{abs}} = \sigma_T - \sigma_n - \sigma_{n'} - \sigma_{2n} - \sigma_{3n} \quad ?$$

- Is the transport cross section  $\sigma_{\text{tr}}$  consistent with

$$\sigma_{\text{tr}} = \sigma_T - \bar{\mu} \sigma_n \quad ?$$

( $\bar{\mu}$ : mean cosine of lab scattering angle)

- Is the capture-to-fission ratio  $\alpha$  consistent with

$$\alpha = \frac{\sigma_\gamma}{\sigma_f} \quad ?$$

- Is the quantity  $\eta$  consistent with

$$\eta = \frac{\bar{\nu} \sigma_f}{\sigma_f + \sigma_\gamma} = \frac{\bar{\nu}}{1 + \alpha}$$

( $\bar{\nu}$ : average number of emitted neutrons per fission) ?

- In the case of cross sections for process like  $(n, n'\alpha)$ ,  $(n, \alpha 2n)$  it is checked whether

$$\sigma_{n'\alpha} < \sigma_{n'}$$

$$\sigma_{\alpha 2n} < \sigma_{2n}$$

- It is checked whether the total inelastic cross section,  $\sigma_{n'}$ , is equal to the sum over all partial inelastic cross sections

$$\sigma_{n'} = \sum_i \sigma_{n'}^{(i)} + \sigma_{n'}^{(c)}$$

$\sigma_{n'}^{(i)}$ : contribution from the  $i$ -th excited level of the target nucleus,  
 $\sigma_{n'}^{(c)}$ : contribution from excited levels in the continuum). Threshold energies are checked against the discrete level energies.

### 2.1.3 Input/output specification

A brief description of the more computer-technical aspects of the program including input/output options is appended. This description is based on an unpublished internal report written by R. Meyer, the author of the program.

Appendix

Brief User's Guide to the  
KEDAK print and test program PRTSTKED

BRIEF USER'S GUIDE TO THE KEDAK  
PRINT AND TEST PROGRAM PRTSTKED

1) Introduction

At the panel on evaluation problems in Vienna in September 1971 a number of testing programs were presented for ENDF/B and UKNDL and it was felt that such testing programs were essential in ensuring the correctness of large amounts of data. No testing program for KEDAK was available at that time.

At the end of 1971 a number of ENDF/B2 materials were converted with a very preliminary version of BRIGITTE /3/ and the lack of a testing program again turned out to be a drawback.

At about the same time it became obvious, that it would be impossible to maintain the KEDAK conventions of tabulating neutron cross section data at one and the same energy scale for all reaction types of one material. Also, this convention would be violated by the converted ENDF/B materials. Therefore, the available printing codes became obsolete.

For this reason, in December 1971 the author started programming of the printing and testing code, the first

step of which after thorough testing was operational in August 72.

This code, the input and operation of which is described in the succeeding paragraphs, was designed for printing of a large number of KEDAK reaction types and for performing a number of consistency checks and physical tests.

The version referred to hereafter is of a limited scope in that it does provide printout only of the most important data types plus a few more, but still a number of quantities cannot be handled. Also, tests are restricted to cross section data, and testing of resonance data and distribution data has not been implemented.

Since these extensions of the program had been planned from the very beginning, the program is ready to accept them. And by using the program resources already implemented, as error handling, Condition handling, editing, retrieval etc., these extensions are fairly easy to make.

However, since presently it cannot be foreseen at what time work on the program can be resumed it was decided to provide a preliminary issue of input and operation description.

A detailed description of the individual program sections is in preparation and will be published separately.

## 2) Sectioning

The program is sectioned into the following modules:

- CALPAC
- ERRPAC
- ISØPAC
- LDFPAC
- PRIPAC
- RETPAC
- TSTBL1
- TSTBL2
- TSTKED
- CØNPAC

where TSTKED contains the control program and control input processing program, LDFPAC contains an overlayable system of library-oriented retrieval routines.



RETPAC contains an overlayable system of user-oriented retrieval programs.

PRIPAC contains the editing routines for the printing option.

ISØPAC performs testing of nuclear-property data. For certain detected errors ISOPAC performs corrective actions needed for the resonance testing package (which is not yet finished).

TSTBL1 performs physics checks and numeric checks on the individual, energy-dependent cross section types.

TSTBL2 performs consistency checks on redundant quantities.

CALPAC is a package of support routines for consistency tests permitting arithmetic operations on complete cross section sets.

ERRPAC is a routine package to save detected errors and to edit error messages.

CONPAC is a routine package for handling of condition codes set by detected errors.

### 3) Unresolved references:

Routines not in the package and supplied separately:

- DEFI (A) : A routine to make the DEFINE FILE statement dynamical. A description and a FORTRAN replacement routine to simulate its functions are available.
- CONVY (A) : a routine for conversion of character strings to numeric data and vice versa. A description can be supplied on request.
- XTAREA (A) (+ Entry REXTAR): A routine to dynamically allocate core storage to the program via GETMAIN and FREEMAIN macros. A description is available. Advice how to simulate the functions of the routine with FORTRAN statements can be given on request.
- FREESP (A) : A routine to determine the free core storage available for allocation of data sets. A description can be distributed. Advice how to replace the routine by a FORTRAN routine simulating its function can be given.

DATUM (A) : A routine to retrieve the current date and time. A description is available and can be distributed on request.

Advice how to replace the routine by a FORTRAN routine simulating its function can be given.

A8FØRM (with entries RAMANF, RAMØUT) (F) : routines for printing text with capital letters (12 lines high).

All simulator routines offered do only simulate the functions of the above routines for the special purposes of the testing and printing program.

Note : (A) denotes Assembler routines  
(F) Fortran routines.

#### 4) Overlay:

An overlay package is included. Together with the dynamic allocation of core storage it ensures that the program can be run within 122K region on an IBM - 370 computer if the compilation is done by the H - extended compiler. The program requires 84K core storage with overlay and 256K without, to which 10K system+buffers and core storage for dynamic allocation must be added, amounting to at least 38K.

#### 5) JCL

The following JCL cards are required (for IBM-360 users only) 7

//JOB card

//\_EXEC\_FHLG,PARM.L = (MAP, LIST, ØVLY) if the program must be linked or

//\_EXFC\_FHG,LIB = ...

if a catalogued library is used to store the load module.

in the former case also

## 6) REGION AND TIME

The minimum region in which the program can be run (IBM-370, H-extended compiler) is about 122K with overlay and about 300K without. Execution however may be speeded up appreciably if the region is increased. The maximum region that can be used is  $8 \times NP + 16 \times NCS$  (K) where NP is the maximum number of points for a single reaction type for that material and NCS is the length of the complete energy scale (both in thousands) eg. to test Ta-181 converted from ENDF/B2 takes 48 sec of CPU time with the maximum region and 90 sec of CPU time with the minimum region, core times changing quite similarly (IBM-360/65). The running time for the printing option is not influenced by the region parameter and the option will always use 106K exactly with overlay. The use of the overlay slows down the program execution (CPU and incore time) by a factor of 1.5 - 2. Because the overlay is a multiple region overlay calls across region boundaries may occur causing additional delay. The overlay has been optimized to keep the number of such calls low. Nevertheless it may happen in very exceptional cases that execution time exceeds reasonable limits. The only advice that can be given is to cancel the overlay option in such cases.

## 7) Printing with stand alone code.

In case the printing option is the only one desired, the printing part can be run on a stand-alone basis. The overlay option may then be cancelled and the program will occupy exactly 120K.

The stand-alone code can be easily constructed by replacing routines TSTKED and TSTINP in module TSTKED by a dummy, discarding modules CALPAC, ERRPAC, ISOPAC, TSTBL1, TSTBL2 and CONPAC and selecting from RETPAC only those routines actually requested. It is suggested that the dummy routine TSTKED issues an error message if called by the control program due to erroneous control input.

```
//L.LIB_DD_defining the program library
//L.SYSIN_DD_ *
  INCLUDE cards
  ENTRY MAIN
/ *

go step:
//G.FTO1FOO1_UNIT=2314, DISP=SHR,VOL=SER=..., DSN=...
  Kedak direct access library containing the
  material to be tested/printed
//G.FTO3FOO1_UNIT=SYSDA,SPACE=(2008, 100),
  DCB=(RECFM=VBS,BLKSIZE=2008)
//G.FTO4FOO1_DD_ like FTO3
  two scratch data sets required by the testing program
  in case the incore storage is not sufficient to hold
  all data.
//G.FTO9FOO1_DD_UNIT=SYSDA,SPACE=(80,99),DCB=(RECFM=F,
  BLKSIZE=80)
  FTO9 is a scratch data set to accept the control input
  data read from SYSIN. The latter is closed immediately
  after copying.
//G.SYSIN_DD_ *
  control input
/*
//job end card
```

Notes: -An underscore in the above text is used to denote a blank.

-Do not alter any block sizes in case you do not wish to risk an 80A or an 806.

-Karlsruhe users: The load modules of the program sections can be included from the partitioned data set INR.STEIN.LOEAAD on GFKO29. Note, that TSTKED and CONPAC have been merged into one member TSTKED. The overlay package is a sequential data set INR.STEIN.OVERLAY on GFKO29.

Load modules of the complete program are stored in STEINPDS on KAPRØS under the names PRTSTKED (version without overlay) and PRTSTKEQ (version with overlay). The overlay version should only be used if the printing option is needed. If the testing option is chosen, however, input/output becomes very expensive with the overlay version and the version without overlay is recommended.

8) Control input. (++)

Control input is provided by Fortran namelist input. Two namelists are used:

PRINT for printing  
TSTKED for testing

The syntax rules are shortly summarized here.

Input is of the form:

```
_&'NAMELISTNAME' _ 'PARAMETERLIST' &END
```

where: 'NAMELISTNAME' is to be replaced by the respective namelist name (PRINT, TSTKED).

'PARAMETERLIST' is optional and may be replaced by a list of keys and associated parameters in the form:

```
'KEY1'=PARM11, PARM12,..., 'KEY2'=PARM2,...
```

where 'KEY1', 'KEY2' are to be replaced by the names of the respective keys and PARM1<sub>1</sub>, PARM1<sub>2</sub>,... are numeric or character data assigned to them. Character data must be enclosed within quotes. Numeric data are coded as usual with the trailing comma delimiting the field width (trailing blanks are treated as zeroes).

Each key is optional and if not coded its value(s) remain(s) unchanged.

&END indicates the end of the respective namelist.

Continuation cards may be used and must start in column two. However, neither keys nor parameter values may start on one card and be continued on the next:

---

(++) An underscore indicates a blank hereafter.

A) PRINT (+) :

```
_&PRINT_NAMZ='I1', NAMEN="MAT", "TYP", 'X', 'Y', EMIN='Emin',  
EMAX='Emax', FAST= $\begin{Bmatrix} T \\ F \end{Bmatrix}$ , PAGE='I2', &END
```

or

```
_&PRINT_NAMZ='I1', MAT="MAT", TYP="TYP", EXC='X', .....as  
above....., &END
```

where

NAMZ gives the number of names for the reaction type to be printed

default: NAMZ = 2

NAMEN gives the names of the reaction types to be printed.

'MAT' ist the material name

'TYP' is the reaction type name

'X', 'Y' are eventual further (numeric) names, e.g. 'X' might be an excitation energy

alternatively

MAT= "MAT" may be used to enter a material name

TYP= "TYP" may be used to enter a reaction type name

EXC= 'X1' may be used to enter a numeric name, e.g. an excitation energy.

In case more than one single data type is to be printed, alternatively the following keys may be used:

MATS="MAT1", "MAT2", ... up to ten material names my be entered.

TYPES="TYP1", "TYP2", ... up to ten reaction names may be entered.

EXCS='X1', 'X2', ... up to ten (excitation) energies may be entered.

Printing in this case will loop upon the specified reaction types as follows.

---

(+) The material and reaction type names are given in Ref. /1/, /2/.

```
'MAT1' : 'TYP1'  
        'TYP2'  
        .  
        .  
        .
```

```
'MAT2' : 'TYP1'  
        'TYP2'  
        .  
        .  
        .  
        .
```

and if one of the reaction types is SGIZ, the inelastic level excitation, printing will loop on 'X1', 'X2' given.

[Default: See note 3) ]

FAST =  $\begin{cases} T & \text{fast printing option} \\ F & \text{slow printing option: data will be edited with} \\ & \text{exponents in multiples of three (eV, keV, MeV} \\ & \text{and } \mu\text{b, mb, b, etc.).} \end{cases}$

default: FAST=F

PAGE='I<sub>2</sub>' indicates that page numbering should start with 'I<sub>2</sub>'.  
Page is used only, if a new material name is encountered.

default: PAGE=1

EMIN='Emin' are the energy limits (in eV) between which  
EMAX='Emax' the data shall be printed. Printing will start  
with the last energy  $\leq$  'Emin' and will stop  
with the first energy  $\geq$  'Emax'.  
If 'Emin' > 'Emax' no energy limits are applied.

default: EMIN=0. EMAX=-1.

- Notes: 1) The printing program only can access the following data types presently:  
AASTATUS, ISØT1, ISØT2, ISØT3, all energy-dependent tabulated cross section data and resolved-resonance parameters.  
Not included are: SGNC, LEGNC, (elastic distribution data), or unresolved-resonance data.
- 2) To print ISØT1 - ISØT3 code TYP='ISØT' or TYP = ....., 'ISØT', ....., coding 'ISØT1' or 'ISØT2' or 'ISØT3' will cause an error.
- 3) If TYP= 'ALL' is entered, all data types for the specified material are printed which the printing program can access.

NAMZ is ignored for TYP = 'ALL'.

If EXC = 'ALL' is coded and reaction type 'SGIZ' has been specified explicitly or is implied by TYP = 'ALL' printing will loop on all level energies for that material.

To print all informations for a given material accessible for the program, TYP='ALL', EXC = 'ALL' may be coded.

These are the default values for TYP and EXC. No default value for the material name is supplied.

- 4) The current value of a parameter is default and is used if the respective key is not in the input namelist. The initial defaults at program start were given above.
- 5) NAMZ is only required, if not the retrieval package LDFPAC included in the program is used. Else it will be ignored.



Alias: FROM for EMIN, TO for EMAX

Abbreviations accepted:

&P for &PRINT  
M for MAT  
T for TYP  
F for FRØM

B) TSTKED (+)

&TSTKED\_MAT='MAT', TSTESC =  $\left\{ \begin{array}{l} \text{'NØ'} \\ \text{'TYP1', 'TYP2', ...} \\ \text{'EXCLUDE', 'TYP1', ...} \\ \text{'ALL'} \end{array} \right.$

MAX='I1', EPSIL='EPS', BØUNDS='B1', 'B2', ~~&END~~  
FROM='EFROM', TO='ETO', &END

where

MAT='MAT' gives the material name. Only one per list may be entered.

MAX='I1' gives the maximum number of points for a single KEDAK type.

If the retrieval routine package of the testing program is used, this parameter is ignored and will be determined by the program.

Else if set to 0 the program will assume a number of 20 000. MAX is only important for program efficiency optimization and an erroneous specification will not lead to an error provided it is large enough. Otherwise an 80A or 806 error might occur.

default: MAX = 0

EPSIL='EPS' specifies a boundary value for the tests on the validity of linear interpolation and gives the requested accuracy of linear representation in percent.

The program will print a warning message when the shape of the (assumed) true curve deviates from the linear interpolated curve by more than 'EPS'.

---

(+) For references of material and reaction type names see /1/, /2/

default: EPSIL = 5.,  
BOUNDS = 'B1', 'B2'

gives the percent accuracy requested for the consistency of redundant quantities (in percent). Violation of consistency by more than 'B1' will cause a warning, by more than 'B2' an error message to be issued. ('B1' < 'B2')

default: B = 0.1, 5.,

FROM='EFROM' and  
TØ='ETØ' specify lower and upper boundary of the energy range in which testing is requested. For ETØ-values less than or equal to zero the data are tested at all available energies.

defaults: FROM=0.0, TØ=-1.0

TSTESC= {  
'NØ' no test on completeness of energy scale  
'TYP1', 'TYP2', ... test whether energy scales of these types are complete. Up to 5 reaction types may be entered.  
'EXCLUDE', 'TYP1', ... test completeness of energy scales of all types but 'TYP1', ... up to 4 reaction types may be entered.  
'ALL' test completeness of energy scale of all types.

Since one of the former conventions of the KEDAK library was that all reaction type data be tabulated along the same energy mesh points, TSTESC offers a possibility to test this convention. Violation of this convention in present files may be an indication also of other errors.

It is however of no use to test more recent files or files converted from ENDF/B on this convention.

default: TSTESC = 'NØ',

Abbreviation accepted:

M	for	MAT
TE	for	TSTESC
E	for	EPSIL
B	for	BØUNDS

Note:

For use of defaults see Note 4) of 8B).

9) Output:

It is hoped that the output in all cases is self-explanatory. In most cases input errors are recognized, either by the program or by the IBM namelist processing routine and will give rise to an error message.

printing module:

large letters indicating the beginning of a new material allow easy separation. The date of the Kodak version printed together with the date of printout is displayed for each material on a heading page.

Each reaction type will be preceded by a heading to identify the data being printed. The heading will be repeated on continuation pages. Normally each reaction type will start with a new page, except for SGIZ when the level data to be printed still would fit into the current page.

For easier reference data points are numbered. Numbering will restart with 1 after 9999 had been reached.

Pages are numbered consecutively, the starting number being specified by the user. The page number is preceded by the name of the material currently being printed.

At the bottom of each page a line is displayed giving the physical units of the data printed.

testing module:

Three levels of messages are maintained according to the associated degree of error: warning, error and serious error messages. To avoid too much printout, each error has been assigned a maximum number of error messages; if exceeded the user will be told but no further messages for this error are printed.

Because large amounts of data will not usually be held in storage at any given time, error messages would be mixed if on-line printing were used. Instead, errors are collected and after finishing a class of tests, the error messages are sorted and then printed.

The limitation of the number of error messages of each type, although in most cases very useful, has also some disadvantages, e.g. violation of linear interpolation within the resonance range may happen quite frequently and indicates that the number of points chosen to represent that region does not fulfill the accuracy requirements. The individual messages however will not have much importance. In the fast energy range such a violation however could indicate a mispunch or a scaling error. Due to a large number of errors detected in the resonance range these violations might not be printed, if testing is done in the whole energy range. But it is possible to limit the tests on the fast energy range by using the input parameters FROM and TO.

Each material is preceded by separating pages containing the material name in large letters and a heading page giving the date of the KEDAK version used, the date and time of test and the status of the current program control parameters.

Each section of tests is preceded by a heading page shortly informing the user of the kind of tests performed in the succeeding section, and is ended by a page giving an account of the number of errors which occurred and their error levels, where a level  $\leq 4$  pertains to warning messages,  $\leq 8$  to error messages and to serious error messages else.

Pages are numbered and the page number is preceded by the name of the material currently under test.

10) Dispatching to foreign users:

Foreign users may receive the code on request from the nuclear data evaluation group in Karlsruhe. Since the printing program may be used as a stand-alone code also, the requested program package should be indicated.

References:

/1/ D. Woll, KFK 88o, 1968

/2/ B. Krieg, KFK 1725/I, 1973

/3/ J.C. Schepers, private communications and report to  
be published.

## 2.2 SELPLO: Checking and processing program for differential elastic-scattering data

---

### Contents:

- 2.2.1 Applications of SELPLØ
- 2.2.2 SELPLØ input
- 2.2.3 SELPLØ output
- 2.2.4 Short description of SELPLØ main program and subroutines
- 2.2.5 Storage and time requirements on IBM/360-65 and /370-168
- 2.2.6 External storage, job control

### References

Appendix: SELPLØ program listing

## 2.2 SELPLO: Checking and processing program for differential elastic-scattering data

The FORTRAN IV program SELPLO was written by R. Meyer for processing, checking and display of elastic-scattering angular distributions from and for the KEDAK library. The following sections contain descriptions of

- applications,
- input (including an example),
- output,
- main program and subroutines,
- storage requirements and running times,
- external storage devices and job control language.

A program listing is appended.

### 2.2.1 Applications of SELPLO

The program permits processing of differential elastic-scattering data in the center-of-mass system. These can be

- differential elastic-scattering cross sections  $\frac{d\sigma_n(E,\Omega)}{d\Omega}$  ( $\frac{\text{mb}}{\text{sterad}}$ ),
- angular distribution functions  $p(\mu)$  ( - ).

Both data types are related by

$$\frac{d\sigma_n(E,\Omega)}{d\Omega} = \frac{\sigma_n(E)}{2\pi} p(\mu),$$

with

$$\int_{2\pi} \frac{d\sigma_n(E,\Omega)}{d\Omega} d\Omega = \sigma_n(E),$$

and

$$\int_{-1}^1 p(\mu) d\mu = 1,$$

where  $\mu$  is the cosine of the scattering angle,

$\sigma_n(E)$  the (angle-integrated) elastic-scattering cross section

KEDAK nomenclature:

SGNC for  $p(\mu)$ , with  $\mu$  given in the center-of-mass system.

SGN for  $\sigma_n(E)$ , with  $E$  given in the lab system.

Note: In the sequel the general notation SGNC will be used for both types of differential (angle-dependent) data, i.e. for  $d\sigma_n/d\Omega$  and for  $p$ .

The differential elastic scattering data may be

- read from the KEDAK library or
- read from other external sources (tape, disc) or
- calculated from Legendre coefficients that were read from cards.

The following options exist:

- plotting,
- integration over all angles and normalization check whether the integral is equal to unity (for  $p(\mu)$ ) or to the angle-integrated elastic-scattering cross section  $\sigma_n$  stored on KEDAK (for  $d\sigma_n/d\Omega$ );
- transformation to laboratory angles, calculation of the average lab cosine  $\bar{\mu}_L$  (KEDAK name: MUEL) and consistency check against MUEL as stored on KEDAK;
- generation of statistics on the deviations between  $\sigma_n$  (or  $\bar{\mu}_L$ ) as calculated from SGNC and  $\sigma_n$  (or  $\bar{\mu}_L$ ) as stored on KEDAK;
- calculation of SGNC in the center-of-mass system from Legendre coefficients for an input cosine grid with check whether this grid is fine enough;
- normalization of SGNC to unity and storage in KEMA input format for later incorporation in KEDAK.



2.2.2 SELPLO input

Input quantities other than SGNC are read according to the FORTRAN NAMELIST conventions into NAMELIST/IN/. The (card) input has the form

  &IN  MAT=...,EMIN=...,EMAX=...,other names, END

where underscores    denote blanks.

The quantities and options which can appear in the IN name list are as follows:

Name	Explanation	Default value	Type
MAT	material name (KEDAK name)		REAL*8
EMIN	lower energy limit in eV	0.	REAL*4
EMAX	upper energy limit in eV	20.E+6	"
PLOT	= $\left\{ \begin{array}{l} \text{T : plot} \\ \text{F : do not plot} \end{array} \right\}$ SGNC	F	LOGICAL*1
SIGMEL	= $\left\{ \begin{array}{l} \text{T : compare} \\ \text{F : do not compare} \end{array} \right\}$ the angular integral of SGNC with SGN as stored on KEDAK	F	"
MUL	= $\left\{ \begin{array}{l} \text{T : compare} \\ \text{F : do not compare} \end{array} \right\}$ MUEL calculated from SGNC with MUEL as stored on KEDAK	F	"
NORM	= $\left\{ \begin{array}{l} \text{T : normalize} \\ \text{F : do not normalize} \end{array} \right\}$ SGNC to unity, write result in KEMA input format on unit 2, prepare punching from unit 3.	F	"
STATIS	= $\left\{ \begin{array}{l} \text{T : generate statistics} \\ \text{F : no statistics} \end{array} \right\}$ on deviations between integrals calculated from SGNC and SGN values on KEDAK if SIGMEL=T, or between corresponding MUEL values if MUL=T.	F	"

Name	Explanation	Default value	Type
INTYP	= { 'KEDA' read SGNC from KEDAK, unit 1 'TAPE' " SGNC " tape 'CARD' " SGNC " cards	'KEDA'	REAL*4
A	atomic weight of material specified by MAT ( <sup>12</sup> C system) if it is not to be read from KEDAK		"
DRUCK	= { T : print F : do not print } Legendre coefficients read from cards and calculated SGNC values	F	LOGICAL*1
NUMKED	number ≤ 101 of cosine grid points only required if INTYP='CARD'		INTEGER*4
ITAP	unit > 8 from which SGNC values are to be read if INTYP='TAPE'.		"
PLOFOR	= { T : ink plot F : pencil plot } if PLOT=T	T	REAL*4

For INTYP='KEDAK' or INTYP='TAPE' another &IN... namelist can follow. If not, the end of input is reached. For INTYP='CARD' the name list NAMELIST/CD/ must be specified for each incident energy by

&CD\_MODE=...,ENERGY=...,other names, END.

The names that can appear in the &CD name list are as follows:

Name	Explanation	Default value	Type
MODE	= { 'LEG' : read Legendre coefficients 'TAB' : read tabulated SGCN values (not active in present code verison)	'LEG'	REAL*8
ENERGY	incident-neutron energy, EMIN<ENERGY<EMAX		REAL*4

Name	Explanation	Default value	Type
NUMLC	number of Legendre coefficients, ≤ 20, if MODE='LEG'		INTEGER*4
NEWTAB	= { 'NO' : use old cosine grid, 'READ' : read cosine grid values 'EQI' : calculate grid of equi- distant cosines	'NO'	REAL*8
NUMCOS	number ≤ 101 of cosine grid values for NEWTAB='READ' or ='EQI'		INTEGER*4
COSTAB	list of NUMCOS cosine grid values separated by commas, in descending order, for NEWTAB='READ'.		REAL*4 array(101)
LEGC	list of NUMLC Legendre coefficients separated by commas, in descending order, for MODE='LEG'		

If another material or the present material with different input parameters is to be treated a new `&IN_MAT=...` name list follows, otherwise the end of input is reached. Input of tabulated SGNC values from cards is not possible with the present version of SELPLO. Input of tabulated SGNC values via tape utilizes KEMA input format. For each energy the following quantities are read:

NUM            INTEGER\*4            number of 4-byte words following NUM in this record, i.e.

$$\text{NUM} = 10 + 2 * \text{NUMCOS}$$

where NUMCOS is the number of cosine values for which SGNC values are to be read

'ADD'	REAL*8	constant
3	INTEGER*4	constant
MAT	REAL*8	material name, e.g. 'FE__'
'SGNC'	REAL*8	constant
ENERGY	REAL*4	incident energy
1	INTEGER*4	constant
1	"	"
(X(I),Y(I), I=1,NUMCOS)	REAL*4	abscissae (cosines) and ordinates (SGNC values)

These data are read on unit ITAP and must be written with LRECL=80. The first byte of the first logical record must be occupied, all following records must begin with a blank.

Input Example:

```
&IN_MAT='FE,EMIN=1.E+4,EMAX=1.E+5,  
-PLOT=T,SIGMEL=T,MUL=T,STATIS=T, END
```

2.2.3 SELPLO output

The output is essentially self-explanatory. If  $\bar{\mu}_L^S$  calculated from SGNC and  $\bar{\mu}_L^K$  stored on KEDAK are compared, the following quantity is printed:

$$\frac{\delta \bar{\mu}_L}{1 - \bar{\mu}_L^K} = \frac{\left| \bar{\mu}_L^K - \bar{\mu}_L^S \right|}{1 - \bar{\mu}_L^K} = \left| \frac{\sigma_{tr}^S - \sigma_{tr}^K}{\sigma_{tr}^K} \right|$$

which gives the relative deviation of the elastic transport cross section

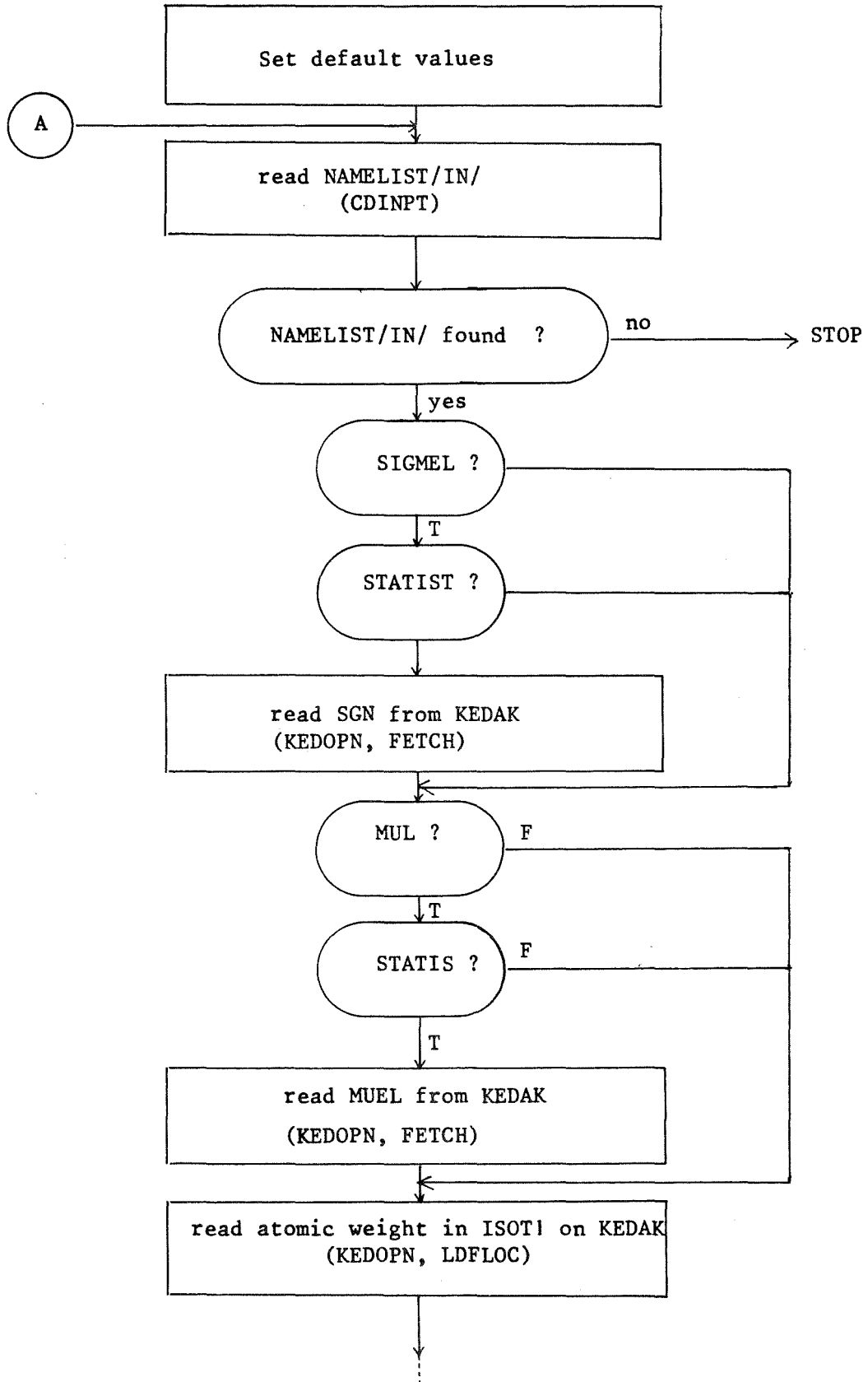
$$\sigma_{tr} = (1 - \bar{\mu}_{el}) \sigma_n$$

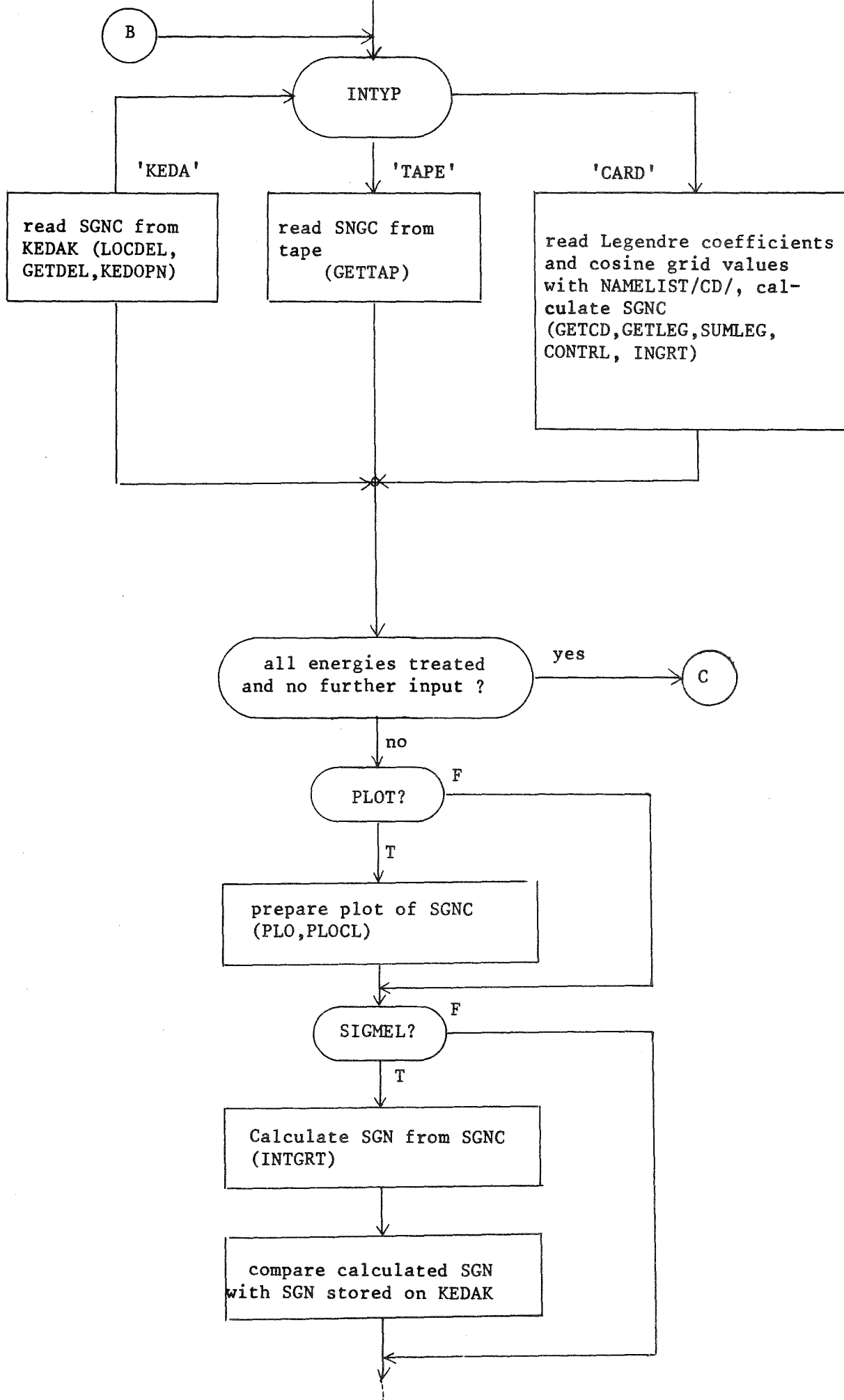
if  $\bar{\mu}_L$  is calculated from SGNC rather than retrieved from KEDAK.

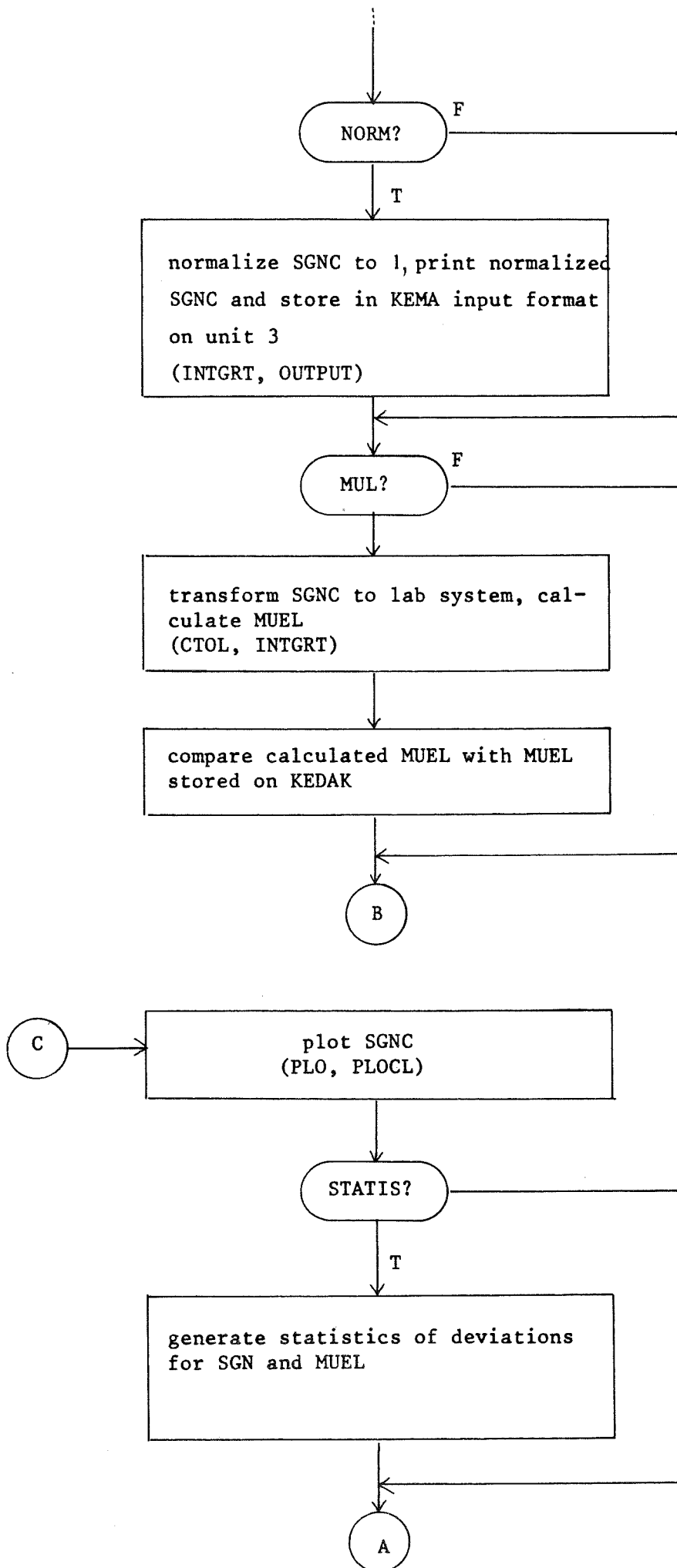
2.2.4 Short description of SELPLO main program and subroutines

Flow diagram of SELPLO main program

(T=TRUE, F=FALSE, subroutines called are shown in brackets)







SELPLO subroutines

CDINPT reads card input from unit 5 and stores it momentarily on unit 4

KEDOPN opens nuclear data file KEDAK, calls LDFOPN

FETCH reads SGN or MUEL from KEDAK, calls LDFLOC, LDFNXT

LOCDEL reads SGNC from KEDAK (ENTRY GETDEL), calls LDFLOC, LDFNXT

GETTAP reads SGNC from tape

GETCD reads SGNC or Legendre coefficients with cosine grid values from cards, calls GETLEG, GETTAB

GETLEG calculates SGNC from Legendre coefficients for all cosine grid values, checks whether cosine grid is fine enough, calls SUMLEG, CONTRL

SUMLEG calculates SGNC from Legendre coefficients for given energy

CONTRL calculates first two Legendre moments of a distribution and compares them with the generating Legendre coefficients

GETTAB dummy in present version of SELPLO

INTGRT integrates according to trapezoidal rule

OUTPUT prints SGNC, stores SGNC in KEMA input format



CTOL                    transforms from center-of-mass to lab system

ERMSG                   prints error messages

PLO                     plots SGNC (ENTRY PLOCL),  
                         calls RASTER, GRIDLI, PLOTA

RASTER                  plots a grid of horizontal and vertical lines,  
                         calls RALIN, RALOG, PLOTA

RALIN                   generates grid for linear scale

RALOG                   generates grid for logarithmic scale

GRIDLI                  generates parameters for linear axis  
                         corresponding to size of plot paper

PLOTA                   is a Karlsruhe general-purpose plot subroutine /1/

LDF                     reads the nuclear data library KEDAK /2/,  
                         (ENTRIES LDFOPN, LDFLOC, LDFNXT)

STRING,CONVX           system routines /3/

FREEFO                  conversion of input data to binary representation /4/

#### 2.2.5 Storage and time requirements on IBM/360-65 and /370-168

SELPLO uses fixed-dimension arrays. At most 2500 SGN and MUEL values can be treated in the energy range EMIN...EMAX specified in the input. The maximum number of SGNC values is 101.

The source program requires about 196 k bytes. An additional 30-50 k bytes are needed for buffers depending on number and blocking of external units. About 0.05 s of CPU time are needed per processed angular distribution.

2.2.6 External storage, job control

Unit	contents	required
1	KEDAK	if INTYP='KEDA' or SIGMEL=T or MUL=T or if atomic weight is read from KEDAK
2	normalized SGNC values in KEMA input format (for output other than printing)	if NORM=T
3	normalized SGNC values (for printing)	if NORM=T
4	intermediate storage of input read from unit 5	always
5	card input	always } { units assigned automatically, if standard procedu- res are used.
6	printed output	
7	punched output for IBM 1130 plotter	if PLOT=1
8	SGNC in binary representation (for reading by unit ITAP)	if INTYP='TAPE'
ITAP	SGNC in KEMA input format	if INTYP='TAPE'
(>8)	(input)	

Utilization of all external units (with ITAP=9) requires the following job control information:

```
//INR901SE JØB(0901,106,P6M1A),GØEL,CLASS=A,
// REGIØN=250k,TIME=1
/*FØRMAT PU,DDNAME=FTO7FOO1,FØRMS=TUSCHE
/*SETUP DEVICE=2314,ID=GFK050
/*SETUP DEVICE=2314,ID=GFK029
// EXEC FGLG
//L.PLIB DD UNIT=2314,VØL=SER=GFK029,
//_DSN=INR.STEIN.LØAD,DISP=SHR
//L.SYSIN DD *
  INCLUDE PLIB(RETPAC,LDFPAC,DEFI,SELPLØ)
  ENTRY MAIN
/*
//G.FTO1FOO1 DD UNIT=2314,VØL=SER=GFK050,
// DSN=KEDAK3,DISP=SHR
//G.FTO2FOO1 DD UNIT=SYSDA,SPACE=(TRK,20),
// DISP=(NEW,DELETE)
//G.FTO3FOO1 DD SYSØUT=A
//G.FTO4FOO1 DD UNIT=SYSDA,SPACE=(TRK,(100,10)),
//_DCB=(RECFM=FB,LRECL=80,BLKSIZE=6400),
//_DISP=(NEW,DELETE)
//G.FTO7FOO1 DD SYSØUT=T
//G.FTO8FOO1 DD UNIT=SYSDA,SPACE=(TRK,10)
//G.FTO9FOO1 DD UNIT=2314,VØL=SER=GFK016,
// DSN=SELPLØ.DATA,DISP=SHR
//G.SYSIN DD *
  &IN...(input parameters)...&END
/*
//
```

References

/1/ S. Heine, PLOTA, Programmbeschreibung  
(1967) and updates, unpublished

/2/ Present Compendium Ch. III.2

/3/ K. Gogg, CONVX, Programmbeschreibung, unpublished  
K. Gogg. STRING, Programmbeschreibung, unpublished

/4/ H. Bachmann, (1970) unpublished

Appendix: SELPLØ program listing

LOGICAL*1 SIGMEL,MUL,NORM,PLOT,STATIS,DRUCK	10	2* PLOT=',L2,', PLOFOR=',A2,	560
REAL*4 MU,INTGRL,INTGR,MUB,MUK,INTYP,KEDA/'KEDA'/	20	X', SIGMEL=',L2,', MUL=',L2,', NORM=',L2,', STATIS=',L2,	570
DIMENSION ARG(4),NARG(3),COS(101),DSIG(101),ENEL(2500),SIGEL(2500)	30	X', DRUCK=',L2/	580
1,ENMU(2500),MU(2500),TXT(20)	40	4* INTYP=',A4,', NUMKED=',I3)	590
DIMENSION XZ(1001),YZ(1001),ZZ(1001),NMU(42),NSL(22)	50	IF(NORM) WRITE(KOUT,622) ITAB	600
RFAL*8 MAT,NAME(3),SGN/'SGN' '//, MUJEL/'MUJEL'	60	IF (INTYP.EQ.TAPE) GO TO 11	610
1'/',BLANK/' '//,ISOT1/'ISOT1' '//,TTEND/'ENDE'//	70	622 FORMAT(' ITAB=',I2)	620
COMMON INPUT,KOUT,ITAP,ITAB	80	NAME(1)=MAT	630
DATA CARD/'CARD'//,TAPE/'TAPE'//,TIN/' &IN'//,TEND/'ENDE'//,I2/2/	90	NAME(3)=BLANK	640
LOGICAL FIRST//	100	GOTO 14	650
DATA TUSCHE/'T'//	110	C	660
NAMLIST/IN/ MAT,EMIN,EMAX,PLOT,SIGMEL,MUL,NORM,STATIS,INTYP,A,	120	C FEHLERZWEIG.	670
1 DRUCK,NUMKED,ITAP,PLOFOR	130	12 IF(TXT(1).EQ.TEND) GOTO 99	680
DATA ATN/1.0086654/,PI/3.14159265/	140	CALL ERMSG(TIN,TXT)	690
CALL FSPIE	150	GOTO 6	700
KEKONT=0	160	C	710
LENX=101	170	11 WRITE(KOUT,631)ITAP	720
LAXX=1001	180	631 FORMAT(' TAPE INPUT UNIT: ITAP=',I3)	730
PLOFOR=TUSCHE	190	IF (ITAP.GT.8) GO TO 13	740
MAXNUM=2500	200	WRITE(KOUT,632)	750
ITAP=8	210	632 FORMAT(' UNITS 1 - 8 ARE INTERNALLY RESERVED USE ANOTHER UNIT	760
ITAB=2	220	1 FOR TAPE INPUT')	770
INPUT=4	230	GO TO 999	780
KIN=5	240	13 CALL FREE72(ITAP,8,6,MU,MU,MU)	790
KCUT=6	250	C	900
CALL CDINPT(KIN)	260	C ITAP INPUT UNIT ANGULAR DISTRIBUTION KEAINPUT FORMAT	810
EMIN=0.	270	C 8 UNIT ON WHICH THE INPUT INFORMATION IS WRITTEN IN	820
EMAX=20.E+6	280	C BINARY FORM	830
PLOT=.FALSE.	290	C 6 PROTOCOL UNIT	840
SIGMEL=.FALSE.	300	C MU WORKING SPACE	850
MUL=.FALSE.	310	C	860
NORM=.FALSE.	320	C SGN WERDEN VON KEDAK NACH (ENEL,SIGEL) GELESEN	870
STATIS=.FALSE.	330	14 IF(.NOT.SIGMEL.OR..NOT.STATIS) GOTO 20	880
DRUCK=.FALSE.	340	NAME(2)=SGN	890
INTYP=KEDA	350	NARG(1)=2	900
	360	IF(KEKONT.EQ.0) CALL KEDOPN(KEKONT,895)	910
	370	CALL FETCH(NARG,NAME,EMIN,EMAX,ENEL,SIGEL,MAXEL,MAXNUM,&15)	920
	380	GO TO 20	930
	390	15 STATIS=.FALSE.	940
	400	C	950
	410	C MUJEL WERDEN VON KEDAK NACH (ENMU,MU) GELESEN	960
	420	20 IF(.NOT.MUL.OR..NOT.STATIS) GOTO 30	970
	430	NAME(2)=MUJEL	980
	440	NARG(1)=2	990
	450	IF(KEKONT.EQ.0) CALL KEDOPN(KEKONT,895)	1000
	460	CALL FETCH(NARG,NAME,EMIN,EMAX,ENMU,MU,MAXMU,MAXNUM,&25)	1010
	470	GOTO 30	1020
	480	25 STATIS=.FALSE.	1030
	490	C	1040
	500	C ISOT1 WIRD GEHOLT.ATOMGEWICHT WIRD EINGELESEN.	1050
	510	30 MEL=1	1060
	520	MMU=1	1070
	530	IF(A.NE.0.) GOTO 28	1080
	540	NARG(1)=2	1090
	550	NAME(1)=MAT	1100

NAME(2)=ISOT1	1110	IF(.NOT.PLOT) GOTO 38	1660
IF(KEKONT.EQ.0) CALL KEDOPN(KEKONT,895)	1120	CALL PLO(COS,DSIG,MAX,EL,MAT,3)	1670
CALL LDFLOC(NERR,NARG,NAME,ARG)	1130	NPLO=NPLO+1	1680
IF(NERR.NE.0) GO TO 27	1140	IF(NPLO.LT.4) GOTO 38	1690
WRITE(KOUT,633)NAME(1)	1150	CALL PLOCL(PLOFOR)	1700
633 FORMAT(' ***** ATOMIC WEIGHT FOR ',A8,' NOT AVAILABLE ON KEDAK E	1160	NPLO=0	1710
INTER IT VIA INPUT PARAMETER A *****')	1170	GOTO 38	1720
GO TO 999	1180	C	1730
27 A=ARG(1)	1190	C SGNC WIRD UEBER 4PI INTEGRIERT	1740
28 WRITE(6,650) A	1200	38 IF(.NOT.SIGMEL) GOTO 47	1750
650 FCRMAT(' A=',F12.6)	1210	CALL INTGRT(COS,DSIG,MAX,INTGRL)	1760
A=A/ATN	1220	SN=2.*PI*INTGRL	1770
C	1230	C	1780
C ELASTISCHE WINKELVERTEILUNGEN WERDEN VON KEDAK GEHLT (SGNC)	1240	C ZUGFUEHRIGER ELASTISCHER QUERSCHNITT WIRD GEHLT	1790
BZW VON BAND IN AUFNAHMEFORMAT FUER KEDAK.	1250	IF(.NOT.STATIS) GOTO 43	1800
WRITE(KOUT,624)	1260	39 IF(MEL.GT.MAXEL) GOTO 43	1810
624 FORMAT('//' ENERGIE IN EV,SIGMAS IN BARN//')	1270	IF(ENEL(MEL)-EL) 40,42,44	1820
IF(.NOT.STATIS.AND.(SIGMEL.OR.MUL)) WRITE(KOUT,623)	1280	40 MEL=MEL+1	1830
623 FORMAT(3X,'ENERGIE',25X,'SIGMA EL',29X,'MUE//')	1290	GOTO 39	1840
IF(STATIS) WRITE(KOUT,621)	1300	42 SNK=SIGEL(MEL)	1850
621 FORMAT(	1310	GOTO 46	1860
13X,'ENERGIE',15X,'SGN',7X,'AUS SGNC',3X,'DIFF(%)',9X,'MUEL',5X,'AU	1320	43 WRITE(KOUT,608) SN	1870
25 SGNC',4X,'DIFF',3X,'D(MU)/(1-MU)//')	1330	608 FORMAT('+',22X,'-----',2X,F10.4,3X,'----')	1880
EL=EMIN	1340	GOTO 47	1890
31 MAX=NUMKED	1350	44 SNK=SIGEL(MEL-1)+(SIGEL(MEL)-SIGEL(MEL-1))*(EL-ENEL(MEL-1))/(ENEL(	1900
IF(INTYP.EQ.KEDA) GOTO 34	1360	1MEL)-ENEL(MEL-1))	1910
IF(INTYP.EQ.TAPE) GOTO 33	1370	46 SN=SN*SNK	1920
IF(INTYP.EQ.CARD) GOTO 36	1380	DIFF=100.*A3S(SNK-SN)/SNK	1930
WRITE(KOUT,620)	1390	WRITE(KOUT,609) SNK,SN,DIFF	1940
620 FORMAT(' ALS EINGABEMEDIEN SIND DERZEIT NUR BAND,KEDAK UND KARTEN	1400	609 FORMAT('+',18X,F10.4,2X,F10.4,2X,F5.1)	1950
1 ERLAUBT')	1410	IF(DIFF.GE.1.) GOTO 41	1960
GOTO 999	1420	DIFF=DIFF+1.	1970
33 CALL GETTAP(MAT,EL,COS,DSIG,MAX,EMIN,EMAX,NUMKED,LENX,880)	1430	GOTO 45	1980
IF(FIRST.AND.NUMKED.EQ.0) NUMKED=MAX	1440	41 DIFF=DIFF/5.+2.	1990
IF(FIRST) FIRST=.FALSE.	1450	IF(DIFF.GT.21.) DIFF=22.1	2000
IF(MAX.NE.NUMKED) GOTO 94	1460	45 NSL(DIFF)=NSL(DIFF)+1	2010
GOTO 35	1470	C	2020
34 IF(FIRST) GOTO 341	1480	C NORMIEREN DER ELASTISCHEN VERTEILUNGEN AUF 1.	2030
CALL GETDEL(EL,COS,DSIG,MAX,880,880)	1490	47 IF(.NOT.NDRM) GOTO 50	2040
GOTO 342	1500	IF(SIGMEL) GOTO 49	2050
341 IF(KEKONT.EQ.0) CALL KEDOPN(KEKONT,895)	1510	CALL INTGRT(COS,DSIG,MAX,INTGRL)	2060
CALL LOCDEL(MAT,EL,COS,DSIG,MAX,880,880)	1520	SN=2.*PI*INTGRL	2070
IF(FIRST.AND.NUMKED.EQ.0) NUMKED=MAX	1530	49 FNORM=1./SN	2080
FIRST=.FALSE.	1540	DO 48 M=1,MAX	2090
GOTO 342	1550	48 DSIG(M)=FNORM*DSIG(M)	2100
342 IF(EL.GT.EMAX) GOTO 80	1560	CALL OUTPUT(MAT,EL,COS,DSIG,MAX)	2110
IF(MAX.NE.NUMKED) GOTO 94	1570	C	2120
GOTO 35	1580	C BERECHNUNG VON MUEL AUS SGNC	2130
36 CALL GETCD(COS,DSIG,MAX,EL,ZZ,LENX,DRUCK,880)	1590	50 IF(.NOT.MUL) GOTO 31	2140
IF(EL.GT.FMAX) GOTO 80	1600	CALL CTOL(COS,DSIG,MAX,XZ,YZ, LAX,LAXX,A)	2150
35 WRITE(KOUT,607) EL	1610	CALL INTGRT(XZ,YZ,LAX,INTGRL)	2160
607 FORMAT(/2X,1PE12.3)	1620	DO 52 I=1,LAX	2170
NUMELD=NUMELD+1	1630	52 ZZ(I)=XZ(I)*YZ(I)	2180
C	1640	CALL INTGRT(XZ,ZZ,LAX,INTGR)	2190
C PLOTTEN DER ELASTISCHEN VERTEILUNGEN	1650	MUB=INTGR/INTGRL	2200

C		2210			2760
C	KEDAK-MUEL HOLEN	2220	DC 84 M=1,41		
	IF(.NOT.STATIS) GOTO 59	2230	MA=M+1		2770
54	IF(MMU.GT.MAXMU) GOTO 59	2240	DO 83 J=MA,42		2780
	IF(ENMU(MMU)-EL) 56,58,60	2250	83 NMU(M)=NMU(M)+NMU(J)		2790
56	MMU=MMU+1	2260	DIFF=.05*FLOAT(M-2)		2800
	GOTO 54	2270	IF(M.EQ.1) DIFF=0.		2910
58	MUK=MU(MMU)	2280	IF(M.EQ.2) DIFF=.01		2820
	GOTO 62	2290	WRITE(KOUT,629) DIFF,NMU(M)		2830
59	WRITE(KOUT,610) MUB	2300	629 FCRMAT(' DIFF>',F4.2,I6)		2840
610	FORMAT(' '+,59X,'-----',2X,F8.4,2(4X,'-----'))	2310	84 CONTINUE		2850
	GOTO 31	2320	DIFF=2.		2860
60	MUK=MU(MMU-1)+(MU(MMU)-MU(MMU-1))*(EL-ENMU(MMU-1))/(ENMU(MMU)-ENMU(MMU-1))	2330	WRITE(KOUT,629) DIFF,NMU(42)		2870
		2340	GOTO 999		2880
62	DIFF=ABS(MUK-MUB)	2350	89 WRITE(KOUT,615)		2890
	DIFF1=DIFF/(1.-MUK)	2360	615 FORMAT('IENDE DER EINGABE ERREICHT')		2900
	WRITE(KOUT,611) MUK,MUB,DIFF,DIFF1	2370	IF(.NOT.NORM) GOTO 90		2910
611	FCRMT(' '+,57X,3(F8.4,2X),G12.4)	2380	WRITE(ITAB) I2,TTEND		2920
	IF(DIFF.GE.1.E-2) GOTO 63	2390	WRITE(3,634) I2,TTEND		2930
	DIFF=DIFF+1.01	2400	634 FORMAT(19,3X,A4)		2940
	GOTO 64	2410	ENDFILE ITAB		2950
63	DIFF=DIFF*20.+2.	2420	WRITE(3,635) ITAB		2960
	IF(DIFF.GT.41.) DIFF=42.1	2430	635 FORMAT(' END OF FILE WRITTEN ON UNIT',I3)		2970
64	NMU(DIFF)=NMU(DIFF)+1	2440	90 STOP		2980
	GOTO 31	2450	C		2990
C		2460	C FEHLERZWEIGE		3000
C	STATISTIK	2470	94 WRITE(KOUT,616) MAT,EL,NUMKED		3010
C		2480	616 FCRMAT(' FUER ',A8,' SGAC EL=',E12.4,' IST DIE ANZAHL DER SAETZE N		3020
	80 CALL PLOCL(PLOFOR)	2490	IICHT = ',I3)		3030
	FIRST=.TRUE.	2500	MAX=NUMKED		3040
	WRITE(KOUT,626) NUMELD	2510	GOTO 35		3050
626	FORMAT('I'/' ANZAHL DER BEARBEITETEN ELASTISCHEN VERTEILUNGEN = ',	2520	C		3060
	I15)	2530	95 WRITE(KOUT,617)		3070
	IF(.NOT.STATIS) GOTO 999	2540	617 FORMAT(' LDFOPN-AUFRUF BLIEB OHNE ERFOLG')		3080
	WRITE(KOUT,618)	2550	GOTO 80		3090
618	FORMAT(/2X,'VERTEILUNG DER ABWEICHUNGEN ZWISCHEN KEDAK TABFLIFRT	2560	C		3100
	1 UND KEDAK GERECHNET AUS DEN DIFFERENTIELLEN VERTEILUNGEN')	2570	END		3110
	IF(.NOT.SIGMEL) GOTO 85	2580			
	WRITE(KOUT,619)	2590			
619	FORMAT(/' FUER SGN:')	2600			
	DO 82 M=1,21	2610			
	MA=M+1	2620	SUBROUTINE FETCH(NARG,NAME,EMIN,EMAX,ARG,WERT,MAX,MAXE,*)		10
	DO 81 J=MA,22	2630	COMMON INPUT,KOUT		20
81	NSL(M)=NSL(M)+NSL(J)	2640	DIMENSION ARG(2500),WERT(2500),Z(2),NARG(3)		30
	MM=5*(M-2)	2650	REAL *8 NAME(3)		40
	IF(M.EQ.2) MM=1	2660	I=1		50
	IF(M.EQ.1) MM=0	2670	CALL LDFLOC(NERR,NARG,NAME,Z)		60
	WRITE(KOUT,627) MM,NSL(M)	2680	IF(NERR.EQ.0) GOTO 30		70
627	FORMAT(' DIFF>',I4,' %',I6)	2690	IF(Z(1).LT.EMIN) GOTO 20		80
82	CONTINUE	2700	IF(Z(1).GT.FMAX) GOTO 32		90
	MM=100	2710	ARG(I)=Z(I)		100
	WRITE(KOUT,627) MM,NSL(22)	2720	WERT(I)=Z(2)		110
85	IF(.NOT.MUL) GOTO 999	2730	I=I+1		120
	WRITE(KOUT,628)	2740	20 CALL LDFNXT(NERR,NARG,NAME,Z)		130
628	FORMAT(/' FUER MUEL:')	2750	IF(NERR.EQ.0) GOTO 22		140
			IF(Z(1).LT.FMIN) GOTO 20		150



```

IF(Z(1).GT.EMAX) GOTO 24
ARG(I)=Z(1)
WERT(I)=Z(2)
I=I+1
IF(I.GT.MAXE) GOTO 34
GOTO 20
22 WRITE(KOUT,605) (NAME(J),J=1,2)
605 FORMAT(' EMAX > OBERSTER ENERGIEWERT AUF KEDAK FUER ',2A8)
24 MAX=I-1
WRITE(KOUT,609) (NAME(J),J=1,2),MAX
RETURN
30 WRITE(KOUT,606) (NAME(I),I=1,2),EMIN,EMAX
606 FORMAT(' FUER ',2A8,' WURDEN ZWISCHEN',E12.4,' FV UND',F12.4,' FV
1AUF KEDAK KEINE DATEN GEFUNDEN')
RETURN 1
32 WRITE(KOUT,607) (NAME(I),I=1,2),EMAX,EMIN
607 FORMAT(' FUER ',2A8,' IST ANGEGEBENES EMAX=',E12.4,' FV KLEINER ALS
11. KEDAKENERGIE ODER ANGEGEBENES FMIN=',E12.4,' EV')
RETURN 1
34 WRITE(KOUT,608) (NAME(J),J=1,2),MAXE,EMIN,ARG(I-1)
608 FORMAT(' ANZAHL DER ENERGIEN AUF KEDAK FUER ',2A8,' GROESSER ALS '
1,15 '/' ES WIRD DAS EINLESEN VON KEDAKWERTEN BEI DIESER ZAHL ABGEB
2ROCHEN.'/' EINGELESEN WURDEN ALLE WERTE ZWISCHEN',F12.4,' EV UND ',
3E12.4,' EV')
MAX=I-1
WRITE(KOUT,609) (NAME(J),J=1,2),MAX
609 FORMAT(' ANZAHL DER VON KEDAK GEHOLTEN DATENPUNKTE FUER ',2A8,' =
1',I5)
RETURN
END

SUBROUTINE PLO(X,Y,MAX,EL,MAT,NTI)
COMMON INPUT,KOUT
REAL*8 NSKB(11)/'-1.0.. ', '-.8.. ', '-0.6.. ', '-0.4.. ', '-0.2.
1. ', ' 0.0.. ', '+0.2.. ', '+0.4.. ', '+0.6.. ', '+0.8.. ', '+1.0.
2. ', ' SGNC/SGNC ', ' MAT
DIMENSION XBS(11),YBS(11),NDIRSB(10),NSCSB(10),XBSDUM(2),YBSDUM(2)
DIMENSION X(2),Y(2)
DIMENSION XZ(404),YZ(404)
DIMENSION NTEXT(15),XB(2),YB(2),NDIR(2),NSC(2),XT(2),YT(2),NTE(10)
DIMENSION XR(200),YR(200)
DIMENSION MAXN(10),ELA(10),NT(10),NP(10)
LOGICAL XLOG
DATA NUM/9/,ID/0/,YMAX,YMIN/0.,-2./,
INLGY/+1/,YA,YE/-1.,+1./,NFX/'F9.2'/,NFY/'F5.1'/,NSCX,NSCY/1,1/,
2 NUOX,NUOY/+1,+1/,NPSX,NPSY/-1,+1/,NX,NY/1,1/,NT/3/,
3NPG/1/,IND/1/,NPA/1/,DX,DY/1.,.2/,ELA/0./,NTE(1)/'EN= '/,
4 MN/0/,NP/0,1,2,3,4,5,6,7,8,9/,SY/2.54E-3/,NDIR/3,3/,NSC/1,1/
DATA NSCSB/10*1/,NDIRSB/10*3/,NTSB/10/,NTSBI/1/
DATA TUSCHE/'T'/
DATA DYY/.2/
C EINSPEICHERN DER ZU PLOTTENDEN DATEN NACH XZ,YZ

```

```

160 IF(MN+MAX.GT.404) GOTO 89
170 NUM=NUM+1
180 DO 101 J=1,MAX
190 MN=MN+1
200 XZ(MN)=Y(J)
210 101 YZ(MN)=X(J)
220 ELA(NUM)=EL
230 MAXN(NUM)=MN
240 NT(NUM)=NTI
250 GOTO 99
C
C ES WERDEN ALLE SEIT DEM LETZTEN PLOCL-AUFRUF DURCH PLO ANGELIEFERTEN
C KURVEN GEZEICHNET.
C
ENTRY PLOCL(PLOFOR)
IF(NUM.LT.1) GOTO 99
SY=2.54E-3
IF(PLOFOR.NE.TUSCHE) SY=2.5E-3
C
C TITEL UND NUMMER DER ZEICHNUNG
CALL STRING(NTEXT(1),' ')
CALL STRING(NTEXT(2),NTEXT(1),56)
CALL STRING(NTEXT(1),MAT,8)
CALL STRING(NTEXT(3),SGNC,9)
ID=ID+1
C
C PLOTGRENZEN UND PLOTART FESTLEGEN.
XMAX=XZ(1)
XMIN=XZ(1)
DO 3 M=2,MN
IF(XZ(M).GT.XMAX) GOTO 2
IF(XZ(M).GE.XMIN) GOTO 3
XMIN=XZ(M)
GOTO 3
2 XMAX=XZ(M)
3 CONTINUE
IF(XMIN.NE.0) GOTO 4
XLOG=.FALSE.
GOTO 6
4 IF(XMAX.GT.20.*XMIN) GOTO 5
XLOG=.FALSE.
GOTO 6
5 XLOG=.TRUE.
6 IF(XLOG) GOTO 7
GOTO 30
C
C ERSTER PLOT FUER LOGARITHMISCHE SIGMA-ACHSE.
7 XMAX=ALOG10(XMAX)
XMIN=ALOG10(XMIN)
XRES=.05*FLOAT(NUM+2)
DO 9 M=1,MN
YZ(M)=YZ(M)-1.
9 XZ(M)=ALOG10(XZ(M))
MA=INT(XMAX)
IF(FLOAT(MA).LT.XMAX) MA=MA+1

```

```

M=INT(XMAX+XRES)
IF(FLOAT(M).LT.XMAX+XRES) M=M+1
MI=INT(XMIN)
IF(FLOAT(MI).GT.XMIN) MI=MI-1
NRES=INT(XMIN-XRES)
IF(FLOAT(NRES).GT.XMIN-XRES) NRES=NRES+1
IF(M.EQ.MA) GOTO 11
IF(NRES.NE.MI) GOTO 10
NRES=1
GOTO 12
10 MA=M
11 NRES=2
GOTO 12
12 XMIN=FLOAT(MI)
XMAX=FLOAT(MA)
INDZ=INT((XMAX-XMIN)/2.5)+1
XA=10.**MI
XE=10.**MA
XALOG=XMIN
XELOG=XMAX
SX=2.54E-3
NLGX=-1
CALL PLOTA(XZ,YZ,MAXN(1),NT(1),NP(1),NPG,IND,NPA,INDZ,XMAX,XMIN,
1 SX,YMAX,YMIN,SY,NTEXT,1D,
2 NLGX,XA,DX,XE,NFX,NSCX,NUOX,NPSX,NX,
3 NLGY,YA,DY,YE,NFY,NSCY,NUOY,NPSY,NY,0)
INDZ=0
CALL RASTER(-1,+1,XMAX,XMIN,SX,YMAX,YMIN,SY,INDZ,XR,YR,DX,DY)
GOTO 20
C
C ERSTER PLOT IM FALL LINEARER SIGMA-ACHSE
30 CALL GRIDLI(XMIN,XMAX,RMIN,RMAX,SX)
IF(PLOFOR.NE.TUSCHE) SX=SX*2.5/2.54
DX=20.*SX/.254
IF(PLOFOR.NE.TUSCHE) DX=80.*SX
XRES=FLOAT(NUM+2)*DX/4.
NRES=INT(XRES/DX)+1
XRES=FLOAT(NRES)*DX
XMAX=RMAX+XRES
NRES=2
XMIN=RMIN
INDZ=INT((XMAX-XMIN)/800./SX)+1
NLGX=+1
XA=XMIN
XE=XMAX
DO 32 M=1,MN
32 YZ(M)=YZ(M)-1.
CALL PLOTA(XZ,YZ,MAXN(1),NT(1),NP(1),NPG,IND,NPA,INDZ,XMAX,XMIN,
1 SX,YMAX,YMIN,SY,NTEXT,1D,
2 NLGX,XA,DX,XE,NFX,NSCX,NUOX,NPSX,NX,
3 NLGY,YA,DY,YE,NFY,NSCY,NUOY,NPSY,NY,0)
INDZ=0
CALL RASTER(1,1,XMAX,XMIN,SX,YMAX,YMIN,SY,INDZ,XR,YR,DX,DY)
GOTO 20
C

```

```

770 C BESCHRIFTUNG DER MUE-ACHSE 1320
780 20 XX=XMIN-24.*SX 1330
790 YBS(1)=YMIN+16.*SY 1340
800 DO 21 I=1,11 1350
810 21 XBS(I)=XX 1360
820 DC 22 I=2,11 1370
830 22 YBS(I)=YBS(I-1)+DYY 1380
840 XBSDUM(1)=XMIN-1.E+3*SX 1390
850 XBSDUM(2)=XBSDUM(1) 1400
860 YBSDUM(1)=YMIN-1.E+3*SY 1410
870 YBSDUM(2)=YBSDUM(1) 1420
880 CALL PLOTA(XBSDUM,YBSDUM,2,2,1,1,IND,1,INDZ,XMAX,XMIN,SX,YMAX,YMIN 1430
890 1,SY,1,1D,0,0, 1440
900 2NLSB,XBS,YBS,NDIRSB,NSCSB,NSKB(1),NSKB(2),NSKB(3),NSKB(4),NSKB(5), 1450
910 3NSKB(6),NSKB(7),NSKB(8),NSKB(9),NSKB(10)) 1460
920 CALL PLOTA(XBSDUM,YBSDUM,2,1,1,1,IND,1,INDZ,XMAX,XMIN,SX,YMAX,YMIN 1470
930 1,SY,1,1D,0,0, 1480
940 2NLSB1,XBS(11),YBS(11),NDIRSB,NSCSB,NSKB(11)) 1490
950 C 1500
960 C ZEICHENERKLAERUNG FUER DIE ERSTE KURVE 1510
970 IF(NRES.EQ.1) XT(1)=XMIN+XRES 1520
980 IF(NRES.EQ.2) XT(1)=XMAX-10.*SX 1530
990 XT(2)=XT(1) 1540
1000 YT(1)=YMAX-60.*SY 1550
1010 YT(2)=YT(1) 1560
1020 XB(1)=XT(1)-8.*SX 1570
1030 YB(1)=YT(1)-25.*SY 1580
1040 CALL STRING(NTE(2),' @') 1590
1050 CALL STRING(NTE(3),NTE(2),32) 1600
1060 CALL STRING(NTE(5),' EV@') 1610
1070 CALL STRING(NTE(6),' @') 1620
1080 CALL CONVX(ELA(1),NTE(2),'E12.5') 1630
1090 CALL PLOTA(XT,YT,2,1,NP(1),NPG,IND,NPA,INDZ,XMAX,XMIN,SX, 1640
1100 1 YMAX,YMIN,SY,1,1D,0,0,1, 1650
1110 2 XB,YB,NDIR,NSC,NTE) 1660
1120 C 1670
1130 C JETZT WERDEN DIE UEBRIGEN KURVEN GEPLOTTET ZUSAMMEN MIT 1680
1140 C ZEICHENERKLAERUNG. 1690
1150 IF(NUM.LT.2) GOTO 90 1700
1160 DC 25 N=2,NUM 1710
1170 M=MAXN(N-1)+1 1720
1180 LMAX=MAXN(N)-MAXN(N-1) 1730
1190 CALL PLOTA(XZ(M),YZ(M),LMAX,NT(N),NP(N),NPG,INC,NPA,INDZ, 1740
1200 1 XMAX,XMIN,SX,YMAX,YMIN,SY,1,1D,0,0,0) 1750
1210 XT(1)=XT(1)-20.*SX 1760
1220 XT(2)=XT(1) 1770
1230 XB(1)=XT(1)-8.*SX 1780
1240 YB(1)=YT(1)-25.*SY 1790
1250 NDIR(1)=3 1800
1260 NSC(1)=1 1810
1270 CALL STRING(NTE(2),' @') 1820
1280 CALL STRING(NTE(3),NTE(2),32) 1830
1290 CALL STRING(NTE(5),' EV@') 1840
1300 CALL STRING(NTE(6),' @') 1850
1310 CALL CONVX(ELA(N),NTE(2),'E12.5') 1860

```

CALL PLOT4(XT,YT,2,1,NP(N),NPG,IND,NPA,INDZ,XMAX,XMIN,SX,	1870	SX=.2*EH*.254	390
1YMAX,YMIN,SY,1,ID,0,0,1,XB,YB,NDIR,NSC,NTF)	1880	GOTO 20	400
25 CONTINUE	1990	16 IF(LSKAL.LE.10) GOTO 18	410
GOTO 90	1900	ISTART=(IMIN/2)*2	420
89 WRITE(KOUT,689)	1910	IEND=(IMAX/2)*2	430
689 FORMAT(/' DIE ANGELIEFERTEN DATEN FINDEN IM PLOTFFLD KEINEN PLATZ	1920	IF(IEND.LT.IMAX) IEND=IEND+2	440
1'/)	1930	SX=.1*EH*.254	450
GOTO 99	1940	GOTO 20	460
90 MN=0	1950	18 IMIN=IMIN*10	470
NUM=0	1960	IMAX=IMAX*10	480
GOTO 99	1970	EH=EH/10.	490
99 RETURN	1980	LSKAL=IMAX-IMIN	500
END	1990	GOTO 10	510
		20 RMIN=FLOAT(ISTART)*EH	520
		RMAX=FLOAT(IEND)*EH	530
		RETURN	540
		END	550
SUBROUTINE GRIDLI(XMIN,XMAX,RMIN,RMAX,SX)	10		
C	20		
C GRIDLI BESTIMMT DIE ACHSEN-KENNGROESSEN FUER LINEARE ACHSEN SO,	30		
C DASS DEREN LAENGE 10-20CM BETRAEGT.	40		
XLX=ALOG10(XMAX)	50	SUBROUTINE INTGRT(X,F,MAX,S)	10
LX=INT(XLX)-1	60	DIMENSION X(2),F(2)	20
IF(XLX.LT.0.) LX=LX-1	70	S=0.	30
EH=10.**LX	80	DC 10 I=2,MAX	40
IMAX=INT(XMAX/EH)	90	10 S=S+(F(I)+F(I-1))*(X(I)-X(I-1))	50
IF(XMAX.GT.(FLOAT(IMAX)*EH)) IMAX=IMAX+1	100	S=S*.5	60
IMIN=INT(XMIN/EH)	110	RETURN	70
2 LSKAL=IMAX-IMIN	120	END	80
IF(LSKAL.LE.100) GOTO 4	130		
IMIN=IMIN/10	140	SUBROUTINE CTOL(COSC,DSIGC,MAX,COSL,DSIGL, LMAX,LLMAX,A)	10
IM=IMAX/10	150	DIMENSION COSC(1),DSIGC(1),COSL(1),DSIGL(1)	20
IF((10*IM).LT.IMAX) IM=IM+1	160	A22=A**2	30
IMAX=IM	170	A2=2.*A	40
EH=EH*10.	180	IF(A.LT.2.AND.A.GT.1.) GOTO 80	50
GOTO 2	190	MA=1	60
4 IF(LSKAL.NE.0) GOTO 10	200	L=1	70
IMAX=IMAX+50	210	MMAX=MAX-1	80
IMIN=IMIN-50	220	LC=(LLMAX-1)/MMAX	90
10 IF(LSKAL.LE.50) GOTO 12	230	DO 12 M=MA,MMAX	100
ISTART=(IMIN/10)*10	240	COSL(L)=COSC(M)	110
IEND=(IMAX/10)*10	250	DSIGL(L)=DSIGC(M)	120
IF(IEND.LT.IMAX) IEND=IEND+10	260	DC=(COSC(M+1)-COSC(M))/FLOAT(LD)	130
SX=.5*EH*.254	270	DS=(DSIGC(M+1)-DSIGC(M))/FLOAT(LD)	140
GOTO 20	280	LA=L+1	150
12 IF(LSKAL.LE.40) GOTO 14	290	LE=L+LD-1	160
ISTART=(IMIN/5)*5	300	L=L+LD	170
IEND=(IMAX/5)*5	310	DO 10 LL=LA,LE	180
IF(IEND.LT.IMAX) IEND=IEND+5	320	COSL(LL)=COSL(LL-1)+DC	190
SX=.25*EH*.254	330	DSIGL(LL)=DSIGL(LL-1)+DS	200
GOTO 20	340	10 CONTINUE	210
14 IF(LSKAL.LE.20) GOTO 16	350	14 LMAX=LE+1	220
ISTART=(IMIN/4)*4	360		
IEND=(IMAX/4)*4	370		
IF(IEND.LT.IMAX) IEND=IEND+4	380		

```

COSL(LMAX)=COSC(MAX)
DSIGL(LMAX)=DSIGC(MAX)
IF(A.LE.1.) GOTO 30
W=A-1.
DC=-1.
DSTGL(1)=W*DSIGL(1)/(A*(1.-DC/W))
CCSL(1)=DC
LA=2
19 DO 20 L=LA,LMAX
DC=COSL(L)
W=SQRT(A22+A2*DC+1.)
DC=(A*DC+1.)/W
DSIGL(L)=W*DSIGL(L)/(A*(1.-DC/W))
20 CCSL(L)=DC
GOTO 100
30 IF(A.LT.1.) GOTO 40
CCSL(1)=0.
DSIGL(1)=0.
LA=2
GOTO 19
C
C      TO EACH MU(L) THERE EXIST TWO MU(C).
C      PROCEEDING: CALCULATE MU(L) TO GIVEN MU(C).FIND SECOND MU(C)
C      CORRESPONDING TO THIS MU(L).SUM CONTRIBUTIONS FROM BOTH.
40 DC 42 L=2,LMAX
IF(COSL(L).LE.-A) GOTO 42
LA=L
GOTO 43
42 CONTINUE
43 DS=COSL(LA)
DC=COSL(LA-1)
C
C      CALCULATE DSIGMA IN L-SYSTEM FOR LOWEST INTERVAL OF MU(L)
DC=(-.5*((DSIGL(LA)-DSIGL(LA-1))/(DS-DC)*(A+DS))+DSIGL(LA))*(A+DS)
W=SQRT(A22+A2*DS+1.)
DS1=(A*DS+1.)/W
W1=SQRT(A22+DS1**2-1.)
W=(DS1**2-DS1*W1-1.)/A
DC 45 L=1,LMAX
IF(COSL(L).LE.W) GOTO 45
LE=L
GOTO 46
45 CONTINUE
46 DS=COSL(LE)
DS2=COSL(LE-1)
DC=DC+(0.5*((DSIGL(LE)-DSIGL(LE-1))/(DS-DS2)*(W-DS))+DSIGL(LE))*(D
1S-W)
DS2=SQRT(1.-A22)
CCSL(LA)=DS1
DS=DC/(DS1-DS2)
DC=DS2
C
C      DO THE OTHER INTERVALS
LE=LE-1
DSIGL(LA)=(DSIGL(LA))*(2.*DS1+W1+DS1**2/W1)-DSIGL(LE)*(2.*DS1-W1-DS

```

```

230 11**2/W1))/A
240 LA=LA+1
250 IF(LA.GT.LMAX) GOTO 70
260 DC 60 L=LA,LMAX
270 DS1=COSL(L)
280 W=SQRT(A22+A2*DS1+1.)
290 DS1=(1.+A*DS1)/W
300 W1=SQRT(A22+DS1**2-1.)
310 W=(DS1**2-DS1*W1-1.)/A
320 51 IF(COSL(LE).LE.W) GOTO 52
330 LE=LE-1
340 GOTO 51
350 52 W=DS1**2/W1
360 DS2=2.*DS1
370 DSIGL(L)=(DSIGL(L)*(DS2+W1+W)-DSIGL(LE)*(DS2-W1-W))/A
380 CCSL(L)=DS1
390 60 CCNTINUE
400 70 LA=LA-1
410 DSIGL(LA)=DS
420 COSL(LA)=DC
430 DO 72 L=LA,LMAX
440 J=L+1-LA
450 DSIGL(J)=DSIGL(L)
460 72 CCSL(J)=COSL(L)
470 LMAX=LMAX+1-LA
480 GOTO 100
490 80 W=1./A
500 LE=LLMAX/4
510 DC=(1-W)/(LE-1)
520 DO 81 L=1,LE
530 81 COSL(L)=(L-1)*DC-1.
540 LA=2
550 DO 83 L=1,LE
560 82 IF(COSL(L).LT.COSC(LA)) GOTO 83
570 LA=LA+1
580 GOTO 82
590 83 DSIGL(L)=(DSIGC(LA)-DSIGC(LA-1))/(COSC(LA)-COSC(LA-1))*(COSC(LA)-C
ZOSL(L))+DSIGC(LA)
600 DS2=DSIGC(LA)
610 DC2=COSC(LA)
620 DS1=DSIGC(LA-1)
630 DC1=COSC(LA-1)
640 W1=W+DC2
650 DC=0.001
660 85 MA=W1/DC
670 IF(MA.LE.LLMAX/10) GOTO 86
680 DC=2.*DC
690 GOTO 85
700 86 L=LE+1
710 DS=COSL(LE)
720 COSL(L)=DS+(L-LE)*DC
730 IF(COSL(L).GE.0.999*DC2) GOTO 89
740 DSIGL(L)=(DS2-DS1)/(DC2-DC1)*(DC2-COSL(L))+DS2
750 L=L+1
760 GOTO 87
770
900 880
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

```

```

89 MA=LA
   MMAX=MAX-1
   IF(MAX.EQ.MA) GOTO 14
   LD=(LLMAX-L)/(MAX-MA)
   GOTO 8
100 RETURN
   END

SUBROUTINE OUTPUT(MAT,EL,X,Y,MAX)
COMMON INPUT,KOUT,ITAP,ITAB
DIMENSION X(21),Y(21),NM(3)
REAL*8 ADD/'ADD      '/,NAME(3)
REAL*8 SGNC/'SGNC    '/,MAT
DATA NM/3,1,1/
NUM=10+2*MAX
NALP=2
NNAM=NM(1)
NARG=NM(2)
NWERT=NM(3)
NAME(1)=MAT
NAME(2)=SGNC
NAME(3)=DBLE(EL)
WRITE(3,600) ITAB,NUM,ADD,NNAM,(NAME(I),I=1,NALP),EL,NARG,
INWERT,(X(I),Y(I),I=1,MAX)
600 FORMAT(' OUTPUT WRITTEN ON UNIT',I3,/,
1      I3,1X,A8,I2,1X,2A8,E12.4,1X,I2,I2/(7(F8.4)))
WRITE(ITAB) NUM,ADD,NNAM,(NAME(I),I=1,NALP),EL,NARG,NWERT,
1(X(I),Y(I),I=1,MAX)
RETURN
END

SUBROUTINE LOCDEL(MAT,ENERGY,X,Y,MAX,*,*)
REAL*8 NAME(3),          SGNC/'SGNC      '/, ENGOLD, MAT
DIMENSION NARG(3),ARG(2),X(MAX),Y(MAX)
COMMON INPUT,KOUT
NARG(1)=3
NAME(1)=MAT
NAME(2)=SGNC
NAME(3)=DBLE(ENERGY)
CALL LDFLOC(NERR,NARG,NAME,ARG)
CALL LDFLOC(NERR,NARG,NAME,ARG)
IF(NERR.EQ.0) GOTO 91
8 X(1)=ARG(1)
  Y(1)=ARG(2)
  J=2
  ENGOLD=NAME(3)
  ENERGY=SNGL(ENGOLD)
9 CALL LDFNXT(NERR,NARG,NAME,ARG)
  IF(NERR.EQ.0) GOTO 20

```

```

1330 X(J)=ARG(1)
1340 Y(J)=ARG(2)
1350 J=J+1
1360 GOTO 9
1370 ENTRY GETDEL(ENERGY,X,Y,MAX,*,*)
1380 CALL LDFLOC(NERR,NARG,NAME,ARG)
1390 IF(NERR.EQ.0) GOTO 92
   IF(ENGOLD.NE.NAME(3)) GOTO 8
   WRITE(KOUT,610) MAT,NAME(3)
610 FORMAT(' LETZTER SGNC-SATZ AUF KEDAK FUER ',A8,' IST',IPE14.5)
   RETURN 1
20 MK=J-1
   IF(MAX.EQ.0) GOTO 22
   IF(J-1.EQ.MAX) GOTO 22
   WRITE(KOUT,620) NAME,MAX,MK
620 FORMAT(' ANZAHL DER ERWARTETEN SAETZE FUER ',2A8,D12.4/
1' = ',I3,' GEFUNDEN WURDEN ',I3)
22 MAX=MK
   RETURN
91 WRITE(KOUT,691) NAME
691 FORMAT(' LDFLOC-AUFRUF BLIEB ERFOLGLOS FUER ',2A8,D12.4)
   RETURN 2
92 WRITE(KOUT,692)
692 FORMAT(' PROGRAMMIERER RUFEN ')
   RETURN 2
   END

SUBROUTINE GETTAP(MAT,ENERGY,X,Y,MJ,EUNT,EOB,NUMKFD,LENX,*)
COMMON INPUT,KOUT,ITAP,ITAB
DIMENSION X(2),Y(2)
REAL*8 NAME(3),MAT,SGNC/'SGNC      '/,OPT,ADD/'ACC      '/
REAL*8 MATOLD/'      '/
LOGICAL PRINT/F/,RD/F/
DATA NAME/'      ','      ','      '/
DATA TEND/'ENDE'/
DIMENSION W(212),M(1),NARG(3)
EQUIVALENCE(W(1),M(1))
C WRITE(6,700) MAT,EUNT,EOB
C 700 FORMAT('/ GETTAP MAT=',A8,' EUNT=',E13.6,' EOB=',E13.6)
M1=2*NUMKFD+10
IF(MATOLD.EQ.MAT) GOTO 2
RD=.FALSE.
MATOLD=MAT
IF(NAME(1).NE.MAT) GOTO 2
IF(NAME(2).NE.SGNC) GOTO 2
GOTO 4
2 READ(8,END=10) NUM,(W(I),I=1,NUM)
C WRITE(6,701) (W(I),I=1,11)
C 701 FORMAT(' NEUES RECORD GELESEN ',11(Z8,1X))
IF(NUMKFD.EQ.0) GOTO 3
IF(NUM.NE.M1) GOTO 13
3 M2=(NUM-10)/2

```

```

IF(M2.GT.LENX) GOTO 19
NARG(1)=M(3)
J=M(3)+6
NARG(2)=M(J)
NARG(3)=M(J+1)
CALL STRING(NAME(1),W(4),16)
4 CALL STRING(ENERGY,W(8),4)
CALL STRING(OPT,W(1),8)
NAME(3)=DBLE(ENERGY)
IF(OPT.NE.ADD) GOTO 12
IF(NAME(1).NE.MAT) GOTO 11
IF(NAME(2).NE.SGNC) GOTO 11
RD=.TRUE.
IF(ENERGY.LT.EUNT) GOTO 2
IF(ENERGY.GT.EOB) GOTO 17
J=M(3)+8
MJ=0
DO 6 I=J,NUM,2
MJ=MJ+1
X(MJ)=W(I)
6 Y(MJ)=W(I+1)
IF(PRINT) WRITE(KOUT,600) NARG,NAME,(X(I),Y(I),I=1,MJ)
600 FORMAT(' NARG=',I2,' NAMES=',2A8,D12.4/(5(F10.2,F12.4)))
RETURN
10 WRITE(KOUT,604) ITAP
604 FORMAT(' ENDE DER DATEN AUF FT',I2,' (8) ERREICHT')
14 RETURN 1
11 IF(.NOT.RD) GOTO 2
WRITE(KOUT,601) NAME
601 FORMAT(' NAMEN STIMMEN MIT DEN VORGEgebenEN NICHT UEBEREIN.'/
1' SIE LAUTEN:',2A8,D12.4)
GOTO 14
12 WRITE(KOUT,602) OPT
602 FORMAT(' OPTION=',A8,' NICHT GLEICH ADD')
GOTO 10
13 IF(NUM.EQ.2) GOTO 15
16 WRITE(KOUT,603) NUM,M1
603 FORMAT(' ANZAHL DER WORTE AUF DEM BAND =',I3,' UNGLEICH ',I3)
GOTO 10
15 IF(W(1).NE.TEND) GOTO 16
WRITE(KOUT,605) ITAP
605 FORMAT(' 1 ENDE-RECORD AUF BAND',I2,' ERREICHT')
GOTO 14
17 WRITE(KOUT,606) ITAP,EUNT,EOB
606 FORMAT(' ENDE DER DATEN AUF FT',I2,' ZWISCHEN EMIN=',F13.6,'EV
IUND EMAX=',F13.6,'EV')
GOTO 14
19 WRITE(KOUT,607) NUM,LENX
607 FORMAT(' ANZAHL DER DATENPAARE IM SATZ =',I3,' UEBERSTIEGT DATEN
FELDLAENGE =',I3)
GOTO 14
END

```

```

260 SUBROUTINE RASTER(KREX,KREY,XMAX,XMIN,SX,YMAX,YMIN,SY,INDZ,X,Y,DX, 10
270 I,DY) 20
280 DIMENSION X(2),Y(2) 30
290 DATA NRAST/9999/ 40
300 IF(INDZ.NE.0.AND.INDZ.NE.10) GOTO 50 50
310 IF(KREY.EQ.-1) GOTO 10 60
320 CALL RALIN(YMIN,YMAX,DY,XMIN,XMAX,M,X,Y) 70
330 GOTO 20 80
340 10 CALL RALOG(YMIN,YMAX,SY,XMIN,XMAX,M,X,Y) 90
350 20 CALL PLOTA(X,Y,M,2,0,1,1,1,INDZ,XMAX,XMIN,SX,YMAX,YMIN,SY,1, 100
360 INRAST,0,0,0) 110
370 IF(KREX.EQ.-1) GOTO 30 120
380 CALL RALIN(XMIN,XMAX,DX,YMIN,YMAX,M,Y,X) 130
390 GOTO 40 140
400 30 CALL RALOG(XMIN,XMAX,SX,YMIN,YMAX,M,Y,X) 150
410 40 CALL PLOTA(X,Y,M,2,0,1,1,1,INDZ,XMAX,XMIN,SX,YMAX,YMIN,SY,1, 160
420 INRAST,0,0,0) 170
430 50 RETURN 180
440 END 190

```

```

520 C
530 C DIESE VERSION ERZEUGT GITTERLINIEN IM ABSTAND DY
540 C
550 YL=YMAX-YMIN
560 NLIN=INT(YL/DY)
570 IF((YL-FLOAT(NLIN)*DY).GT.DY/2.) NLIN=NLIN+1
580 NLIN=NLIN+1
590 NPOINT=2*NLIN
600 IF(NPOINT.GT.200) NPOINT=200
610 Y1=YMIN
620 Y2=YMIN+DY
630 DO 10 I=1,NPOINT,4
640 X(I)=XMIN
650 X(I+1)=XMAX
660 X(I+2)=XMAX
670 X(I+3)=XMIN
680 Y(I)=Y1
690 Y(I+1)=Y1
700 Y(I+2)=Y2
710 Y(I+3)=Y2
720 Y1=Y2+DY
730 Y2=Y1+DY
740 10 CONTINUE
750 RETURN
760 END
770

```

```

SUBROUTINE RALOG(YMIN,YMAX,SY,XMIN,XMAX,NPCINT,X,Y)
DIMENSION X(2),Y(2),DYL(9)
DATA DYL / .0458, .3010, .1761, .1250, .0969, .0792, .0669, .0580, .0511 /
YL=YMAX-YMIN
NDEK=INT(YL)
IF((YL-FLOAT(NDEK)).GT.0.8) NDEK=NDEK+1
IF(NDEK.GT.6) GOTO 25
NLIN=9*NDEK+1
NPCINT=2*NLIN
J=0
Y1=YMIN
Y2=YMIN+DYL(2)
DO 20 I=1,NPCINT,4
X(I)=XMIN
X(I+1)=XMAX
X(I+2)=XMAX
X(I+3)=XMIN
Y(I)=Y1
Y(I+1)=Y1
Y(I+2)=Y2
Y(I+3)=Y2
J=J+2
Y1=Y2+DYL(MOD(J,9)+1)
Y2=Y1+DYL(MOD(J+1,9)+1)
20 CONTINUE
25 RETURN
END

```

```

10 GOTO 999
20 300 IF(TXT(1).EQ.TIN.OR.TXT(1).EQ.TEND) GOTO 310
30 CALL ERMSG(TCD,TXT)
40 GOTO 10
50 310 BACKSPACE INPUT
60 RETURN 1
70 999 RETURN
80 501 FORMAT(20A4)
90 600 FORMAT(/' ***** MODE CAN ONLY BE TAB OR LEG'/)
100 601 FORMAT(/' ENERGY=',E14.6,' MODE=',A8,' NUMLC=',I6,' NEWTA
110 18=',A8,' NUMCOS=',I6/)
120 END
130
140
150
160
170
180
190
200
210
220
230
240
250
260
270

```

```

SUBROUTINE GETTAB
COMMON INPUT,KOUT
WRITE(KOUT,600)
600 FORMAT(/' ***** NO TABULATED INPUT VIA CARDS ALLOWED'//)
RETURN
END

```

```

SUBROUTINE GETCD(X,Y,NX,EL,ZZ,LENX,DRUCK,*)
LOGICAL*1 DRUCK
DIMENSION X(1),Y(1),ZZ(1)
DIMENSION TXT(20),COSTAB(101)
COMMON INPUT,KOUT
REAL*8 MODE/'LEG'/,LEG/'LEG'/,TAB/'TAB'/,NEWTAB/'NO'/,FMT(10)
REAL*4 LEGCOF(20)
NAMELIST/CD/ MODE,ENERGY,NUMLC,NEWTAB,NUMCOS,
1 LEGCOF,COSTAB
DATA FMT/'(8F10.2)',9*' '/,TCD/' &CD'/,TIN/' &IN'/,TEND/'ENDE'/'
10 READ(INPUT,501) TXT
IF(TXT(1).NE.TCD) GOTO 300
BACKSPACE INPUT
READ(INPUT,CD)
WRITE(KOUT,601) ENRGY,MODE,NUMLC,NEWTAB,NUMCOS
EL=ENERGY
IF(MODE.EQ.LEG) GOTO 100
IF(MODE.EQ.TAB) GOTO 200
WRITE(KOUT,600)
GOTO 999
100 CALL GETLEG(X,Y,NX,NUMCOS,NUMLC,NEWTAB,ZZ,LENX,DRUCK,LEGCOF,
1 COSTAB)
GOTO 999
200 CALL GETTAB

```

```

SUBROUTINE GETLEG(X,Y,NX,LTAB,ND,NEWTAB,ZZ,LENX,DRUCK,CL,CT)
C
C GETLEG CONVERTS LEGENDRE COEFFICIENT INTO TABULATED DATA
C
DIMENSION X(1),Y(1),ZZ(1),CL(20),CT(1)
LOGICAL*1 DRUCK
REAL*8 NEWTAB,EQUI/'EQUI'/,RED/'READ'/,NO/'NO'/,FMT(1)
COMMON INPUT,KOUT
MAXCL=20
IF(ND.GT.MAXCL) GOTO 48
9 IF(NEWTAB.EQ.NO) GOTO 50
IF(LTAB.GT.LENX) GOTO 49
10 IF(NEWTAB.EQ.EQUI) GOTO 30
IF(NEWTAB.EQ.RED) GOTO 20
GOTO 47
20 DO 21 I=1,LTAB
21 X(I)=CT(I)
WRITE(KOUT,500) (X(I),I=1,LTAB)
500 FORMAT('NEW COSTAB READ IS',16F5.2)
NX=LTAB
GOTO 50
30 X(1)=-1.
DX=LTAB-1
DX=2./DX
DO 32 I=2,LTAB
32 X(I)=X(I-1)+DX
X(LTAB)=+1.
NX=LTAB
GOTO 50

```

```

47 WRITE(KOUT,600)
600 FORMAT(/' ***** NEWTAB MUST BE = EQUI OR = READ'/)
    NX=0
    GOTO 999
48 WRITE(KOUT,601) MAXCL
601 FORMAT(/' *****NUMDAT MUST BE LESS THAN ',I3/)
    ND=MAXCL
    GOTO 9
49 WRITE(KOUT,602) LENX
602 FORMAT(/' ***** LENGTH OF TABULATION MUST NOT EXCEED ',I4/)
    NX=LENX
    GOTO 10
50 CONTINUE
    IF(.NOT.DRUCK) GOTO 58
    WRITE(KOUT,603)
603 FORMAT(/' LEGENDRE COEFF. IN INPUT'/)
    WRITE(KOUT,604) (I,CL(I),I=1,ND)
604 FORMAT((5(I9,E14.6)))
58 DO 60 I=1,NX
60 CALL SUMLEG(Y(I),X(I),CL,ND-1)
    IF(.NOT.DRUCK) GOTO 68
    WRITE(KOUT,605)
605 FORMAT(/' TABULATED DISTRIBUTION:'/)
    WRITE(KOUT,606) (X(I),Y(I),I=1,NX)
606 FORMAT((5(5X,F6.3,E14.6)))
68 CALL CONTRL(X,Y,NX,CL,ND,ZZ)
999 RETURN
    END

```

```

SUBROUTINE CONTRL(X,Y,NX,CL,ND,ZZ)
C
C   CONTRL CHECKS CONVERTED DISTRIBUTION.
C   IT CALCULATES P0 AND P1 AND COMPARES IT WITH GIVEN LEGCOEFF.
C
    DIMENSION X(1),Y(1),CL(1),ZZ(1)
    DATA PI/3.14159265/,KOUT/6/
    P1=0.
    IF(ND.LT.2) GOTO 20
    DO 10 I=1,NX
10  ZZ(I)=X(I)*Y(I)
    CALL INTGRT(X,ZZ,NX,P1)
20  CALL INTGRT(X,Y,NX,P0)
    P1=P1/P0
    IF(ABS(2.*CL(1)-P0).LT.1.E-4*P0) GOTO 30
    PC=2.*PI*P0
    POS=4.*PI*CL(1)
    WRITE(KOUT,600) POS,P0
600 FORMAT(/' ***** FROM LEG SIGMA INTEGRATED=',F14.6,' FROM TAB SIG
1#A INTEGRATED=',E14.6)
    WRITE(KOUT,601)
601 FORMAT(/'10X' IF YOU WISH TO IMPROVE AGREEMENT INCREASE NUMBER OF P
POINTS IN TABULATION'/)

```

```

300
310
320
330
340
350
360
370
380
390
400
410
420
430
440
450
460
470
480
490
500
510
520
530
540
550
560
570
30 IF(ND.LT.2) GOTO 999
    POS=CL(2)/3./CL(1)
    IF(ABS(POS-P1).LT.1.E-4) GOTO 999
    WRITE(KOUT,602) POS,P1
602 FCRMAT(/' ***** FROM LEGCOEFF. MUEL(CM.)=',E14.6,' FROM TABULATI
10N MUEL(CM.)=',E14.6)
    WRITE(KOUT,601)
999 RETURN
    END

```

```

SUBROUTINE ERMSG(T1,T2)
C
C   PRINTS ERRORMESSAGE.
C
    COMMON INPUT,KOUT
    DIMENSION T2(20)
    WRITE(KOUT,601) T1,T2
601 FORMAT(/' ***** ',A4,' EXPECTED.FOUND :',4X,'*',20A4,'**'/)
    RETURN
    END

```

```

SUBROUTINE SUMLEG(Y,X,C,N)
C
C   SUMLEG BERECHNET DEN WERT EINER LEGENDRE-ENTWICKLUNG Y AN DER STELLE X
C   C(1) BIS C(N+1) SIND DIE KOEFFIZIENTEN DER ENTWICKLUNG NACH
C   P0 BIS PN.
C
    DIMENSION C(1)
    IF(N.LT.0) GOTO 9
    Y=C(1)
    IF(N.GE.1) GOTO 1
    GOTO 9
1  Y=Y+C(2)*X
    IF(N.GE.2) GOTO 2
    GOTO 9
2  Y1=1.
    Y2=X
    DO 3 I=2,N
    Y0=Y1
    Y1=Y2
    Y2=X*Y1
    Y2=Y2-Y0+Y2-(Y2-Y0)/FLOAT(I)
3  Y=Y+C(I+1)*Y2
9  RETURN
    END

```



SUBROUTINE KEDOPN(K,*)	10
IF(K.NE.0) GOTO 100	20
K=1	30
CALL LDFOPN(1, IDATUM, &200)	40
100 RETURN	50
200 RETURN	60
END	70

SUBROUTINE CDINPT(KIN)	10
DIMENSION A(20), TEND(20)	20
DATA TTEND/' ENDE' /, BLANK/' ' /	30
COMMON INPUT, KOUT	40
TEND(1)=TTEND	50
DC 100 I=2,20	60
100 TEND(I)=BLANK	70
N=1	80
1000 READ(KIN,500,END=99,ERR=9) A	90
WRITE(INPUT,500) A	100
500 FCRMAT(20A4)	110
WRITE(KOUT,501)A	120
501 FCRMAT(10X,20A4)	130
N=N+1	140
GOTO 1000	150
9 WRITE(KOUT,601) N	160
601 FORMAT('/ FEHLER BEIM LESEN DER',I5,'-TEN KARTE'/)	170
GOTO 1000	180
99 WRITE(INPUT,500) TEND	190
REWIND INPUT	200
WRITE(KOUT,602) INPUT,N	210
602 FORMAT('/ DER DATASET AUF G.FT',I2,' BESTEHT AUS',I5,' RECORDS.'/	220
15X,'ALS LETZTES RECORD IST FIN *ENDE*-RECORD HINZUGEFUEGT WORDEN')	230
RETURN	240
END	250

### 3. PROGRAMS PERFORMING ARITHMETIC OPERATIONS ON NUCLEAR DATA

In the course of nuclear data evaluation work it is quite frequently necessary to perform simple arithmetic operations on data sets such as

- multiplication with a constant (e.g. renormalisation) or
- summing with or without weighting (e.g. calculation of averages, of best estimates or of cross sections for isotopic mixtures).

Two programs for calculations of this type, CALCUL and the Karlsruhe version of SCORE, are described in the sequel. Other programs like PRTSTKED, SELPLO, SIGPLO perform also a certain amount of arithmetic calculations, but since this is not their main purpose, they are described under different headings in the present compendium (vol. V 2.1, 2.2, 4.1 respectively)

### 3.1 Simple Averages (AVERAGE)

#### Contents:

3.1.1 Purpose of AVERAGE

3.1.2 Operating description for AVERAGE

#### 3.1.1 Purpose of AVERAGE

Calculation of simple averages provides a quick and convenient way to compare data sets (e.g. data from different measurers) or to check modified (e.g. reformatted) data against the original data. The FORTRAN IV program AVERAGE was written to calculate unweighted averages of point data from three types of input sources:

- KEDAK,
- files in NEUDADA "transmission" format,
- card images.

The point data are assumed to be given as abscissae  $x_i$  (put by the program in ascending order, if necessary) and ordinates  $y_i$ ,  $i = 1, 2, \dots, n$ .

The average  $\bar{y}$  in a specified interval  $(x_A \dots x_B)$  is defined as

$$\bar{y} = \frac{1}{x_B - x_A} \int_{x_A}^{x_B} y(x) dx,$$

the function  $y(x)$  being obtained from the point data by

$$y(x) = y_i + (y_{i+1} - y_i) \frac{x - x_i}{x_{i+1} - x_i} \quad \text{if } x_i \leq x \leq x_{i+1};$$

$$y(x) = y_1 \quad \text{if } x < x_1,$$

$$y(x) = y_n \quad \text{if } x > x_n,$$

i.e. by linear interpolation within the interval of point data, and by utilization of the nearest ordinate outside this interval. The average is set equal to zero if the range of point data  $(x_1 \dots x_n)$  and the averaging interval  $(x_A \dots x_B)$  do not overlap.

The full capability of the program can be seen from the following operating description written by the author of AVERAGE in 1973. A program list is appended.

3.1.2 Operating description for AVERAGE

by R. Meyer

Contents

- 1) Introduction.
- 2) How to get the program.
- 3) How the program is controlled by the user.
- 4) Input of energy limits.
- 5) Requesting KEDAK data.
- 6) How to get NEUDADA data.
- 7) Input of card image data.
- 8) Concatenating data sets.
- 9) Quick reference for the experienced user.
- 10) Notes, Region.
- 11) Example

References

Program list

1) Introduction

AVERAGE is a program designed as a convenient means to calculate unweighted averages for neutron cross section data from three types of input sources: KEDAK, NEUDADA-format files, card image.

Its main purpose is to retrieve averages quickly and conveniently for comparison of different data sets and/or checking of derived - e.g. reformatted - data.

Since averaging is done straight forward, AVERAGE is not intended to replace any of the current group constants programs. Also, resultant averages should be used only in the manner described above. The user should not hope to get accurate results for further application purposes by this program and is referred to the respective averaging capabilities of MIGROS or other similar programs to evaluate more accurate results.

2) How to get the program 1),2)

The program is part of the program services for the nuclear data library KEDAK.

At present it is member of the dataset `INR.STEIN.LOAD` on the set-up dis `GFKO29` the mounting of which may be requested by job control cards. The name of the member is `AVERAGE`.

You may use the following job control cards to obtain the program:

```
// JOB card
/*SETUP_DEVICE=2314, ID=GFKO29
//_EXEC_FHLG, PARM.L=MAP
//L.LIB_DD_UNIT=2314, VOL=SER=GFKO29, DISP=SHR,
   DSN= INR.STEIN.LOAD
//L.SYSIN_DD_*
   _INCLUDE_LIB(AVERAGE, LDFPAC, DEFI, PRIEIN)
   _ENTRY_MAIN
/*
.
.   JOB control cards for running
.   the program - see following paragraphs.
.
// Job end card
```

Chapter 9) contains a summary of input instructions to operate `AVERAGE`. It may be used for quick reference.

Experienced programmers may read that section only and use the other paragraphs for reference in cases where chapter 9) is not sufficient.

- 1) In the following text an underscore (`_`) will be used to denote blanks.
- 2) This paragraph is intended for Karlsruhe users only.

3) How the program is controlled by the user

The user controls the program by supplying

program control input.

Program control input may be supplied by punched cards. These punched cards must immediately follow the job control card

```
//G.SYSIN_DD_*
```

This job control card together with the succeeding program control input may be placed anywhere among the job control cards for running the program (see preceding paragraph).

Program control input is supplied by `namelist input` and must follow its rules /1/. These rules will not be repeated here. However the user is invited to follow the examples given and look into reference /1/ only in crucial cases.

Only one namelist name is used: `IN`. Its parameters are explained in detail in the following sections.

Examples:

```
//G.SYSIN_DD_*  
_&IN_INPUT='KEDAK',NAMES='U 235','SGG',EMIN=1.24E+3,  
_7.86E+3,EMAX37.86E+3,19.07E+3,&END  
_&IN_INPUT='CARD',UNIT=9,FMT='(2E 14.6)',NUM=1240,  
_&END  
/*
```

The above example shows the program control input for averaging two data sets, one from KEDAK and one from card image input.



Note the general format of namelist input data

```
&IN, parameterlist, &END
```

where parameterlist is a list of names and values separated by commas:

```
Parameterlist = PARM1 = value1, PARM2 = value21, value22,...
```

Note, that one or several values may be assigned to a named variable. This is explained in more detail later.

You will also notice, that each input line starts in column two. It may be extended up to column 80 and continuation lines are allowed. An input list is terminated by &END.

You also will see, that some data are enclosed within apostrophes. These data will be interpreted as text, whereas all other data will be taken to be numeric.

The line containing /\* is optional. If used it terminates the program control input. If missing, program control input will be terminated by the next job control card encountered after // G.SYSIN...

Although not recommended, program control input may be mixed with data input. This will be shown later.

#### 4) Input of energy limits.

Upper and lower energy limits may be defined using the parameters EMIN and EMAX.

Several values may be assigned to each of them.

Averaging will be done within the energy intervals

```
EMIN (1) - EMAX (1)
```

```
EMIN (2) - EMAX (2)
```

```
⋮
```

A maximum of 100 averaging intervals may be used.

Energy limits are assumed to be given in eV.

Example:

```
//G.SYSIN_DD_*  
  &IN_INPUT = 'KEDAK', NAMEN= 'FE', 'SGT'.  
  EMIN = 1.E6, 2.E6, 3. E6, 10.E6 ,  
  EMAX = 2.E6, 3.E6, 4. E6, 12.E6,  
  &END
```

```
_ &IN_NAMEN (2) = 'SGG', &END  
_ &IN_NAMEN(2) = 'SGN', &END  
/*
```

In this example the first input list requests averaging of KEDAK data for the total cross section of iron in the intervals <1MeV, 2MeV> ,<2MeV, 3MeV>, <3MeV,4MeV>, <10MeV,12MeV>. The second input list requests averaging for the capture cross section of iron. Since no new energy limits are defined, the previous ones will be used. The same holds for the third line, where averaging for the elastic cross section of the same material will be done.

Note that the intervals need not be continuous. Therefore minimum and maximum energy both are required. A limiting energy need not be contained in the tabulation energies but interpolation will be performed. The requested data however must cover the specified energy intervals.

Attention:

- linear interpolation will be used always. If linear interpolation is not sufficient, a suitable data set may be prepared using the interpolation capabilities of SIGPLO /2/ or any other interpolating program. Note however, that AVERAGE is only intended to give good estimates of averages for quick comparison or check.

- It is not possible to redefine energy limits by the following method:

```
//G.SYSIN_DD_x  
_ &IN_INPUT='KEDAK',NAMEN='FE','SGT',  
_ EMIN=1.E6,2.E6,EMAX=2.E6,3.E6,&END  
_ &IN_TYP='SGG',EMIN(2)=3.E6,EMAX(2)=4.E6,&END  
/*
```

In fact, the second input list would not redefine any energy limits. If energy limits are given, all of them must be

specified from the very beginning.

They will then make up the new set of averaging intervals.

That is a complete set of energy intervals must be input.

It is not possible to just modify the current one.

The correct procedure for the above example would be:

```
//G.SYSIN_DD_*  
_&IN_INPUT='KEDAK',NAMEN='FE','SGT',  
_EMIN=1.E6,2.E6,EMAX=2.E6,3.E6,&END  
_&IN_TYP='SGG',EMIN=1.E6,3.E6,EMAX=2.E6,4.E6,&END  
/*
```

5) Requesting KEDAK data

The preceding paragraphs already gave a few examples of retrieving data from the KEDAK library.

Use the following parameters for this purpose:

|| INPUT, NAMES, NAMZ, PLEH ||

a) With INPUT you specify the kind of input source. In the present case you code INPUT = 'KEDAK'.

If the current value of INPUT is already 'KEDAK' you need not respecify it.

b) PLEH allows you to specify the FORTRAN data set reference number for the KEDAK library to address. Thus it is possible to address two or more KEDAK libraries in one job.

For each KEDAK library addressed you must supply the corresponding DD-card. Do not forget that you will need 8K buffer for each KEDAK library usually.

Example:

If you code PLEH = 1 the KEDAK library is expected on logical unit 1. A DD-card describing the KEDAK library must be present:

```
//G.FTO1FOO1_DD_UNIT=2314,VOL=SER=GFK050,DISP=SHR,DSN=KEDAK3
```

This DD-card would address the present released version of KEDAK. You insert this DD-card anywhere among the job control cards for running the program (see paragraph 1). But you must not mix it with your program control data. PLEH = 1 is the default value: you need not specify PLEH if it is to be 1.

c) You use NAMES to address a specific cross section set on KEDAK, in the order: material name, data type name,

and if applicable also an excitation energy, e.g.:

```
NAMES='FE','SGT'  
NAMES='U_238','SGF'  
NAMES='PU239','SGIZ',0.  
NAMES='FE___EN3','SGG'
```

You will notice, that each non-numeric name must be enclosed in apostrophes. For definition and explanation of KEDAK names you should refer to /3/.

d) NAMZ is a parameter to specify the number of names necessary to address a KEDAK data set. Its default value is 2.

Example:

```
//G.SYSIN_DD_*  
_&IN_NAMES='FE','SGT',EMIN=1.E+6,EMAX=1.E+7,&END  
_&IN_NAMES(2)='SGN',&END  
_&IN_TYP='SGP',&END  
_&IN_MAT='FE___EN3',TYP='SGT',NAMZ=2,PLEH=2,&END  
_&IN_TYP='SGN',&END  
_&IN_TYP='SGP',&END  
_&IN_TYP='SGIZ',EXC=0.,NAMZ=3,&END  
_&IN_MAT='FE',TYP='SGX',PLEH=1,NAMZ=2,&END  
/*  
//G.FTO1FOO1_DD_UNIT=2314,VØL=SER=NUSYSO,DISP=SHR,DSN=KNDF  
//G.FTO2FOO1_DD_UNIT=2314,VØL=SER=GFKO29,DISP=SHR,DSN=ENDFB3.CONV  
/*SETUP_DEVICE=2314,ID=NUSYSO
```

In this example averages will be calculated for the interval <1MeV, 10MeV> for  $\sigma_T$ ,  $\sigma_n$ ,  $\sigma_p$  and the excitation cross section of the first inelastic level for two materials: 'FE' and 'FE\_\_\_EN3', where the former material is obviously to be retrieved from the released version of the KEDAK library

via // G.FT01FO01- since no PLEH is given and therefore is one by default - and the second material is some conversion from ENDF/B3 and is taken from a data library that is specified by // G.FT02FO01...

- . Note, that in the last input line it is necessary to specify PLEH = 1 although this was the default value, since its value had been changed from default one to two by input line 5.
- . Also note, that it is necessary to redefine NAMZ = 2 in input lines 5 and 9.
- . Note the alternative use of MAT, TYP and EXC for NAMES (1) , NAMES (2), NAMES (3) respectively. Coding NAMES (1) or MAT is completely identical in effect. The same holds for the other alternatives.

Attention:

- . A maximum of five KEDAK libraries may be used in one run. The addressed materials must be uniquely defined, i.e. no equally named materials must be contained in two different libraries, or confusion might occur. 8K buffer area will be required by each addressed library normally.
- . If a different package of retrieval routines is used for the KEDAK library than given in paragraph 2) PLEH = 1 may be used only.
- . Energies are expected to be in eV.

6) How to get NEUDADA-data

The following parameters control retrieval of data from NEUDADA-tapes<sup>x)</sup>:

||INPUT, UNIT, NEUFOR, NAMES, SCALE||

a) INPUT

You request data to be retrieved from a data set in NEUDADA-format, if you code INPUT = 'NEUDADA'. If the current value of INPUT already is 'NEUDADA' you need not respecify it. Simply omit parameter INPUT in this case.

b) UNIT

You must specify the Fortran data set reference number which applies to the data to be retrieved, e.g. UNIT = 3 would imply that the data shall be fetched from a data set described by // G.FT03F0001\_DD \_ .....

You will observe that each unit number is linked with a DD-card. This DD-card must describe the features of the NEUDADA-format data set. See below for examples.

It is absolutely necessary to specify the unit number, since its default value (UNIT = 5) must not be applied to NEUDADA data. 2K buffer area will be needed for each NEUDADA file normaly.

c) NEUFOR

NEUDADA tapes are distributed by CCDN/Saclay usually in the NEUDADA Transmission Format. On request it is also possible to obtain data in Internal Expanded Format. The latter format reduces the amount of storage and retrieval time needed to keep and access the data.

Code NEUFOR = 1 for data in Internal Expanded Format,  
NEUFOR = 2 for data in Transmission Format.  
NEUFOR = 2 is the default value. If not previously  
respecified by the user, the parameter may be omitted if  
the default value should be applied.

d) NAMES

Data sets in both NEUDADA formats get an identifier of  
six characters in length. With each NEUDADA tape you will  
receive a list displaying the contents of the tape and the  
identifiers given to the various data sets on the tape.  
If you wish to retrieve a given data set, you should look  
up its corresponding identifier on that list. Specify  
this identifier in your program control input, e.g.

NAMES = 'E00001' or

NAMES = 'U00049'

Make sure, that the desired data have received a unique  
identifier. If not, use the capability of AVERAGE to  
concatenate these two or more sets. Refer to paragraph 8)  
for this purpose.

Attention:

The current contents of NAMES are destroyed when requesting  
NEUDADA input, i.e. although the user actually only fills  
NAMES (1), the current values of NAMES (2) , .... are  
destroyed since they will be used by the program when  
it retrieves the data.

Therefore do not expect in the following example:

```
&IN INPUT='KEDAK',NAMES='FE','SGT',&END  
&IN INPUT='NEUDADA',NEUFOR=2,NAMES='E00022',&END  
&IN INPUT='KEDAK',NAMES='CR',&END
```

to obtain the total cross section of chromium in the third  
input line.



e) SCALE

Energies on NEUDADA format data sets are expected to be given in MeV, cross section values in barn.

In many time-of-flight measurements however cross sections are given in barns  $\times \sqrt{E}$ . These cases will be indicated on the list you receive from CCDN/Saclay together with the tape. The program AVERAGE optionally permits you to have cross section data converted from this scale to the normal scale (barns) by specifying SCALE = 'SQE'. You may nullify this specification by coding SCALE = '-', which is the default as long as no other value has been specified.

Example:

```
//G.SYSIN_DD_*
  _&IN_INPUT='NEUDADA',UNIT=2,NAMES='E00013',NEUFOR=2,
  _EMIN=0.25,1.02,2.66,4.81,
  _EMAX=1.02,2.66,4.81,9.67,
  _&END
  _&IN_NAMES='E00016',&END
  _&IN_NAMES='U00027',&END
  _&IN_NAMES='E00001',UNIT=3,NEUFOR=1,&END
/*
//G.FTO2FOO1_DD_UNIT=TAPE7,VOL=SER=NDO607,LABEL=(2,NL,,IN),
//_DCB=(RECFM=FB,LRECL=132,BLKSIZE=1320,TRTCH=ET,DEN=1)
//_DISP=(OLD,PASS)
/*SETUP_DEVICE=TAPE7,ID=(NDO607,NORING,,NL)
//G.FTO3FOO1_DD_UNIT=TAPE9,VOL=SER=NDO657,LABEL=(2,NL,,IN),
//_DCB=(RECFM=FB,LRECL=72,BLKSIZE=720),DISP=(OLD,PASS)
/*SETUP_DEVICE=TAPE9,ID=(NDO657,NORING,NL)
```

In this example averaging is requested in the four intervals <0.25eV,1.02 eV> , <1.02eV, 2.66eV> , <2.66eV,4.81eV> , <4.81eV,9.67eV> for four data sets, all residing on NEUDADA-tapes.

The first three data sets 'E00013', E00016', '000027' are on the 7-track tape NDO607 in transmission format.

This tape is accessed by // G.FT02F001\_DD \_ .. and the respective Setup card.

The fourth data set 'E00001' is on 9-track tape NDO657 in internal expanded format. This tape is accessed by // G.FT03F001 \_ DD \_ ... and the respective Setup card.

Note the following:

- . The above two DD-cards are two typical examples of how the DD-cards for NEUDADA-tapes must look like. If you do not have any experience in coding DD-cards you should take the paper label which is distributed with each tape and contact some more experienced person for coding the DD-statements. Bring the above lines as example with you. The paper label will contain all information to build up the necessary DD-card.  
Since CCDN/ Saclay presently is operating its IBM 360/30 in DOS the difference in tape marks between DOS and OS must be considered usually leading to the result that DOS-files 1,2,3, .... will be found in OS as files 2,4,6, .....
- . Be aware of the fact, that energy limits are always specified in eV, no matter what units the energies of the requested data points are. The program will take care of the necessary conversion.  
Conversion always will be to eV.

Program execution may be slowed down appreciably, if data sets are not specified in the order they reside on tape. Backspacing and rewinding however will be provided by the program automatically.

Note that the program expects the tape to be numerically sorted in ascending order, i.e. it expects E00013 to be physically behind E00012 and U00002 behind U00001.

It does not have any expectation concerning the sorting of E and U.

7) Input of card image data

A third input possibility is offered for all those data sets to which neither of the preceding formats apply. Data may also be input from card image data sets with the applicable format being defined by the user. The following parameters are available for card image input:

||INPUT, UNIT, FMT, NUM,C||

a) INPUT

To request card image input you must code INPUT = 'CARD'. It is not necessary to repeat this specification if the current value of INPUT already is 'CARD'.

b) UNIT

By UNIT = nn you specify the data set reference number to be used for reading the data input, e.g. coding UNIT = 9 will cause the requested data to be read in from  
// G.FTO9FOO1 \_ DD \_ ... Do not forget to supply a DD-card for each data set you reference. You may insert this DD-card anywhere among the job control cards for running the job (see paragraph 2). You must not mix this DD-card however with any of your data input or program control input cards.

If you omit the UNIT-specification, data will be read from logical unit five (unless no different specification preceded). Logical unit five is the program control data set. Although it is not recommended it is thus possible to mix data- and program control input.

c) FMT

With FMT you specify the input format to be used for reading the data. Since no default value is available you must specify a format.

Note that input energies are expected to be in eV.  
However by using a suitable scaling factor in your format (/1/) you may select different energy units for your input data.

d) NUM

Gives the number of data points to be read. Be sure to use the correct number since an error in NUM may cause the job to fail. For the maximum number of data points permitted see the respective paragraph.

e) C

C may be set .True. or .False. respectively if a comment card precedes the data set to be read or not.

Code C = .TRUE. or C = T if so ,

or C = .FALSE. or C = F else.

It is highly recommended that such a card precedes the data set. It will be read and printed by the program and offers a convenient means to identify the data set.

Output from SIGPLO /2/ will contain such heading identifying cards.

If it was not respecified by the user, C will be .TRUE. by default.

Example:

```
//G.SYSIN_DD_*
_&IN_INPUT='CARD',UNIT=9,FMT='(2E10.5)',NUM=5,
_&EMIN=1.,EMAX=5.,&END
_&IN_UNIT=5,NUM=4,&END
NOTE_THAT_INPUT_READ_FROM_UNIT_5_MUST_IMMEDIATELY_FOLLOW.
```

1.0		1.0
3.0		1.8
4.0		2.5
6.2		1.9

↑  
----- Column 10

```
_ &IN_UNIT=9,NUM=2,&END
_ &IN_UNIT=10,NUM=4,FMT='(10F8.5)',&END
_ &IN_NUM=3,&END
/*
//G.FT10FOO1_DD_*
COMMENT:THIS_IS_THE_FIRST_INPUT_DATA_SET_FROM_UNIT_10.
    0.5      1.9      2.8      0.97      2.4      1.01      5.08      0.42
COMMENT:THIS_IS_THE_SECOND_INPUT_FROM_UNIT_10.
    6.0      1.87     2.0      1.97      0.99      1.88
/*
//G.FT09FOO1_DD_*
HEADING_CARD_1
    0.97      1.87
    1.62      0.98
    3.05      1.17
    4.99      1.86
    5.17      2.88
HEADING_CARD_2
    1.0      1.0
    5.0      1.0
/*
```

From this example you may learn:

- data are read in this order  $E(1)$ ,  $\sigma(1)$ ,  $E(2)$ ,  $\sigma(2)$ , ...  
using the specified format.
- Each card may contain one or more data points.  
This is exclusively a matter of format.
- If data are to be read from the control unit 5, they must  
immediately follow the requesting control card, as is the case  
for input list number two.
- Data need not be ordered. The program will reorder them if  
necessary.
- It is essential that you observe your format specifications  
and punch the data in the appropriate columns.

Notes

Although not used in the above example, the data set may of course reside also on tape or disc and its line length may exceed 80 columns. Comment will always however be read in 80 columns length. Therefore no shorter line length should be used or the job will fail.

8) Concatenating data sets.

In some cases it might be necessary to combine two or more sets of data into one before averaging.

This can be done by using the parameter ADD.

If you specify ADD = .TRUE. or ADD = T the next set of data to be read will be added to the one currently read. Averaging will be postponed.

Example:

```
//G.SYSIN_DD_*
  &IN_INPUT='NEUDADA',NAMES='E00014',UNIT=1,ADD=T,
  &EMIN=1.67E+4,EMAX=2.36E+4,&END
  &IN_NAMES='E00015',&END
  &IN_NAMES='E00016',&END
  &IN_INPUT='CARD',NUM=1,FMT='(2F10.5)',C=F,ADD=F,&END
    1.67E+4  0.0
  &IN_INPUT='KEDAK',NAMES='FE','SGT',UNIT=2,&END
/*
//G.FTO1FOO1_DD_UNIT=TAPE9,VOL=SER=ND2108,DISP=(OLD,PASS),
//_LABEL=(2,NL,,IN),
//_DCB=(RECFM=FB,LRECL=132,BLUSIZE=1320)
//G.FTO2FOO1_DD_UNIT=2314,VOL=SER=GFK050,DISP=SHR,
//_DSN=KEDAK3
/*SETUP_DEVICE=2314,ID=GFK050
/*SETUP_DEVICE=TAPE9,ID=(ND2108,NORING,,NL)
```

In this example the three data sets 'E0014', 'E00015', 'E00016' on NEUDADA-tape and the single data point from card input are combined into one set of data. Note that this is possible since the data are being reordered before they are used. Averaging is performed after the last data set to be added has been read in, which is indicated by ADD = F.



Attention:

Do not forget to turn off the ADD-function when requesting input of the last set of data to be added, or unwanted results will be obtained.

9) Quick reference for the experienced user.

The program AVERAGE is controlled by program control input. This input is supplied via // G.SYSIN \_ DD \_ .... and is assigned logical unit 5. It is in the form of namelist data. The only namelist name being used is IN. The following input parameters may be used:

EMIN }  
EMAX } to specify averaging intervals

INPUT = { 'CARD' for requesting card image - ,  
'KEDAK' KEDAK - , or NEUDADA input  
'NEUDADA'

NAMES }  
NAMZ } to control retrieval from KEDAK  
UNIT }

NAMES }  
UNIT } to control retrieval from NEUDADA files.  
NEUFOR }  
SCALE }

UNIT }  
FMT } to control card image input.  
NUM }  
C }

ADD - to combine two or more data sets before averaging.

Averaging intervals are given by entering values for EMIN, EMAX. Up to 100 values may be entered for each. Each pair {EMIN (1), EMAX (1)} sets up one averaging interval . The values must be given in eV. The set of intervals remains in effect until a new set is entered. Only complete sets can be entered at a time. If an EMAX (1) < EMIN (1), the corresponding average will be set 0.

INPUT source is requested by parameter INPUT. Its value remains in effect until a new one is specified.

KEDAK data may be retrieved by specifying the appropriate names of the requested data set by NAMES =, giving the number of names by NAMZ (if different from two) and specifying the logical unit to access the Kedak library by UNIT = nn. For each accessed KEDAK library a DD-card must be available. Up to five KEDAK libraries may be accessed in one job step. Each KEDAK library accessed will eat up 8K of buffer area.

If a retrieval routine package other than LDFPAC is used, the UNIT-number must be 1 and only one library can be accessed within one job step.

Energies are expected to be in eV.

NEUDADA data may be retrieved from both transmission format or internal expanded format files. This is controlled by specifying NEUFOR = 2 or NEUFOR = 1 respectively, where the former value is default. The 6-character data set name is specified by NAMES = '.....'. Note that all locations of NAMES are used and therefore previous contents of NAMES (2), NAMES (3), ... are not kept. The program expects the identifier to consist of a leading character followed by five decimal digits, and that within each character group data sets are ordered according to increasing numeric part. Backspacing and Rewind is controlled by the retrieval program on the basis of this assumption. The logical unit assigned to the NEUDADA file is given by UNIT = nn.

Optionally, the user may switch from cross section values in barn to barn  $\times \sqrt{E}$  by specifying SCALE = 'SQE'.

The program then will calculate the correct cross section values.

This specification is nullified by SCALE = ' \_ ' ,

Energies are expected to be in MeV and are converted to eV accordingly.

For data set attributes and data set identifiers consider the labels and listings accompanying each tape.

Card image input for all other kinds of data requires specification of the input format via FMT. The format specification may not exceed 40 characters. The minimum line length is eighty. An optional heading card for reading and printing only is read to 80 characters length only. If this length is exceeded, the text will be truncated. The heading card will be read if logical variable C is true (default), else it will be assumed that the starting record of the data set already contains numerical data. Scaling of data to eV/barns must be done by format scaling factors, if necessary.

UNIT specifies the logical unit to be used for reading. If 5, the data records must immediately follow the namelist requesting them. If not 5, an appropriate DD-card must be available. NUM specifies the number of data points to be read.

Data points need not be sorted.

Combination of data sets is requested by setting logical variable ADD to .True. . If so, the data requested by the current namelist input are not averaged but kept and combined with the data requested by the succeeding namelist input. This is repeated until ADD = F terminates this process. ADD = F must be coded on the last data set to be concatenated and will cause averaging of the combined data. Data sets need not be concatenated in their numerical order.

Defaults if mentioned apply to program start.

Values assigned to any parameter (except NUM) remain in effect until respecified by the user and thus form a set of "current defaults".

10) Notes, Region

- Data will be retrieved from KEDAK/NEUDADA within the energy range defined by the lowest valued EMIN and the maximum value of EMAX. Card image data will be read as specified by NUM.
- Interpolation to limiting energies is done linear in x/linear in y. The same interpolation rule is used for integration.
- Data need not be sorted, but will be sorted by the program if necessary.
- Running times are difficult to estimate since very different storage media may be used. It should be emphasized however, that the averaging procedure usually will require least of the running time.
- Allocation of working area for the cross section data is dynamical. 30 % of the available free region will be reserved for buffers, up to a maximum of 40K buffer area. Jobs requiring more buffer space cannot be run but must be separated to reduce buffer request. Allocation of buffer and working area is displayed to the user, so he may after an initial run easily estimate the actual REGION parameter he must apply provided he knows the maximum number of data points to be averaged. Note that you always may reduce the required amount of core by dividing one set of averaging intervals into several.
- Output is hoped to be self- explanative and is not discussed here.

11) Example:

This last example is designed to give a quick overview of the job control cards and program control input necessary to run the program.

```
//INR613AV_JOB_(0613,101,PO000),MEYER,REGION=24OK,TIME=3
//_EXEC_FHLG,PARM.L=MAP
/XSETUP_DEVICE=2314,ID=GFKO29
/*SETUP_DEVICE=2314,ID=GFKO50
/*SETUP_DEVICE=TAPE7,ID=(ND5321,NORING,,NL)
//L.LIB_DO_UNIT=2314,VOL=SER=GFKO29,DISP=SHR,DSN=INR.STEIN.LOAD
//L.SYSIN_DD_*
_INCLUDE_LIB(AVERAGE,LDFPAC,DEFI,PRIEIN)
_ENTRY_Main
/*
//G.FTO1FOO1_DD_UNIT=2314,VOL=SER=GFKO50,DISP=SHR,DSN=KEDAK3
//G.FTO2FOO1_DD_UNIT=2314,VOL=SER=GFKO29,DISP=SHR,DSN=MEYER.ENDFB3
//G.FTO3FOO1_DD_UNIT=TAPE7,VOL=SER=ND5321,DISP=(OLD,PASS),
//_LABEL=(2,NL,,IN),
//_DCB=(RECFM=FB,LREC=132,BLKSIZE=1320,TRTCH=ET)
//G.FTO9FOO1_DD_UNIT=2314,VOL=SER=GFKO29,DISP=SHR,
//_DSN=MEYER.AVERG.CARDIM
//G.SYSIN_DD_*
_IN_UNIT=1,INPUT='KEDAK',NAMES='NA-23','SGG',
_EMIN=0.5E+3,1.E+3,2.E+3,2.5E+3,3.E+3,4.E+3,6.E+3,8.E+3,10.E+3
_EMAX=1.E+3,2.E+3,2.5E+3,3.E+3,4.E+3,6.E+3,8.E+3,10.E+3,15.E+3,
_END
_IN_UNIT=2,MAT='NA_23EN3',&END
_IN_UNIT=3,NAMES='EOOOO1',INPUT='NEUDADA',ADD=T,
&END
_IN_NAMES='EOOO14',ADD=F,&END
_IN_UNIT=9,INPUT='CARD',FMT='(2E14.6)',NUM=1487,&END
/*
//job end card
```

The reader is invited to interpret the above example himself.

References

- /1/ IBM 05/360 FORTRAN IV Language manual
- /2/ R. Meyer, SIGPLO program description to be published
- /3/ B. Krieg, Program services of the evaluated nuclear data library KEDAK, Part I, KFK 1725/I

AVERAGE: Program List

DIMENSION X(1),EQ(100),EU(100),S(100)	10	WRITE(KOUT,611) UNIT,NEUFOR	560
DIMENSION EMIN(100),EMAX(100),NARG(5)	20	611 FORMAT(10,'UNIT=',I2,' NEUFOR=',I2/)	570
REAL*8 NAMES(3),NAMOLD(3)/3*' ' /, INPUT, CARD/'CARD',KEDAK/'	30	19 IF(ADAT) GOTO 35	580
IKEDAK/' /,FMT(5)/5*' ' /,TXT(10),NEUDAD/' NEUDADA' /,SQ/'SQE' /,	40	DO 20 I=1,MAXE	590
2SCALE/' ' /,MAT,TYP,EXC	50	IF(EMIN(I).NE.0.OR.EMAX(I).NE.0) GOTO 20	600
COMMON/INOUT/KOUT	60	NI=I-1	610
LOGICAL C/T/,ADD/F/,ADAT/F/,STANZ/F/	70	GOTO 30	620
INTEGER UNIT,PLEH	80	20 CONTINUE	630
LOGICAL*1 KEKONT(99)/99*F/	90	NI=MAXE	640
EXTERNAL LDFLOC,LDFNXT,NEULOC,NEUNXT,PRIEIN	100	30 IF(NI.NE.0.OR.NNI.NE.0) GOTO 32	650
EQUIVALENCE (KE,UNIT),(NAMES(1),MAT),(NAMES(2),TYP),(NAMES(3),EXC)	110	WRITE(KOUT,601)	660
NAMELIST/IN/NAMES,NAMZ,EMIN,EMAX,INPUT,FMT,KE,NLM,C,NEUFOR,UNIT,	120	601 FORMAT(' NO ENERGY LIMITS GIVEN')	670
1 SCALE,ADD,MAT,TYP,EXC,STANZ,KSTANZ,DIFV,MAXINT,PLEH	130	GOTO 1000	680
DATA MAXE/100/,KIN/5/,NNI/0/,NAMZ/2/	140	32 IF(NI.EQ.0) GOTO 35	690
KOUT=6	150	NNI=NI	700
KSTANZ=7	160	DO 34 I=1,NI	710
MAXINT=25	170	EC(I)=EMAX(I)	720
DIFV=0.0	180	EU(I)=EMIN(I)	730
PLEH=1	190	34 CONTINUE	740
CALL PRIEIN(KIN,KOUT,60)	200	EA=XMIN(NNI,EU)	750
NEUFOR=2	210	ER=XMAX(NNI,EO)	760
KE=KIN	220	35 IF(INPUT.EQ.KEDAK) GOTO 36	770
CALL SPACE(X(1),LX,NMAX)	230	IF(INPUT.EQ.NEUDAD) GOTO 50	780
5 DO 10 I=1,MAXE	240	IF(.NOT.C) GOTO 31	790
EMIN(I)=0	250	READ(KE,500) TXT	800
EMAX(I)=0	260	500 FORMAT(10A8)	810
10 CONTINUE	270	WRITE(KOUT,607) TXT	820
READ(KIN,IN,END=1010)	280	607 FORMAT(LX,'COMMENTCARD=***',10A8,'***'/)	830
MAXINT=MAX(1,MIN(MAXINT,100))	290	31 IF(NUM+NTOT.GT.NMAX) GOTO 990	840
IF(DIFV.LE.0) GOTO 2	300	NTOX=NTOT+LX-1	850
EMAX(1)=EMIN(1)+DIFV	310	NTOY=NTOT+LX+NMAX-1	860
IF(MAXINT.LE.1) GOTO 6	320	READ(KE,FMT) (X(NTOX+I),X(NTOY+I),I=1,NUM)	870
DO 3 I=2,MAXINT	330	NUMX=NUM	880
EMIN(I)=EMAX(I-1)	340	NAMOLD(1)=CARD	890
3 EMAX(I)=EMIN(I)+DIFV	350	GOTO 39	900
6 II=MAXINT+1	360	36 NARG(1)=NAMZ	910
IF(II.GT.100) GOTO 2	370	DC 37 I=1,NAMZ	920
DC 7 I=II,100	380	IF(NAMOLD(I).NE.NAMES(I)) GOTO 38	930
EMIN(I)=0.0	390	37 CONTINUE	940
7 EMAX(I)=0.0	400	GOTO 39	950
2 IF(.NOT.ADAT) NTOT=0	410	38 CALL RETXS(NARG,NAMES,EA,EB,X(NTOT+LX),X(NTOT+LX+NMAX),NUMX,	960
WRITE(KOUT,605) INPUT	420	INMAX-NTOT,NR,LDFLOC,LDFNXT)	970
605 FORMAT(1H1,' TYPE OF INPUT REQUESTED=',A8/)	430	DO 33 I=1,NAMZ	980
IF(INPUT.NE.KEDAK) GOTO 12	440	33 NAMOLD(I)=NAMES(I)	990
WRITE(KOUT,600) (NAMES(I),I=1,NAMZ)	450	IF(NR.GE.2.AND.NR.NE.10) GOTO 990	1000
600 FORMAT(' AVERAGING FOR ',2(A8,2X),1P4E13.5/)	460	GOTO 39	1010
IF(.NOT.KEKONT(PLEH)) CALL LDFOPN(PLEH,IDAT,&1C1C)	470	50 NARG(1)=3	1020
KEKONT(PLEH)=.TRUE.	480	NARG(2)=1	1030
GOTO 19	490	NARG(3)=1	1040
12 IF(INPUT.EQ.NEUDAD) GOTO 14	500	NARG(4)=UNIT	1050
WRITE(KOUT,606) NUM,KE,FMT	510	NARG(5)=NEUFOR	1060
606 FORMAT(' AVERAGING FOR CARD INPUT REQUESTED:*/	520	NAMES(2)=UNIT	1070
1' NUM=',I5,' KE=',I2,' FMT=',5A8/)	530	NAMES(3)=EA	1080
GOTO 19	540	DC 51 I=1,3	1090
14 WRITE(KOUT,600) NAMES(1)	550	IF(NAMES(I).NE.NAMOLD(I)) GOTO 52	1100



```

51 CONTINUE
   GOTO 39
52 EX=EA*1.F-6
   EY=EB*1.F-6
   CALL RETXS(NARG,NAMES,EX,EY,X(NTOT+LX),X(NTOT+LX+NMAX),NUMX,
   1NMAX-NTOT,NR,NEULOC,NEUNXT)
   DC 53 I=1,3
53 NAMOLD(I)=NAMES(I)
   IF(NR.GE.2.AND.NR.NE.10) GOTO 990
   J=NTOT+LX
   DC 54 I=1,NUMX
   X(J)=X(J)*1.E+6
54 J=J+1
   IF(SCALE.NE.SQ) GOTO 56
   J=NTOT+LX+NMAX
   DC 55 I=1,NUMX
   X(J)=X(J)/SQRT(X(J))
55 J=J+1
56 CONTINUE
39 NTOT=NTOT+NUMX
   IF(.NOT.ADD) GOTO 57
   ACAT=.TRUE.
   GOTO 5
57 ACAT=.FALSE.
   CALL ORDREN(NTOT,X(LX),X(LX+NMAX))
   CALL EQUEN(NTOT,X(LX),X(LX+NMAX))
   WRITE(KOUT,602)
602 FORMAT(/6X,'FROM',T24,'TO',T36,'AVERAGE'/)
   DO 40 I=1,NNI
   CALL AVERG(X(LX),X(LX+NMAX),NTOT,EU(I),EO(I),S(I))
   WRITE(KOUT,610) I,EU(I),EO(I),S(I)
610 FORMAT('0',I5,2X,1P3E15.5)
   IF(.NOT.STANZ) GOTO 40
   EUO=0.5*(EU(I)+EO(I))
   WRITE(KSTANZ,612) EUO,S(I)
612 FORMAT(2E14.6)
40 CONTINUE
   GOTO 1000
990 IF(NR.NE.2) GOTO 992
   WRITE(KOUT,603)
603 FORMAT(' NOT ENOUGH INCORE STORAGE FOR THIS DATA-TYPE.*/)
   GOTO 1000
992 WRITE(KOUT,604)
604 FORMAT(' NO DATA IN REQUESTED ENERGY INTERVAL(S).*/)
   GOTO 1000
1000 ACAT=.FALSE.
   IF(ADD) ADAT=.TRUE.
   GOTO 5
1010 STCP
   END

```

```

1110 FUNCTION XMIN(N,X)
1120 DIMENSION X(1)
1130 XMIN=X(1)
1140 IF(N.LT.2) GOTO 100
1150 DO 10 I=2,N
1160 IF(X(I).LT.XMIN) XMIN=X(I)
1170 10 CONTINUE
1180 100 RETURN
1190 END
1200
1210
1220
1230
1240
1250 FUNCTION XMAX(N,X)
1260 DIMENSION X(1)
1270 XMAX=X(1)
1280 IF(N.LT.2) GOTO 100
1290 DO 10 I=2,N
1300 IF(X(I).GT.XMAX) XMAX=X(I)
1310 10 CONTINUE
1320 100 RETURN
1330 END
1340
1350
1360
1370
1380
1390
1400 C SUBROUTINE AVERG(X,Y,N,XA,XB,YAV)
1410 C
1420 C CALCULATION OF UNWEIGHTED AVERAGE IN A GIVEN RANGE WHERE
1430 C A FUNCTION Y(X) IS GIVEN BY A DISCRETE NUMBER OF POINTS.
1440 C LINEAR INTERPOLATION IS USED BETWEEN THESE POINTS.
1450 C X(K): ABSCISSAE
1460 C Y(K): ORDINATES
1470 C N : NUMBER OF CO-ORDINATE PAIRS
1480 C XA : LOWER LIMIT OF AVERAGING INTERVAL
1490 C XB : UPPER " " " "
1500 C YAV : UNWEIGHTED AVERAGE
1510 C
1520 C DIMENSION X(1),Y(1)
1530 C
1540 C YAV=0.
1550 C RETURN UNLESS THE RANGES XA...XB AND X(1)...X(N) OVERLAP
1560 C IF(XA.GE.XB.OR.XA.GE.X(N).OR.XB.LE.X(1))RETURN
1570 C RETURN IF NOT ALL X(K) ARE DIFFERENT AND IN ASCENDING CRDR
1580 C IF(N.EQ.1) GO TO 2
1590 C DO 1 K=2,N
1600 C IF(X(K).LE.X(K-1))RETURN
1610 C 1 CONTINUE
1620 C FIND SUBSCRIPT KA OF LARGEST X(K) NOT EXCEEDING XA
1630 C 2 DO 3 K=1,N
1640 C KA=K-1
1650 C IF(X(K).GT.XA)GO TO 4
1660 C 3 CONTINUE
1670 C FIND SUBSCRIPT KB OF LARGEST X(K) NOT EXCEEDING XB

```

```

4 DO 5 K=1,N
  KB=N+1-K
  IF(X(KB).LT.XB)GO TO 6
5 CONTINUE
C   CALCULATE ABSCISSAE YA,YB FOR ORDINATES XA,XB
6 IF(KA.EQ.0)YA=Y(1)
  IF(KA.GT.0)YA=Y(KA)+(Y(KA+1)-Y(KA))*(XA-X(KA))/(X(KA+1)-X(KA))
  IF(KB.EQ.N)YB=Y(N)
  IF(KB.LT.N)YB=Y(KB)+(Y(KB+1)-Y(KB))*(XB-X(KB))/(X(KB+1)-X(KB))
C   CALCULATE AVERAGE
C   IF(KA.LT.KB)GO TO 7
C   (1) NO MESH POINTS BETWEEN XA AND XB
  YAV=0.5*(YA+YB)
  RETURN
C   (2) AT LEAST ONE MESH POINT BETWEEN XA AND XB
7 S=(Y(KA+1)+YA)*(X(KA+1)-XA)+(YB+Y(KB))*(XB-X(KB))
  IF(KB.EQ.KA+1)GO TO 9
  K1=KA+1
  K2=KB-1
  DO 8 K=K1,K2
8 S=S+(Y(K+1)+Y(K))*(X(K+1)-X(K))
9 YAV=0.5*S/(XB-XA)
  RETURN
END

```

300  
310  
320  
330  
340  
350  
360  
370  
380  
390  
400  
410  
420  
430  
440  
450  
460  
470  
480  
490  
500  
510  
520  
530

```

10 IF(NREW.NE.0) GOTO 1001
  REWIND IFIL
  NREW=NREW+1
15 READ(IFIL,500,END=10) AREQ,IREQ
500 FCRMAT(T42,A1,I5)
  IF(AR.NE.AREQ.OR.IR.NE.IREQ) GOTO 15
  BACKSPACE IFIL
  READ(IFIL,FMT) TXT,AREQ,IREQ,XX
16 WRITE(KOUT,600) TXT,AREQ,IREQ,IFIL
600 FCRMAT(//10X,42A1,I5,' FOUND ON UNIT ',I2//)
  X(1)=XX(1)
  J=2
  IF(NARG(2).NE.2) GOTO 22
  X(2)=XX(2)
  J=3
22 X(J)=XX(6)
  IF(NARG(3).NE.2) GOTO 24
  J=J+1
  X(J)=XX(7)
24 RETURN
1001 CONTINUE
  WRITE(KOUT,601) NAMES(1),IFIL
601 FORMAT(' ',A8,' NOT FOUND ON UNIT=',I2,' OR')
  NR=0
  RETURN
END

```

280  
290  
300  
310  
320  
330  
340  
350  
360  
370  
380  
390  
400  
410  
420  
430  
440  
450  
460  
470  
480  
490  
500  
510  
520  
530

```

SUBROUTINE NFULCC(NR,NARG,NAMES,X)
COMMON/INOUT/KOUT
REAL*8 NAMES(1),NAM,FMT(2)/'(42A1,I5',' ',FMT2/,'7A4)'/,FMT3/,'7F
112.5)'/,FMTF(2)
DIMENSION NARG(1),X(1),XX(7),XXA(7)
LCGICAL*1 TXT(41),LNAME(1),LR
INTEGER*2 AR,BL/' ',AREQ
EQUIVALENCE(NAM,LNAME(1)),(AR,LR),(FMTF(1),FMT2),(FMTF(2),FMT3)
NREW=0
NR=1
NAM=NAMES(1)
AR=BL
LR=LNAME(1)
CALL CONVY(LNAME(2),IR,3HI5 ,LHI)
WRITE(6,699) IR,AR
699 FORMAT(/T10,'IR=',I5,' AR=',A2)
  IFIL=NARG(4)
  IFOR=NARG(5)
  WRITE(6,698) IFIL,IFOR
698 FCRMAT(T10,'IFIL=',I2,' IFOR=',I2)
  BACKSPACE IFIL
  IF(NARG(1).GE.3) BACKSPACE IFIL
  FMT(2)=FMTF(IFOR)
  READ(IFIL,FMT,END=10) TXT,AREQ,IREQ,XX
  IF(AR.NE.AREQ.OR.IR.NE.IREQ) GOTO 15
  IF(NARG(1).LT.3.OR.XX(1).GT.NAMES(3)) GOTO 10
  GOTO 16

```

10  
20  
30  
40  
50  
60  
70  
80  
90  
100  
110  
120  
130  
140  
150  
160  
170  
180  
190  
200  
210  
220  
230  
240  
250  
260  
270

```

SUBROUTINE NEUNXT(NR,NARG,NAMES,X)
REAL*8 NAMES(1),REQ,FMT1(2)/'(T42,A6,' ',7E12.5) '/
IFMT2(2)/'(T42,A6,' ',7E12.5) '/
DIMENSION NARG(1),X(1),XX(7)
IFIL=NARG(4)
NR=1
IF(NARG(5).EQ.2) GOTO 5
READ(IFIL,FMT1,END=1001) REQ,XX
GOTO 10
5 READ(IFIL,FMT2,END=1001) REQ,XX
10 IF(REQ.NE.NAMES(1)) GOTO 1002
  X(1)=XX(1)
  J=2
  IF(NARG(2).NE.2) GOTO 22
  X(2)=XX(2)
  J=3
22 X(J)=XX(6)
  IF(NARG(3).NE.2) GOTO 24
  J=J+1
  X(J)=XX(7)
24 RETURN
1001 CONTINUE
1002 CCNTINUE
  NR=0
  RETURN
END

```

10  
20  
30  
40  
50  
60  
70  
80  
90  
100  
110  
120  
130  
140  
150  
160  
170  
180  
190  
200  
210  
220  
230  
240  
250  
260

		I=0	520
		NAMZ=NARG(1)	530
		IF(NAMZ.LE.2) GOTO 21	540
		DO 19 J=3,NAMZ	550
		19 NAMS(V(J-2))=NAMES(J)	560
		GOTO 21	570
C	SUBROUTINE RETXS(NARG, NAMES, EMIN, EMAX, X, Y, NUMX, MAXNUM, NR, LDFLOC, LD	1C	580
C	IFNXT)	20	590
C	COMMON/SIGSAV/Z	30	600
C	DIMENSION X(1),Y(1),Z(2),NARG(1),W(2)	40	610
C	REAL*8 NAMES(1),NAMSV(4)	50	620
RETXS RETRIEVES KEDAK-DATA.		60	630
		70	640
ASSIGN 20 TO NST		80	650
I=0		90	660
CALL LDFLOC(NERR,NARG,NAMES,Z)		100	670
IF(NERR.EQ.0) GOTO 30		110	680
NAMZ=NARG(1)		120	690
IF(NAMZ.LE.2) GOTO 3		130	700
DO 2 J=3,NAMZ		140	710
2	NAMSV(J-2)=NAMES(J)	150	720
3	IF(Z(1).LE.EMIN) GOTO 5	160	730
IF(Z(1).GE.EMAX) GOTO 32		170	740
GOTO 21		180	750
5	CALL LDFNXT(NERR,NARG,NAMES,W)	190	760
IF(NERR.EQ.0) GOTO 23		200	770
IF(W(1).LE.EMIN) GOTO 10		210	780
IF(W(1).GE.EMAX) GOTO 36		220	790
7	I=I+1	230	800
X(I)=Z(1)		240	810
Y(I)=Z(2)		250	820
I=I+1		260	830
X(I)=W(1)		270	840
Y(I)=W(2)		280	850
GOTO NST,(20,200)		290	860
10	CALL LDFNXT(NERR,NARG,NAMES,Z)	300	870
IF(NERR.EQ.0) GOTO 26		310	880
IF(Z(1).LE.EMIN) GOTO 5		320	890
IF(Z(1).GE.EMAX) GOTO 38		330	900
11	I=I+1	340	910
X(I)=W(1)		350	920
Y(I)=W(2)		360	930
12	I=I+1	370	940
X(I)=Z(1)		380	950
Y(I)=Z(2)		390	960
GOTO NST,(20,200)		400	970
20	CALL LDFNXT(NERR,NARG,NAMES,Z)	410	980
IF(NERR.EQ.0) GOTO 22		420	990
IF(Z(1).GE.EMAX) GOTO 24		430	1000
IF(I.EQ.MAXNUM) GOTO 34		440	1010
21	I=I+1	450	1020
X(I)=Z(1)		460	
Y(I)=Z(2)		470	
GOTO 20		480	
		490	
C	ENTRY REPSX(NARG, NAMES, EMIN, EMAX, X, Y, NUMX, MAXNUM, NR)	500	
		510	
		520	
		530	
		540	
		550	
		560	
		570	
		580	
		590	
		600	
		610	
		620	
		630	
		640	
		650	
		660	
		670	
		680	
		690	
		700	
		710	
		720	
		730	
		740	
		750	
		760	
		770	
		780	
		790	
		800	
		810	
		820	
		830	
		840	
		850	
		860	
		870	
		880	
		890	
		900	
		910	
		920	
		930	
		940	
		950	
		960	
		970	
		980	
		990	
		1000	
		1010	
		1020	

	SUBROUTINE LOCXS (NARG, NAMES, X, Y, NUMX, MAXNUM, NR, LDFLOC, LDFNXT)	10	10 M=M-1	120
	COMMON/SIGSAV/ Z	20	IF(M.EQ.0) GOTO 20	130
	DIMENSION X(1), Y(1), Z(2), NARG(1)	30	IF(R.LT.FELD(M)) GOTO 10	140
	REAL*8 NAMES(1), NAMS(4)	40	IA=M+1	150
C		50	IE=K-1	160
C	LOCXS : ARE RETRIEVAL SUBROUTINES	60	J=IE	170
C	NXTXS : FOR KEDAK DATA.	70	DO 30 I=IA, IE	180
C		80	FELD(J+1)=FELD(J)	190
	I=0	90	WERT(J+1)=WERT(J)	200
	CALL LDFLOC(NERR, NARG, NAMES, Z)	100	J=J-1	210
	IF(NERR.EQ.0) GOTO 30	110	FELD(IA)=R	220
19	NAMZ=NARG(1)	120	WERT(IA)=S	230
	IF(NAMZ.LE.2) GOTO 21	130	99 CONTINUE	240
	DO 10 J=3, NAMZ	140	100 RETURN	250
10	NAMS(J-2)=NAMES(J)	150	END	260
	GOTO 21	160		
20	CALL LDFNXT(NERR, NARG, NAMES, Z)	170		
	IF(NERR.EQ.0) GOTO 22	180		
	IF(I.EQ.MAXNUM) GOTO 34	190		
21	I=I+1	200		
	X(I)=Z(1)	210	SUBROUTINE EQUEN(IMAX, E, S)	10
	Y(I)=Z(2)	220	DIMENSION E(2), S(2)	20
	GOTO 20	230	DATA KOUT/6/	30
		240		40
C	ENTRY NXTXS(NARG, NAMES, X, Y, NUMX, MAXNUM, NR)	250	C EQUEN BEHAELT VON MEHRFACH VORKOMMENDEN E-WERTEN NUR DEN JEWEILS	50
	I=0	260	LETZTFN.E MUSS GEORDNET SEIN.	60
	GOTO 19	270		70
		280		80
C		290	IF(IMAX.LT.2) GOTO 20	90
22	NR=1	300	M=0	100
	IF(NAMZ.LE.2) GOTO 200	310	DO 10 I=2, IMAX	110
	DC 23 J=3, NAMZ	320	IF(E(I).NE.E(I-1)) GOTO 10	120
23	NAMES(J)=NAMS(J-2)	330	WRITE(KOUT,600) E(I), E(I-1), S(I-1)	130
	GOTO 200	340	600 FORMAT(20X, 'ZWEI GLEICHE ENERGIEWERTE:', E12.5, ' ERSTES WERTEPAAR W	140
30	NR=3	350	1IRO AUSGESONDERT:', 2E12.5)	150
	GOTO 200	360	M=1	160
34	NR=2	370	E(I-1)=E(I)	170
	GOTO 200	380	S(I-1)=S(I)	180
200	NUMX=I	390	L=I+1	190
	RETURN	400	GOTO 11	200
	END		10 CONTINUE	210
			GOTO 20	220
			11 DO 15 I=L, IMAX	230
			IF(E(I).NE.E(I-1)) GOTO 12	240
			WRITE(KOUT,600) E(I), E(I-1), S(I-1)	250
			M=M+1	260
			12 E(I-M)=E(I)	270
			S(I-M)=S(I)	280
			15 CONTINUE	290
			IMAX=IMAX-M	300
			20 RETURN	310
			END	
	SUBROUTINE ORDNEN(KMAX, FELD, WERT)	10		
	DIMENSION FELD(KMAX), WERT(KMAX)	20		
C		30		
C	ORDNEN ORDNET (FELD, WERT) NACH WACHSENDEN ARGUMENTEN VON FELD	40		
C		50		
	IF(KMAX.LT.2) GOTO 100	60		
	DC 99 K=2, KMAX	70		
	M=K-1	80		
	IF(FELD(K).GE.FELD(M)) GOTO 99	90		
	R=FELD(K)	100		
	S=WERT(K)	110		

SUBROUTINE SPACE(X,LX,NMAX)	10
DIMENSION X(1)	20
C---GET WORKING AREA FOR AVERAGE.	30
REAL*8 DATE,ZEIT	40
COMMON/INOUT/ KOUT	50
CALL FREESP(KB)	60
C===RESERVE 30 % BUFFER AREA, AT MAXIMUM 40KBYTE.	70
LBUF=(3*KB)/10	80
IF(LBUF.GT.40) LBUF=40	90
KF=KB-LBUF	100
C===EACH KBYTE CAN ACCOMODATE FOR 128 DATA POINTS.	110
NMAX=KF*128	120
NBYTE=KF*1024	130
CALL XTAREA(K1,NBYTE,K3,X)	140
IF(NBYTE.EQ.0) GOTO 100	150
LX=(K1-K3)/4+1	160
CALL DATUM(DATE,ZEIT)	170
WRITE(KOUT,601) DATE,ZEIT,LBUF,NMAX,K3	180
601 FCRMAT(1H1//////////2(/10X,70(' '*'))/	190
120X,'YOU ARE USING PROGRAMME AVERAGE'//	200
220X,'DATE:',A8/20X,'TIME:',A8/20X,'YOUR BUFFER AREA IS ',I3,	210
3' KBYTES'/20X,'YOUR WORKING AREA ',I6,' DATAPPOINTS, ALLOCATED AT ',	220
4Z9/2(/10X,70(' '*')))	230
RETURN	240
100 WRITE(KOUT,602)	250
602 FORMAT(//' WHEN ATTEMPTING STORAGE ALLOCATION'/	260
1' AN ERROR OCCURED IN ROUTINE "SPACE".')	270
STOP	280
END	290

3.2 Calculation of composite cross sections (CALCUL)

Contents:

3.2.1 Short introduction

3.2.2 Computer program abstract

### 3.2.1 Short introduction

Composite cross sections are to be understood here as quantities obtained by performing arithmetic operations (addition, subtraction, division etc.) on particular partial cross sections. Important examples are

- the total cross section, i.e. the sum of all partial cross sections.

$$\sigma_T = \sum_x \sigma_x,$$

- the absorption cross section

$$\sigma_a = \sigma_\gamma + \sigma_f + \sigma_p + \sigma_\alpha,$$

- the nonelastic cross section

$$\sigma_{\text{nonel}} = \sigma_T - \sigma_n,$$

- the capture-to-fission ratio

$$\alpha = \sigma_\gamma : \sigma_f.$$

The program CALCUL was written as a tool for the calculation of these and other composite cross sections from angle-integrated partial cross sections. The following Computer Abstract describes its main features. A detailed description will be published as a separate KFK report.

1. NAME OR DESIGNATION OF PROGRAMME: CALCUL is a program system to calculate cross section quantities. To insert into the KEDAK library neutron cross sections data evaluated elsewhere it is necessary to (re-)calculate them and arrange them according to KEDAK conventions.
2. COMPUTER FOR WHICH THE PROGRAMME IS DESIGNED AND OTHERS UPON WHICH IT IS OPERABLE: IBM/370-168 with TSO, MVT
3. NATURE OF PHYSICAL PROBLEM SOLVED: CALCUL processes data from an external source and/or data from the KEDAK library. A control input in form of a command language controls in which manner the data are processed. The output is a data set consisting of ADD records and DROP records that tell the KEDAK Management program which data are to be added to the library and which data are to be deleted from the KEDAK library.
4. METHOD OF SOLUTION: CALCUL assists the user in the general problem of exercising the basic mathematical operations upon functions depending on one variable (energy). Thus CALCUL simulates a desk calculator operating on functions instead of single numerical values; if necessary interpolation is performed.

The operations:

1. operation "+"                    ADD command  
    $R = R + y$                     R - the current result, is equal to zero  
  at the start of the calculation  
  R,y are data arrays
2. operation "-"                    SUBTRACT command  
    $R = R - y$
3. operation "."                    MULTIPLY command  
    $R = R * y$
4. operation "/"                    DIVIDE command  
    $R = R / y$
5. operation                        ETA calculation  
    $R = 1/(1+R)$



Furthermore the additional facility to perform the calculation of composed cross sections is provided in CALCUL. The composed cross sections are calculated with the aid of the single operations but the control in this case is performed by a subprogram instead of the control input for the single commands.

The commands and formulas for calculating of the composed cross section values:

ALPHA1	$\alpha = \sigma_{\gamma} / \sigma_f$
ALPHA2	$\alpha = (\bar{v} / \eta) - 1$
ETA1	$\eta = \bar{v} / (1 + \alpha)$
ETA2	$\eta = \bar{v} * \sigma_f / (\sigma_f + \sigma_{\gamma})$
SGA1	$\sigma_a = \sigma_{\gamma} + \sigma_f + \sigma_p + \sigma_{\alpha}$
SGG1	$\sigma_{\gamma} = \sigma_f * \alpha$
SGG2	$\sigma_{\gamma} = \sigma_f * ((\bar{v} / \eta) - 1)$
SGG3	$\sigma_{\gamma} = \sigma_a - \sigma_f - \sigma_p - \sigma_{\alpha}$
SGI1	$\sigma'_n = \sigma_x - \sigma_{\gamma} - \sigma_p - \sigma_{2n} - \sigma_{\alpha} - \sigma_{3n} - \sigma_f$
SGN1	$\sigma_n = \sigma_T - \sigma_x$
SGT1	$\sigma_T = \sigma_n + \sigma_x$
SGTR1	$\sigma_{trans} = \sigma_T - \sigma_n * \mu_L$
SGX1	$\sigma_x = \sigma_T - \sigma_n$
SGX2	$\sigma_x = \sigma_n + \sigma_{\gamma} + \sigma_f + \sigma_p + \sigma_{\alpha} + \sigma_{2n} + \sigma_{3n}$

5. RESTRICTIONS ON THE COMPLEXITY OF THE PROBLEM: KEDAK library format is prerequisite. The data from the external source are expected in formatted records, each including one energy value and one cross section value. Storage allocation is handled dynamically with the aid of the XTAREA subroutine (see reference 2). No restriction on number of energy mesh points. If the core storage given is not sufficient, CALCUL has the possibility to evade on disk storage (must be made available to the program by the user through DD-statements in the job control language).
6. TYPICAL RUNNING TIME: The running time is dependent on the number of data points (energy points) processed and on the region made available to the program, usually about 1 ms/data point on IBM/370-168.

7. UNUSUAL FEATURES OF THE PROGRAMME: CALCUL could be processed as a background job in batch processing or as a foreground job on a terminal in TSO.

8. RELATED AND AUXILIARY PROGRAMMES: CALCUL is written in modular form with many subroutines to facilitate the addition of new options and to permit the generation of a compact overlay structure.

The subroutines of CALCUL are comprised in the following modules:

1. Control module
2. Operation code definition package - OPDEF
3. Control input processing package - PROCINP
4. Calculation package - CALPAC
5. Cross section formula calculation package - CROSSEC
6. Data management of the auxiliary direct access data set - DATAMAN
7. Output editing package (KEMA-input-format data set) - OUTPUT
8. Processing control input in card image format - CARDIM

The retrieval packages RETPAC and LDFPAC are used to read the data from the KEDAK library. The following subroutines are external references:

- XTAREA and FREESP - to handle the dynamic storage allocation
- CONVY and STRING - to convert floating point and integer data to alphameric representation and v.v.
- DEFI and DINF - to make the DEFINE FILE statement dynamical
- DDTEST - to test for DD-cards

9. STATUS: In use

10. REFERENCES:

1. I. Langner, R. Meyer  
CALCUL - a program system to calculate cross section quantities and arrange data for input to KEMA.  
KFK, to be published.
2. I. Langner, R. Meyer  
IDFPAC/LDFPAC - two retrieval packages for the Karlsruhe Evaluated Nuclear Data Library, KFK, to be published
3. W. Höbel  
XTAREA, REXTAR - dynamische Dimensionierung von FORTRAN-Feldern, KFK, to be published

4. G. Arnecke, H. Bachmann  
DEFI, DINF dynamisches DEFINE FILE, KFK, to be published
5. G. Arnecke  
DDTEST - Benutzte Dateien, KFK, to be published
6. B. Krieg  
Handling and service programs for the Karlsruhe Nuclear  
Data File KEDAK, KFK 1725, June 1973
7. R. Meyer  
RETPAC - A user oriented retrieval package for use with  
the Evaluated Nuclear Data Library KEDAK, GfK, Internal Report
8. H. Blesene  
CONVY - FORTRAN-Unterprogramm für die IBM/360 zur Umwandlung  
von in maschineninterner bzw. in alphanumerischer Darstellung  
vorliegenden Fest- und Gleitkommazahlen in alphanumerische bzw.  
maschineninterne Darstellung, Internal Report, GfK, 1971
  
11. MACHINE REQUIREMENTS: Disk storage on a direct-access device;  
TSO terminal. Without overlay structure the program requires  
240 K core storage.
  
12. PROGRAMMING LANGUAGE USED: FORTRAN IV
  
13. OPERATING SYSTEM OR MONITOR UNDER WHICH PROGRAMME IS EXECUTED:  
OS 360/370 MVT and TSO
  
14. ANY OTHER PROGRAMMING OR OPERATING INFORMATIONS OR RESTRICTIONS:  
With the overlay structure CALCUL requires 130 K bytes core storage.  
The storage in the GO-step depends on the number of data points  
processed. If not enough core storage is available for the temporary  
data sets, the program tries to evade on auxiliary data sets on disk.
  
15. NAME AND ESTABLISHMENT OF AUTHOR:  
I. Langner, R. Meyer  
Gesellschaft für Kernforschung mbH  
Institut für Neutronenphysik und Reaktortechnik  
Postfach 3640  
D-7500 Karlsruhe 1
  
16. MATERIAL AVAILABLE: Source deck. List of available commands, and  
of conventions, and syntax of the control input list for a command.

4. PROGRAMS FOR WEIGHTING AND SMOOTHING OF CROSS SECTION DATA

4.1 Program SIGPLO

Contents:

4.1.1 General characteristics

4.1.2 Data retrieval

4.1.2.1 Retrieval from files in KEDAK format

4.1.2.2 Retrieval from NEUDADA tapes

4.1.2.3 Retrieval of other formatted data

4.1.3 Processing of retrieved data

4.1.3.1 Linear and spline interpolation

4.1.3.2 Curve fitting (smoothing)

4.1.3.3 Fission spectrum averages

4.1.4 Plot production

4.1.4.1 Plotter hardware and software employed by SIGPLO

4.1.4.2 Plot size and scale specification

4.1.4.3 Plot options

4.1.5 Output

4.1.5.1 Output options

4.1.5.2 Specification of new grid for output

4.1.6 Summary table

4.1.7 References

Appendix: SIGPLO Program List

## 4.1 Program SIGPLO

The following program description is based on an unfinished internal report by R. Meyer, the author of the program /1/.

### 4.1.1 General characteristics

Smoothing and fitting, interpolation, tabulation and plotting of numeric data are constantly recurring tasks in nuclear data evaluation. The very large data sets typical for modern neutron data work make utilization of computer methods mandatory. The program SIGPLO was developed in response to this need. The options offered by SIGPLO will be discussed in the following order:

- (1) data retrieval
- (2) processing of retrieved data (interpolation, smoothing)
- (3) output (printing, plotting, punching,...)
- (4) special features

The various options and their default values are summarized for quick reference in Table 1 below.

### 4.1.2 Data retrieval

The SIGPLO user may address three different sources of data:

- (1) data in KEDAK format,
- (2) NEUDADA "transmission" tapes,
- (3) other formatted (FORTRAN-readable) data.

Each set of data thus retrieved forms the "current data set" on which SIGPLO will operate. The current set may be extended by merging it with data from the same or other sources to form a new "current data set". Due to the use of fixed dimensions the size of the "current data set" must not exceed 1000 data points. Data need not be sorted - they are sorted after retrieval, if necessary. Every time a data set is read it either replaces the "current set" or, if merging is requested, it is merged with the "current set". In both cases the old "current set" is lost and cannot be referenced but by retrieving it again.

The examples given in Sect. 4.1.2 are valid for IBM computer installations working under OS and ASP.

#### 4.1.2.1 Retrieval from files in KEDAK format

Data in KEDAK format may be retrieved from up to five different KEDAK-type data libraries, so that for instance several versions of KEDAK can be compared with each other or with data translated from other libraries, e.g. ENDF/B. Retrieval from the KEDAK file is performed with the program package LDFPAC /2/.

In order to address a KEDAK-type library the user must specify the name of the library and the KEDAK name (key) of the data (cf. /3/). He must provide

- (1) a SETUP card for the device that contains the library, as required by IBM-ASP, if necessary;
- (2) DD control card(s) describing the library data set, as specified in the IBM-OS job control language manual.

Examples (underscores denote blanks):

```
//G.FT01FO01_DD_UNIT=2313,VØL=SER=GFK050,DISP=SHR
//          DSN=KEDAK3
```

will address the library KEDAK3 on the disk pack named GFK050. The data will be available to SIGPLO via the ft-number PLEH=1, cf. (3) below;

```
//G.FT02FO01_DD_UNIT=3330,VØL=SER=NUSICE,DISP=SHR,
//          DSN=STRUMAT.ENDFB3
```

will address the data library STRUMAT.ENDFB3 on the disc pack NUSICE; data will be available to SIGPLO through PLEH=2.

- (3) program control input describing names and energy range of the data that are to be retrieved:

```
_ &KEDAK_PLEH=...,NAMZ=...,NAMEN=...,TØ=...,FRØM=...,
_data-processing and output options,&END
```

where PLEH is the input device number (ft-number),

NAMZ is the number of KEDAK names necessary to address the requested data (default option: NAMZ=2), cf. /3/,

NAMEN are these KEDAK names, e.g. for  $\sigma_T$  of Na-23 one would write NAMEN='NA23', 'SGT' or alternatively NAMEN(1)='NA23', NAMEN(2)='SGT'. The KEDAK names must be written in the same sequence as in /3/;

FROM     minimum energy,  
TO        maximum energy.

The data-processing and output options are described in Sects. 4.1.3 - 4.1.5 below.

#### 4.1.2.2 Retrieval from NEUDADA tapes

NEUDADA tapes contain experimental data from the common data base of the "four-centre network" of neutron data centres (National Neutron Cross Section Center, Brookhaven, USA; Centre de Compilation de Données Nucléaires of NEA/OECD, Saclay, France; Nuclear Data Section of IAEA, Vienna, Austria; Centr po Jadernym Dannym, Obninsk, USSR). These (7 or 9 track) tapes are sent by CCDN to users in non-American OECD countries upon request.

In the NEUDADA formats (neutron data under direct access) each data point is coded in one record together with its uncertainty, other related data and a data set identifier. These identifiers (accession numbers) are used by SIGPLO to distinguish the various data sets that may occur on one tape. The user must

- know on which file his data actually reside (tape marks written at CCDN are not consistent with all IBM operating systems) and whether they are written in CCDN "transmission" format or in CCDN internal format;
- check whether his data have got unique accession numbers or whether they are distributed among several accession numbers, which frequently may happen. In view of this possibility SIGPLO offers a special option to combine several sets of data;
- make sure the requested data are complete. If this is not the case SIGPLO allows to add data e.g. on punched cards. At present SIGPLO offers no possibility to delete data points read. The only way to correct faulty data points is to copy-correct the respective data set on the tape.

In order to retrieve data from NEUDADA tapes the user must provide

- (1) a SETUP card as required by IBM-ASP;

- (2) a DD control card for each file to be addressed, as specified in the IBM-OS job control language manual.

Examples:

```
//G.FT10FOO1_DD_UNIT=TAPE9,VOL=SER=ND0501,  
//          LABEL=(2,NL,,IN),DISP=(OLD,PASS),  
//          DCB=(BLKSIZE=1320,LRECL=132)
```

will address data on file 2 of the 9-track tape named ND0501. Data will be available to SIGPLO via ft-number BDEH=10, cf. (3) below;

```
//G.FT04FOO1_DD_UNIT=TAPE7,VOL=SER=ND2354,  
//          LABEL=(2,NL,,IN),DISP=(OLD,PASS),  
//          DCB=(BLKSIZE=1320,LRECL=132,TRTCH=ET)
```

will address data on file 2 of the 7-track tape named ND2354. Data will be available to SIGPLO via BDEH=4, cf. (3) below. The data on these two tapes are available to SIGPLO simultaneously, two tape drives are required.

- (3) Program control input:

```
&NEUDAD BDEH=...,LTR=...,ACCNR=...,SEQN=...,FØDATA=...,  
_FRØM=...,TØ=...,data-processing and output options,&END
```

where

BDEH gives the FT number of the tape unit

LTR specifies the letter, if any, that precedes the accession number supplied by CCDN, e.g. LTR='E' if data set E00012 is desired. The default option is LTR='O', hence LTR need not be specified for data set 000012.

ACCNR is the (numeric) accession number supplied by CCDN, e.g. ACCNR=12 for data set E00012.

SEQN specifies whether the accession number is in line with the expected ascending order on tape (SEQN=T) or not (SEQN=F), so that the tape can be rewound where necessary.

Example

Suppose the sequence of data sets on tape is

```
...U00003,E00002,...E00013,E00014,...U00015,...
```



whereas they are to be retrieved in the order

E00013,E00014,U00015,E00002,U00003.

In this case all sets can be retrieved with SEQN=T except U00003 for which SEQN=F signals that it occurs on tape not in the expected order (i.e. after E00002).

FØDATA indicates whether the NEUDADA tape is written in CCDN "transmission" format (FØDATA=T, default option) or in CCDN internal format (FØDATA=F).

FRØM,TØ are minimum and maximum energy.

Other data-processing and output options are explained below.

#### 4.1.2.3 Retrieval of other formatted data

Formatted input data in the present context means FORTRAN-readable data (as explained e.g. in the IBM FORTRAN IV manuals) with the following restrictions:

- Each set of numeric data records (data set) must be preceded by a comment record with arbitrary free text. This text is read and printed so that the user can check the input sequence.
- Each data point must be completely contained in one record. On the other hand one record may contain several data points.
- The co-ordinates of each data point must be represented in the order  
(1) energy, (2) cross section (or other energy-dependent quantity)  
(3) uncertainty (optional). Other information may be interspersed with these data. The user must specify the reading format accordingly, e.g. by spacing over characters not to be read as data.

Records may be cards or card images on tape or other rigidly structured character sequences on any input device.

For retrieval of a formatted data set the user must provide

- (1) a setup card for the device from which the data are to be read, if necessary,
- (2) a DD card describing the device and the data set.

Examples

```
//G.FT16FOO1_DD_UNIT=2314,VØL=SER=NUSYSO,  
//          DISP=SHR,DSN=FØRM,INP1
```

would provide formatted data with unit reference number (ft-number) 16. Data are read from a disc named NUSYSO and data set FØRM.INP1. Data set characteristics need not be given, they are provided by the operating system.

```
//G.FT11FOO1_DD_UNIT=TAPE9,VØL=SER=TAPE72,  
//          DISP=(ØLD,PASS),LABEL=(1,NL),  
//          DCB=(RECFM=FB,LRECL=30,BLKSIZE=3000)
```

specifies file 1 of the unlabelled 9-track tape named TAPE72 as input. Since the tape has no label, data set characteristics must be supplied.

- (3) program control input describing device, format, energy range and other characteristics of the data set:

```
_&KARTEN_KARTEH=...,FMT=...,ANZ=...,  
_FRØM=...,TØ=...,EINH TX=...,EINH TY=...,  
_REWD=...,data-processing and output options,&END
```

where

KARTEH is the input device reference number. If punched cards are supplied together with control input, KARTEH=9 must be specified (not KARTEH=5) and the cards (preceded by the comment card) must immediately follow the respective control card.

FMT gives the reading format, e.g.

FMT='(15X,E15.5,10X,E15.5)' or

FMT='(8F10.5)'.

ANZ determines the number of data points to be read (as distinct from points actually utilized) as follows:

The requested data set is searched for the first record beginning with an energy  $\geq$  FROM\*0.999999. If this is the first record of the whole data set, reading starts there, otherwise reading starts at the preceding record (because that may contain, in addition to the first energy, other unchecked energies that possibly exceed the limit FROM\*0.999999). Reading stops as soon as ANZ data points are read or an END OF FILE condition occurs. Subsequently SIGPLO searches for the first data point read with energy exceeding T0\*1.000001 and deletes it as well as all subsequent points. Then all data points with energies less than FROM\*0.999999 are discarded. Finally the remaining data points are rearranged in order of ascending energy. Note that ANZ determines the number of data points read, the actual number of data points utilized by SIGPLO can be smaller due to the deletion of points with energies outside the requested range. Since an exact determination of the number of data points to be read may be inconvenient, the author suggests that the user

either supplies only one data point per record, in which case exactly ANZ cards (= data points) are read starting with the card preceding the first data point with energy  $\geq$  FROM, counted from the current position within the data set,

or matches ANZ exactly with the number of data to be read, if the data set contains only points within the given energy limits,

or uses a different unit reference number for each data set, utilizing the END OF FILE condition as a natural delimiter.

The example below illustrates this.

FROM,T0 have their usual meaning of minimum and maximum energy.

EINH TX =  $\left\{ \begin{array}{l} \text{'EV'} \\ \text{'KEV'} \\ \text{'MEV'} \end{array} \right\}$  specifies energy units,

EINH TY =  $\left\{ \begin{array}{l} \text{'MB'} \\ \text{'BARN'} \end{array} \right\}$  specifies cross section units.

Data will always be converted to eV or b internally. Of course, conversion could be obviated by proper scale factors in the FMT input format specification.

REWD may be specified as T (or F) which causes rewinding (or not) before retrieval starts (with the exception of the control input device for which REWD is ignored).

Warning: - ANZ should not exceed 1000.

- Although input data need not be sorted, the user should be aware of the fact that they are sorted only after they have been read. Hence data after the first energy  $\geq T0$  will not be transmitted even though the respective card(s) might have been read.

Example:

```
//G.SYSIN_DD_*
RT_TEST_INPUT
  &KARTEN_KARTEH=9,FMT='(10F10.2)',ANZ=2,
  FROM=0.,TO=100.,&END
C INPUT FROM CONTROL UNIT
1.      0.37   24.2   0.81
  &KARTEN_KARTEH=14,FMT='(2F10.2)',ANZ=50,
  FROM=24.6,TØ=82.9,&END
  &KARTEN_KARTEH=11,FMT='(2F10.2)',ANZ=50,
  FROM=24.6,TØ=82.9,&END
//G.FT11FOO1_DD_*
C INPUT FROM UNIT 11
13.6      0.37
17.77     1.44
21.88     .           (first card read)
24.92     .
23.11     .
41.88     etc.
35.97     .
46.1      .
17.15     .
53.2      .
57.2      .
69.44     .
91.85     .
82.1      .           (last card read)
//G.FT14FOO1_DD_*
C INPUT FROM UNIT 14
25.99     .           (1st card read)
44.12     .
57.31     .
80.50     .           (last card read)
//
```

### Explanation

Let us for the moment disregard the card RT\_\_... which will be explained in Sect. 4.1.3.2 below. The first input data set is read from unit 9. The user took care that only one data card is read by specifying with ANZ=2 the exact number of data to be read. Thus two data points are transferred to SIGPLO:(1.,0.37) and (24.2,0.81).

The second data set is read from unit 11. Reading will start with energy 21.88 eV since this is the last one less than FROM. Reading will stop after energy 82.1 eV caused by the END OF FILE condition. The points with energies 21.88 eV and 23.11 eV are then deleted, these energies being less than FROM. Likewise the point with energy 91.85 eV (the first energy exceeding TØ) and all subsequent points are deleted. Thus the points available to SIGPLO after sorting are those with energies 24.92, 35.97, 41.88, 46.1, 53.2, 57.2, and 69.44 eV.

The third data set is transferred from unit 14 to SIGPLO without deletions.

Note that FMT, FROM and TØ could have been omitted from the last control cards since the previous values are the same and would be kept by default. This, however, does not hold for ANZ. (See Table 1, footnote). Since SIGPLO assumes the units eV and barn by default, EINHTX and EINHTY can be omitted.

The energy-sorting feature allows for easy addition of other data to a punched data set. For instance, if a point at 29.7 eV is to be added to the third set of data one could insert the corresponding card simply at the end of (or some other place within) the card deck containing the third data set.

Merging of data sets is possible with the option

$$\text{DATADD} = \begin{cases} \text{T} & \text{merge "current" set and newly retrieved set,} \\ \text{F} & \text{replace "current" set by newly retrieved set} \\ & \text{(default option)} \end{cases}$$

This option can be used in the control input lists &NEUDAD... or &KARTEN... specifying the newly retrieved set.

### 4.1.3 Processing of retrieved data

SIGPLO offers two possibilities to process retrieved data: interpolation and smooth-curve fitting (smoothing). The results can be plotted, printed and punched on cards (or written onto some other output medium).

#### 4.1.3.1 Linear and spline interpolation

The aim of interpolation is to run a curve through a given set of points known as "nodes" by specifying an interpolation law. The interpolation options of SIGPLO are most useful for drawing a continuous line through a set of data points on a plot or for retabulation on a different set of nodes ("new grid"). Thus initial data set (nodes), interpolation law and new grid must be specified:

- (1) The initial data set is defined by an appropriate retrieval as described in Sect. 4.1.2.
- (2) The interpolation law is defined by the key-word KURVE that may be employed in any of the three input lists &KEDAK..., &NEUDAD..., &KEDAK... (cf. Sect. 4.1.2). The options are

$$\text{KURVE} = \begin{cases} \text{'NEIN'} & \text{no interpolation (default option),} \\ \text{'LINE'} & \text{linear interpolation,} \\ \text{'SPLI'} & \text{cubic spline interpolation.} \end{cases}$$

Linear interpolation means that between adjacent nodes  $(x_i, y_i)$  and  $(x_{i+1}, y_{i+1})$  interpolated values  $(x, y)$  are calculated according to the interpolation law

$$\frac{y - y_i}{x - x_i} = \frac{y_{i+1} - y_i}{x_{i+1} - x_i} .$$

Cubic spline interpolation means that the interpolation law  $y = f(x)$  between each pair of adjacent nodes is a cubic, the different cubics being chosen such that  $f$ ,  $f'$  and  $f''$  are continuous at the matching points  $x_i$  ( $i=2,3,\dots,n-1$ ) and  $f'' = 0$  at the end points  $(x_1, x_n)$ .

It also means that the minimum number of nodes is 3.

(3) Output options are as follows:

- PLØKU =  $\begin{cases} T & \text{plot curve through original points (default option),} \\ F & \text{do not plot,} \end{cases}$
- DRUKU =  $\begin{cases} T & \text{print table of interpolated set, on new grid,} \\ F & \text{do not print (default option),} \end{cases}$
- STAKU =  $\begin{cases} T & \text{"punch" interpolated set, on new grid,} \\ F & \text{do not "punch" (default option)} \end{cases}$   
where punch means generation of any  
FORTRAN-readable output (on cards, tape, disc,...)

For the option PLOKU the new grid is automatically generated by SIGPLO. For DRUKU and STAKU, however, the new grid as well as the printing or "punching" format must be supplied by the user as explained in Sect. 4.1.3.3 below.

Notes:

- PLØKU=T causes plotting of only the initial set and belongs, strictly speaking, to Sect. 4.1.4 below. If the interpolated set would have to be plotted, it would have to be written onto some rereadable computer medium, reread with another "control input" list and plotted.
- If plotted output is desired the plot must be opened first by definition of its characteristics with a suitable "control input" list (cf. Sect. 4.1.4 below).
- The new grid must be defined before interpolation starts, i.e. the control input list defining the new grid must precede the control input list that requests interpolation.
- Similarly the control input list specifying the format for printing or other formatted-output production must precede the control input list that requests the output.
- If  $x_i$  are the initial grid points and  $x_j'$  the new grid points  $x_1 \leq x_1'$  and  $x_{\text{last}} \geq x_{\text{last}}'$  must hold.



- The linear interpolation and spline options are useful if cross sections are to be retabulated on a new or denser grid, e.g. for KEDAK. Splines, however, are sometimes quite sensitive to the exact position of the nodes and a smooth second derivative (curvature) may be difficult to achieve. For further details of spline fitting see /5/.
- If KEDAK cross sections are plotted, the user usually wants a curve without special marking of the nodes. This is achieved by the following combination of options:  

```
&KEDAK NAMEN='material', 'type', FROM=xx, TØ=xx,  
_KURVE='LINE', PLOKU=T, PLO=F, &END
```

  
(PLOKU=T requests the curve, PLO=F suppresses special marking of the nodes, cf. Sect. 4.1.4).

#### 4.1.3.2 Curve fitting (smoothing)

Curve fitting is possible with SIGPLO in the following restricted sense:

A curve  $y=f(x)$  is drawn through a given set of points  $(x_i, y_i)$ ,  $i=1,2,\dots,n$ , which is "as smooth as possible". The user sets a limit  $\chi^2$  which the weighted sum of squared deviations must not exceed:

$$\sum_{i=1}^n w_i [y_i - f(x_i)]^2 \leq \chi^2 \quad (1)$$

Within this limit the fitted curve is determined such that

$$\int_{x_1}^{x_n} f''(x)^2 dx = \min \quad (2)$$

Obviously this kind of fit (smoothing, ref. /6/) requires caution. Of course, one can produce a smooth curve through violently fluctuating data points, but if the fluctuations are genuine the resultant smooth curve is quite unrealistic. Smoothing of discrepant data sets from different experiments can be very difficult. The result will often not be better than an eyeguide curve, the only advantage being that one need not read off numerical values which the computer produces easily. This must be kept in mind before smoothing is requested.

Since frequently a smooth curve is to be drawn through data from different measurements SIGPLO allows merging several input data sets into one set for smoothing. If, for instance, the data sets  $S_1, S_2, \dots$  are to be merged for joint smoothing, one has to indicate this by putting

$$SM\emptyset\emptyset = \chi_j^2$$

into the control input list of  $S_j$ , where the  $\chi_j^2$  are real numbers chosen in such a way that their sum is just the upper limit of the weighted sum of squared deviations in Eq. (1),  $\chi^2 = \sum_j \chi_j^2$ . The joint set  $S = S_1 + S_2 + \dots$  is then retained for final smoothing. Merging a retrieved set into this "retained set" can be prevented by using  $SM\emptyset\emptyset = 'NEIN'$  (default option) in its control input list. Smoothing is possible only for control input lists of type  $\&NEUDAD$  and  $\&KARTEN$ , but not for  $\&KEDAK$  ( $KEDAK$  point data are linearly interpolable and thus smooth by definition).

Smoothing starts as soon as

- the end of control input is reached or
- a card containing  $RT\_$  plus free text is found or
- a  $\&SMOWDH$  control input list is encountered.

An  $RT$  card will empty the retained set after smoothing, a  $\&SMOWDH$  list (explained in more detail below) will not.

The output options for the results of smoothing are as follows:

$$PL\emptyset SM = \begin{cases} T & \text{plot the fitted smooth curve (default option}^*) \\ F & \text{do not plot;} \end{cases}$$
$$DRUSM = \begin{cases} T & \text{print points of the fitted smooth curve,} \\ F & \text{do not print (default option}^*); \end{cases}$$
$$STASM = \begin{cases} T & \text{"punch" points of the fitted smooth curve,} \\ F & \text{do not punch (default option}^*) \end{cases}$$

where "punch" means generation of any  
FORTRAN-readable output (on cards, tape, disc,...)

---

<sup>\*</sup>) Note: The default options are in effect only if the options were not already defined in a preceding control input list!

These output options are not needed until smoothing actually starts. This means they could be coded e.g. within the last control input list that is read before smoothing begins. If plotting is requested (PLØSM=T) the plot must also be opened before smoothing starts. The sum total ( $\chi^2$ , Eq.(1)) of all SMØØ values will be displayed or not as specified by

$$PLØFQU = \begin{cases} T & \text{display the squared-error sum } \chi^2, \\ F & \text{do not display } \chi^2. \end{cases}$$

The weights  $w_i$  in Eq. (1) are specified as follows:

- (1) All points of the "retained" set get unit weight by default, if weights are not defined explicitly.
- (2) Weights are defined explicitly by retrieval of the experimental uncertainties  $\delta y_i$  associated with the data points  $(x_i, y_i)$ . The weights are thus calculated as  $w_i = (\delta y_i)^{-2}$ . The experimental uncertainties can be read together with the data, e.g. from the same tape, or separately. In any case reading of experimental errors must be requested in the same input list as reading of the data. One writes SMGEW=T together with specifications of the medium from which errors are to be read, and error units. The options are as follows:

$$\_ \&NEUDAD \_ \dots SMGEW=T, MEDIUM = \left\{ \begin{array}{l} 'TAPE' \\ 'CARD' \end{array} \right\}, DEINHT = \left\{ \begin{array}{l} 'PERC' \\ 'B' \\ 'MB' \end{array} \right\}, \dots$$

or

$$\_ \&KARTEN \_ \dots SMGEW=T, EINHTD = \left\{ \begin{array}{l} 'PERC' \\ 'B' \\ 'MB' \end{array} \right\}, \dots$$

In case MEDIUM='TAPE' the errors are read from tape together with the  $(x_i, y_i)$  values and the units are converted to barns according to the NEUDADA specifications on tape. Note that DEINHT and EINHTY (see Sect. 4.1.2.2) must then specify the same units. If these specifications are missing, the

user can override the default option by specifying the error units (PERC for percent errors, B or MB for errors in b or mb).

In case MEDIUM='CARD' one must define input unit, input format and number of data to be read by

KARTEH=...,FMT=...,ZAHL=...

as explained in Sect. 4.1.2.3. If ZAHL is less than the number of data points read from tape the last error value read will be applied to all remaining data points. This allows, for example, to give only one error bar if the same error is to be assigned to all data points. In case data points and errors are read from the same medium (same deck of cards, same tape etc.) they must be in the order  $\dots, x_i, y_i, \delta y_i \dots$

Reading of data errors can be suppressed by SMGEW=F.

A user who wishes to modify weights differently in different energy regions can do so with the control input list

&RELGEW\_ZAHLXR=...,XR=...,PR=...,EINHTX=...,&END

where ZAHLXR is an integer  $n \leq 20$  giving the number of modification factors (energy regions);

XR is a list of  $n$  real numbers, separated by commas, giving the lower limits of the energy regions;

PR is a list of  $n$  real numbers, separated by commas, giving the modification factors;

EINHTX =  $\left\{ \begin{array}{l} \text{'EV'} \\ \text{'KEV'} \\ \text{'MEV'} \end{array} \right\}$  indicates in which units the XR values are given.

Relative weights are then applied as follows: Weights of all data points in the "retained data set" are multiplied

- by 1 if  $x < XR(1)$ ,

- by PR(i) if  $XR(i) < x < XR(i+1)$ ,  $i = 1, 2, \dots, n$ .

The first attempt to fit a given set of data is often not very successful, smoothing being very much a trial and error method. Usually a number of fits with varying parameters must be performed before the result is satisfactory. SIGPLO offers a convenient possibility to repeat smoothing, varying only selected parameters given in the following control input list:

```
_&SMØWDH_FQU=...,ZÄHLXR=...,XR=...,PR=...,EINHTX=...,NEUZEI=...,  
_PLØFQU=...,&END
```

where FQU is a real number establishing a new total limit  $\chi^2$  for the weighted squared deviations,

ZÄHLXR,XR,PR,EINHTX may be used to define new weight modification factors as explained before,

NEUZEI =  $\begin{cases} T & \text{start a new plot with the current plot size parameters,} \\ F & \text{draw fitted curve on current plot} \\ & \text{(if plotting is requested)} \end{cases}$

PLØFQU =  $\begin{cases} T & \text{plot the FQU value } (\chi^2) \text{ also,} \\ F & \text{do not plot the FQU value.} \end{cases}$   
(PLØFQU=T can be used only if not more than one smooth curve is drawn onto one plot. Otherwise the numbers would overprint each other).

### Summary

The following control input options are related to smoothing:

(1) current data set options

```
(&NEUDAD,&KARTEN) SMØØ,PLØSM,DRUSM,STASM,SMGEW  
(&NEUDAD) MEDIUM,DEINHT,KARTEN,FMT,ZÄHL  
(&KARTEN) EINHTD
```

(2) weight modification

(&RELGEW) Z AHLXR, XR, PR, EINHTX

(3) repetition of smoothing

(&SMØWDH) FQU, Z AHLXR, XR, PR, EINHTX, NEUZEI, PLØFQU

Note: - The maximum number of data points in the "retained set" is 1000.

- Smoothing should be used only if the point scatter is expected to be statistical (random). Do not smooth if fluctuations are genuine, e.g. in the resonance range.

- If the point scatter is statistical one expects for a good fit

$$\sum_i w_i (f(x_i) - y_i)^2 = \chi^2 \approx N,$$

where N is the degree of freedom of the problem, i.e. the number of points minus the number of fit parameters. Hence N is recommended as a starting value for smoothing.

Example

```
RT__SMØØTH ØF TWØ SETS ØN TAPE
_&NEUDAD_BDEH=2, ACCNR=1, LTR='E', SMØØ=1.2, SMGEW=T, &END
_&PLØTGR_XMIN=0.25, XMAX=1.25, XLG=40, YMIN=0.1, YMAX=0.6, YLG=20., &END
_&NEUDAD_ACCNR=2, SMØØ=2, PLØFQU=T, &END
_&RELGEW_ZAHLXR=1, PR=0.86, &END
_&SMØWDM_FQU=1., NEUZEI=T, &END
_&SMØWDM_FQU=0.1, &END
_&SMØWDM_FQU=0.04, &END
RT__SMØØTH ANØTHER SET, PLØT ØUTPUT
_&PLØTGR_XMIN=2., XMAX=18., XLG=34., YMIN=1, YMAX=7, YLG=24., &END
_&NEUDAD_BDEH=2, ACCNR=5, LTR='U', SMØØ=1.22, SMGEW=T, DEINHT='PERC',
_MEDIUM='CARD', ZAHL=1, FMT='(F5.2)', KARTEH=5, STASM=T, &END
2.6
&STAFØR_STAFMT='(1P6E12.5)', &END
&ØUTINT_ZXINT='KEDA', NAMEN='U235', 'SGT', FROM=2., TØ=18., &END
/*
```

Explanation

The NEUDADA sets E00001 and E00002 are merged and all weights are multiplied by 0.86. The card `&PLØTGR...` defines the dimensions of a plot as explained in Sect. 4.1.4.2 below. Smoothing begins with  $\chi^2 = 1.2+2.0=3.2$  as soon as the first `&SMØWDM` card is encountered (before its content is read). The result is plotted by default. Then smoothing and plotting is repeated with  $\chi^2 = 1.$ , 0.1 and 0.04. For each  $\chi^2$ -value the "retained set" and the resulting smooth curve are plotted onto a new plot (with the initially given plot size specifications). The second RT card causes the retained set to be emptied, whereupon a new plot size is specified and a new data set, U00005, is read in. All error bars are then set equal to 2.6 %. A punching format is given (`&STAFØR...`) and the new grid for the punched output is specified (`&ØUTINT...`) as explained in Sect. 4.1.5.2 below. Smoothing of the last data set (U00005) begins when the end of the file is reached (`/*` card).

#### 4.1.3.3 Fission spectrum average

The SIGPLO program averages cross sections over a standard fission spectrum if the option FISS=T appears in the control input lists &KEDAK..., &NEUDAD... or &KARTEN... (FISS=F is the default option.) The spectrum is taken as

$$\chi(E)dE = 0.4527 \text{ MeV}^{-1} \exp\left(-\frac{E}{0.9650 \text{ MeV}}\right) \sinh \sqrt{\frac{E}{0.4367 \text{ MeV}}} dE$$

which corresponds to  $^{235}\text{U}$  fission at low incident energies. By changing a single statement in subroutine QFISS other spectra could be used. The result,

$$\langle \sigma \rangle_{\chi} = \int \sigma(E)\chi(E)dE$$

is printed as well as the mesh points used for the numerical integration.

#### 4.1.4 Plot production

##### 4.1.4.1 Plotter hardware and software employed by SIGPLO

SIGPLO utilizes the following plotting facilities which are available at Karlsruhe:

- an IBM 1130 plotter producing plots 10 in. high and
- a CALCOMP plotter producing plots 24 in. high.

Access to both devices is controlled by PLØTA /7/, a FORTRAN subroutine extensively used at Karlsruhe.

##### 4.1.4.2 Plot size and scale specification

The user should keep in mind the following rules:

- Only one plot can be produced at a time. This means that output cannot be routed to more than one plot at any given time. If two or more plots are required they must be produced sequentially.



- Before output can be routed to a plot, the size, scaling and other characteristics of the plot must have been established.

These rules can be pictured as the following temporal sequence:

```
[ define plot no. 1
  route output to plot no. 1
[ define plot no. 2
  route output to plot no. 2
  :
  etc.
```

Each plot must be opened by specification of its characteristics in a program control input list as follows:

```
&PLOTGR XMIN=...,XMAX=...,XLG=...,YMIN=...,YMAX=...,YLG=...,
_RAST=...,...,EINHTX=...,EINHTY=...,CALP=...,&END
```

XMIN,XMAX        define minimum and maximum value on the x-axis in the units given by EINHTX.

XLG              is the length of the x-axis in units of 0.4 in.=1.016 cm.

YMIN,YMAX        define minimum and maximum value on the y-axis in the units given by EINHTY.

YLG              is the length of the y-axis in units of 0.4 in.=1.016 cm.

$RAST = \left\{ \begin{array}{l} 'NEIN' \\ 'LINE' \\ 'XLØG' \end{array} \right\}, \left\{ \begin{array}{l} 'NEIN' \\ 'LINE' \\ 'YLØG' \end{array} \right\}$  specifies whether a grid is to be drawn and if so, for which kind of scale. Linear scaling (LINE) produces parallel lines 0.8 in.=2.032 cm apart, while log scaling produces a line at each integer multiple of the current power of ten.

$EINH\text{TX} = \begin{Bmatrix} \text{'EV'} \\ \text{'KEV'} \\ \text{'MEV'} \end{Bmatrix}$  gives the x units ,

$EINH\text{TY} = \begin{Bmatrix} \text{'BARN'} \\ \text{'MB'} \end{Bmatrix}$  gives the y units .

The plot will be drawn in units of eV and b unless the user does not request a conversion by stating EINH TX and EINH TY explicitly.

$CALP = \begin{Bmatrix} \text{T} \\ \text{F} \end{Bmatrix}$  specifies whether a CALCOMP plot is requested (T) or not (F) - in the latter case the IBM1130 plotter is used.

Note that

- $2.*(XMAX-XMIN)/XLG$  and  $2.*(YMAX-YMIN)/YLG$  should be smooth numbers (integers) to produce convenient scales;
- XLG and YLG should be even integers;
- it is not possible to produce logarithmic scales without grid, i.e. RAST='NEIN',... or RAST='...', 'NEIN' is automatically interpreted as meaning a linear scale.

#### 4.1.4.3 Plot options

Having opened a plot by specifying its characteristics in a `&PLØTGR` control input list the user can route output to it by using one of the following options:

$$PLØ = \begin{Bmatrix} \text{T} \\ \text{F} \end{Bmatrix} , \quad PLØKU = \begin{Bmatrix} \text{T} \\ \text{F} \end{Bmatrix} , \quad PLØSM = \begin{Bmatrix} \text{T} \\ \text{F} \end{Bmatrix} .$$

PLØ=T        yields a plot of the "current set" of points,  
 PLØKU=T      "    "    "    "    an interpolated curve and  
 PLØKU=T      "    "    "    "    a fitted smooth curve.

These options may occur in the input lists &KEDAK\_..., &NEUDAD\_..., &KARTEN\_... with the restriction that PLØSM is not allowed for &KEDAK\_. If data points (as distinct from curves) are to be plotted ten different point symbols are available. These symbols are linked to the integers 0,1,...9 as follows:

n	0	1	2	3	4	5	6	7	8	9
IBM 1130 symbol	○	□	◇	+	×	⊗	⊠	△	▽	•
CALCOMP symbol	⊔	⊕	⊖	+	×	◇	⊕	⊗	⊚	⊙

SIGPLO uses the following sequence by default:

9,0,3,4,5,6,7,1,8,2, then repeatedly 0,1,2,7,8.

Up to forty point sets can be plotted onto one plot. The user may overrule the default sequence with

PLOZEI = n,    where n is the integer associated with the desired symbol, and  
 PLOZEI = -1    causes return to the default sequence.

PLOZEI is permitted in any one of the three control input lists &KEDAK, &NEUDAD, &KARTEN.

Example

```
RT__FE-PLØTS
_&PLØTGR_YMIN=1.,YMAX=2.,YLG=10.,XMIN=1.E6,XMAX=5.E6,
_XLG=20.,RAST='LINE','LINE',&END
_&KEDAK_NAMEN='FE','SGN',FRØM=0.998E6,TØ=5.002E6,
_PLØ=F,KURVE='LINE',&END
_&KARTEN_KARTEH=12,ANZ=12,KURVE='NEIN',PLØ=T,
_PLØZEI=5,EINH TX='MEV',&END
_&NEUDAD_BDEH=4,PLØZEI=-1,ACCNR=4,LTR='E',&END
_&PLØTGR_YMIN=2.,YMAX=4.,YLG=20.,&END
_&KARTEN_KARTEH=13,ANZ=108,&END
_&NEUDAD_ACCNR=6,PLØZEI=8,&END
_&NEUDAD_ACCNR=8,PLØZEI=-1,&END
_&NEUDAD_ACCNR=7,LTR='U',&END
_&KEDAK_KURVE='LINE',PLØ=F,NAMEN='FE','SGT',&END
```

Explanation

Convenient scaling is achieved for both plots by the use of (2 cm  $\approx$ ) 0,8 in.  $\hat{=}$  0.2b and  $\hat{=}$  0.4 MeV, respectively. The first plot will contain one curve and two sets of points. The point symbol number 5 is explicitly requested for the first set, symbol number 0 is then assigned to the second set by default. The second plot will contain 4 sets of data points and one curve. The default sequence of point symbols is used except for the first set of points. Note that PLØ=F was used for all KEDAK data so that curves will be shown but not the nodes (the data points through which the curves are drawn). After the first occurrence of PLØ=F marking of points is switched on again by PLØ=T in the next program control input list. Extensive use is made of the default options: CALP=F is omitted; XMIN,XMAX,XLG,RAST are not specified for the second plot so the values of the first plot remain in effect; TØ,FRØM are specified only at the beginning and remain unchanged throughout the job.

#### 4.1.5 Output

A number of output options were mentioned already. The following section complements and summarizes the various possibilities offered by SIGPLO.

##### 4.1.5.1 Output\_options\_

The "current data set" is listed or not with

$$\text{DRUCK} = \begin{cases} \text{T} & : \text{ print} \\ \text{F} & : \text{ do not print (default option),} \end{cases}$$

it is punched (or written on tape or disc etc.) or not with

$$\text{STANZ} = \begin{cases} \text{T} & : \text{ punch} \\ \text{F} & : \text{ do not punch} \end{cases}$$

and it is plotted or not with

$$\text{PL}\emptyset = \begin{cases} \text{T} & : \text{ plot} \\ \text{F} & : \text{ do not plot.} \end{cases}$$

DRUCK, STANZ and PL $\emptyset$  can occur in any one of the control input lists &KEDAK..., &NEUDAD..., &KARTEN\_.... Standard formats are used for printing and punching. In the latter case the user can overrule the default option (2E14.6) by means of the special program control input

&STAFOR='(format)',&END

The options DRUKU, STAKU and PL $\emptyset$ KU were already introduced in Sect. 4.1.3.1 (Interpolation), the options DRUSM, STASM, PLOSM in Sect. 4.1.3.2 (Curve fitting, smoothing). These latter six options refer to processed data and in general require specification of a "new energy grid" for the output.

##### 4.1.5.2 Specification of new grid for output\_

As explained before the options DRUKU, STAKU, DRUSM, STASM permit tabulation of curve points representing the interpolation or smoothing results on various output media. The new grid, i.e. the energy values for the tabulation, will in

general differ from the initial grid of the original data. The user must specify the new grid by the following program control input:

```
_&ØUTINT_FMTØUT='(format)',ZXINT=...,XINT=...,EINHTX=...,  
_FRØM=...,TØ=...,ADD=...,additional information,&END
```

where FMTØUT is the format that is used to print curve points of the smooth curve requested by DRUKU or DRUSM.

ZXINT specifies the medium from which the new grid energies (array XINT) is to be retrieved. There are five possibilities:

ZXINT = n, with integer n, means that in the same program control input list the new grid energies are coded in the form  $XINT=E_1, E_2, \dots, E_n$ , i.e. the array XINT is given as a list of n energy values (real numbers) separated by commas.

ZXINT='KEDAK' means that XINT is to be retrieved from a KEDAK-type library. This requires the following additional information in the &ØUTINT control input list (unless already specified by default):

```
PLEH=...,NAMZ=...,NAMEN=...,FRØM=...,TØ=...
```

(cf. Sect. 4.1.2.1) or an immediately following &KEDAK control input list containing XINTPØ=T.

ZXINT='NEUD' means XINT is to be retrieved from a NEUDADA "transmission" tape. This requires the following additional information in the control input list (unless already specified by default):

```
BDEH=...,ACCNR=...,LTR=...,SEQN=...,FRØM=...,TØ=...,
```

(cf. Sect. 4.1.2.2)

ZXINT='CARD' means XINT is to be read from cards. This requires the following additional information in the &ØUTINT control input list (unless already specified by default):

```
KARTEH=...,FMT=...,ANZ=...,FRØM=...,TØ=...
```

(cf. Sect. 4.1.2.3) or an immediately following &KEDAK control input list containing XINTPØ=T.

ZXINT='DATA' means that the energies of the initial ("current") data set are to be used as "new grid".

EINH TX =  $\left\{ \begin{array}{l} \text{'EV'} \\ \text{'KEV'} \\ \text{'MEV'} \end{array} \right\}$  specifies the unit for XINT

ADD =  $\left\{ \begin{array}{l} \text{T} \\ \text{F} \end{array} \right\}$  causes merging with an already existing XINT array,  
suppresses merging.

If FORTRAN-readable output is to be produced its FORMAT can be specified by the program control input `&STAFOR_STAFMT='(format)',&END`

#### 4.1.6 Summary table

The options available in the various control input lists are shown schematically in Table 1 together with the default options. It should be added that the default option for `&STAFØR_STAFMT=...,&END` is (2E14.6).

Table 1 SIGPLO Options

Option	optional in program control input list						default value	explanatory remarks	
	&KEDAK	&NEUDAD	&KARTEN	&RELGEW	&SMØWDH	&ØUTINT			&PLØTGR
<u>Retrieval</u>									
PLEH	+						+	1	FT number of input device
BDEH		+					+		
KARTEH		+	+				+	2	number of KEDAK names KEDAK names
NAMZ	+						+		
NAMEN	+						+	'0'	letter preceding ACCNR NEUDADA accession number
LTR		+					+		
ACCNR		+					+	0 <sup>+</sup> )	number of data points read
ANZ			+				+		
FRØM	+	+	+				+	0.	minimum energy
TØ							+		
EINHTX			+	+	+		+	1.E+20	maximum energy
EINHTY			+			+	+		
EINHTD			+				+	'EV'	abscissa units
XEINHT		+					+		
YEINHT		+					+	'BARN'	ordinate units
DEINHT		+					+		
FMT		+	+				+	'BARN'	error units
SEQN		+					+		
REWD			+				+	'MEV'	abscissa units
SMGEW			+				+		
MEDIUM		+					+	'BARN'	ordinate units
ZAHL		+					+		
ZAHLXR				+	+		+	'BARN'	error units
DATADD		+	+				+		
ADD							+	'(2E14.6)'	input format
ZXINT							+		
XINTPØ	+	+					+	T	accession number in sequence?
								F	rewinding?
								F	weighting?
								'-----'	medium for uncertainties
								0	number of uncertainties
								0	number of relative weights
								F	merge with "current" set?
								F	merge with current "new grid"?
								0	number of "new grid" points
								F	are these "new grid" points?

+ ) reset for each program control input list



Table 1 (cont.)

Option	optional in program control input list						default value	explanatory remarks
	&KEDAK	&NEUDAD	&KARTEN	&RELGEW	&SMØWDH	&ØUTINT		
<u>Data processing</u>								
KURVE	+	+	+				'NEIN'	interpolation type
SMØØ		+	+				0.+)	maximum chi-squared
FQU					+		0.	new maximum chi-squared?
FISS	+	+	+				F	fission spectrum average?
<u>Plots</u>								
PLØ	+	+	+				T	plot of "current" set?
PLØKU	+	+	+				T	plot of interpolated curve?
PLØSM	+	+	+				T	plot of smooth curve?
PLØZEI	+	+	+				-1	which point symbol?
PLØFQU					+		T	chi-squared displayed?
NEUZEI					+		F	new plot?
XMIN						+	-1.E-60	minimum x-value
XMAX						+	-1.E-60	maximum x-value
XLG						+	-1.E-60	length of x-axis (cm)
YMIN						+	-1.E-60	minimum y-value
YMAX						+	-1.E-60	maximum y-value
YLG						+	-1.E-60	length of y-axis (cm)
RAST						+	'NEIN', '-----'	grid? log scale(s)?
CALP						+	F	CALCOMP plot?

+) reset for each program control input list

Table 1 (cont.)

Option	optional in program control input list					default value	explanatory remarks
	&KEDAK	&NEUDAD	&KARTEN	&RELGEW	&SMØWDH		
<u>Output</u>							
DRUCK	+	+	+			F	"current" set on listing?
STANZ	+	+	+			F	"current" set on cards, tape,...?
DRUKU	+	+	+			F	interpolation results on listing?
STAKU	+	+	+			F	interpolation results on cards, tape,...?
DRUSM		+	+			F	smoothing results on listing?
STASM		+	+			F	smoothing results on cards, tape,...?
FMTØUT					+	'(2E14.6)'	printing format

4.1.7 References

- /1/ R. Meyer, User's Guide to SIGPLO, unpublished (1975)
- /2/ Present Compendium, Ch. III.2
- /3/ Present Compendium, Ch. II.2.3.1.2
- /4/ see CCDN Newsletters for more information on neutron data dissemination by CCDN
- /5/ H. Späth, SPLINE, (1968), unpublished
- /6/ H. Späth, SMØØTH, (1968), unpublished
- /7/ S. Heine, PLØTA, Programmbeschreibung (1967) and updates, unpublished

Appendix: SIGPLO Program List

C-----	SIGPLO.	10
C-----	DATE: 7.4.1973.	20
C-----	PROGRAM TO RETRIEVE, EDIT, PLOT, PRINT, INTERPCLATE,	30
C-----	AND FIT DATASETS THAT ARE CONSIDERED AS EXPERIMENTAL	40
C-----	OR EVALUATED. DATA SOURCES MAY BE ANY FORMATTED INPUT,	50
C-----	KEDAK LEBRARY FORMAT DATA SETS OR DATA SETS WRITTEN IN	60
C-----	NEUDADA TRANSMISSION FORMAT.	70
	RFAL*8 NAMEN(3)	80
	REAL NEIN, LINE, KURVF, MEDIUM, LTR, MOD	90
	INTEGER ACCNR, BDEH	100
	INTEGER ANZ, PLEH, ZAHLXR, ZXINT, WNR, ZAHL, PLOZFI	110
	LOGICAL DRUCK, STANZ, PLO, PLOKU, DRUKU, STAKU, PLOSM, DRUSM, STASM, SMGEW,	120
	INEUZEI, XINTPO, PLOFQU, FISS, SEQN, ADD, CALP, CALPJ, DATADD, TST, REWD, KOK	130
	2, TSTS	140
	DIMENSION NTEXT(15), XPH(2), YPH(2)	150
	DIMENSION RTITEL(20), STEUER(20), NNP(40), TA(2,7), TB(2)	160
	DIMENSION XP(1000), YP(1000), T(1000), S(1000), S2(1000)	170
	DIMENSION Z1(1002), Z2(1002), Z3(1002), Z4(1002), Z5(1002), Z6(1002), Z7	180
	1(1002)	190
	DIMENSION A(1000), B(1000), C(1000), D(1000)	200
	DIMENSION XS(1000), YS(1000)	210
	DIMENSION XINT(1000)	220
	DIMENSION XR(20), PR(20)	230
	DIMENSION P(1000), DELTA(1000), PS(1000)	240
	DIMENSION XL(1000), YL(1000), XSK(200), YSK(200)	250
	REAL*8 BSK(200)	260
	EQUIVALENCE (FROM, VON, EUNT, E1, EMIN), (TO, BIS, EOB, E2, EMAX)	270
	EQUIVALENCE (S2(1), Z1(1), YL(1)), (XP(1), Z3(1)), (YP(1), Z4(1)), (DELTA	280
	1(1), Z5(1)), (T(1), Z6(1)), (S(1), Z7(1))	290
	EQUIVALENCE (Z2(1), XL(1), XSK(1)), (XL(201), YSK(1)), (XL(401), BSK(1))	300
	COMMON XP, YP, ANZ	310
	COMMON/PAR/XMAX, SX, YMAX, SY, XMIN, YMIN, XMAL, XMIL, YMAL, YMIL, NLGGX,	320
	1 NLGGY, NX, NY, INDZ, JNDZ, XMILJ, YMILJ, JP	330
	COMMON/CLP/CALP	340
	COMMON/PARM1/ MAXOPT, PLEH, PLOZFI, ZXINT, IDPLOT, CALPJ	350
	COMMON/PARM2/ IPLOGR, ISTAFO, ISMO, IFMOUT, ISMNP, DISP, EINHTX, EINHTY,	360
	1 EINHTD, XEINHT, YEINHT, DEINHT, LTR, KARTEH, FMTOUT(20), STAFMT(20),	370
	2 FMT(20), DRUCK, STANZ, PLC, KURVE, PLOKU, DRUKU, STAKU, PLOSM, DRUSM,	380
	3 STASM, SMGEW, NEUZEI, XINTPO, PLOFQU, FISS, SEQN, ADD, DATADD, TST,	390
	4 REWD, NANZ, NADD, ZAHL, MEDIUM, WNR,	400
	5 RAST(2), ISMAX, FQU, NAMZ, ZAHLXR, TSTS	410
	COMMON/PARM3/ SMOO	420
	COMMON/PARM4/FROM, TO	430
	COMMON/INOUT/KOUT, KIN, KPU, KINC	440
	NAMELIST/KARTEN/KARTEH, FMT, ANZ, DRUCK, STANZ, PLO, KURVE, PLOKU, DRUKU, S	450
	TAKU, SMOO, PLOSM, DRUSM, STASM, SMGEW, EINHTX, EINHTY, EINHTD, FISS, PLOZEI	460
	2, DATADD, TST, E1, E2, REWD, DISP, TSTS, FROM, VON, EUNT, EMIN, TO, BIS, EOB,	470
	3 EMAX	480
	NAMELIST/KFDAK/PLEH, NAMZ, NAMEN, EUNT, EOB, DRUCK, STANZ, PLO, KURVE, PLOK	490
	1U, DRUKU, STAKU, XINTPC, FISS, PLOZEI, TST, FROM, VON, E1, EMIN,	500
	2 FMAX, TO, BIS, E2	510
	NAMELIST/STAFCR/STAFMT	520
	NAMELIST/RELGFW/ZAHLXR, XR, PR, EINHTX	530
	NAMELIST/OUTINT/FMTOUT, ZXINT, XINT, EINHTX, ADD, PLEH, NAMEN, NAMZ,	540
	1BDFH, ACCNR, LTR, SEQN, VON, BIS, ANZ, FMT, KARTEH, REWD,	550

2 FROM,EUNT,E1,EMIN,TO,ECB,E2,EMAX	560	CALL ECONV(ANZ,YP,EINHXY)	1110
NAMLIST/SMQWDH/FQU,ZAHLXR,XR,PR,NEUZEI,EINHXY,PLOFQU,TST,TSTS	570	IF(.NOT.SMGEW) GOTO 128	1120
DATA NEIN/'NEIN'/,LINE/'LINE'/,SPLI/'SPLI'/,XLOG/'XLOG'/,YLOG/'YLO	580	C-----CONVERT ERROR BARS ALSO.	1130
IG'/,MOD/'MOD'/	590	IF(EINHXY.EQ.PERC) GOTO 113	1140
DATA TEND/'ENDE'/,NEUD/'NEUD'/,KEDA/'KEDA'/,KART/'KARD'/	600	CALL FCONV(ANZ,DELTA,EINHXY)	1150
DATA NNP/9,0,3,4,5,6,7,1,8,2,0,1,2,7,8,0,1,2,7,8,0,1,2,7,8,0,1,2,7	610	GOTO 129	1160
1,8,0,1,2,7,8,0,1,2,7,8/	620	113 CALL PERCNT(ANZ,YP,DELTA)	1170
DATA RT/'RT' /,PERC/'PERC'/	630	GOTO 128	1180
DATA TA/' &KA','RTEN',' &NE','UDAD',' &KE','DAK ',' &PL','OTGR','	640	C-----NEUDADA INPUT IS REQUESTED.	1190
1&ST','AFOR',' &OU','TINT',' &RE','LGEW'/	650	120 CALL NEUDA(XINT,DELTA,&100)	1200
DATA TB/' &SM','QWDH'/	660	C-----ADF APPENDS DATA TO CURRENT SET IF REQUESTED.	1210
DATA NDATA/'DATA'/	670	128 CALL ADF(XL,YL,Z6,NADD,NANZ,ANZ,XP,YP,DELTA,DATAADD,&100)	1220
DATA NTEXT/'FQU=' ,14*' */	680	GOTO 200	1230
C-----INITIALIZATION ONCE FOR WHOLE JOB.	690	C-----KEDAK DATA ARE TO BE FETCHED.	1240
CALL INIT1	700	130 READ(KIN,KEDAK)	1250
C-----PRINT INPUT FROM KINC.	710	KDK=.TRUE.	1260
CALL PRIEIN(KINC,KOUT,60)	720	WRITE(KOUT,636)	1270
C-----PUT CONTROL INPUT GNT0 UNIT FT05.	730	WRITE(KOUT,619) PLEH,NAMZ,(NAMEN(I),I=1,2)	1280
CALL COINPT(RTITEL,STEUER)	740	IF(NAMZ.GE.3) WRITE(KOUT ,656) NAMEN(3)	1290
WRITE(KOUT,638)	750	WRITE(KOUT,657) EUNT,EOB,DRUCK,STANZ,PLO,PLOZEI,KURVE,PLOKU,DRUKU,	1300
C-----RT CARD?	760	1 STAKU,DISP,XINTPO,FISS	1310
1000 READ(KIN,500,END=901,ERR=801) RTITEL	770	CALL LDFOPN(PLEH,IOATUM,&903)	1320
IF(RTITEL(1).EQ.TEND) GOTO 810	780	131 CALL KEDSUB(NR,NAMZ,NAMEN,EUNT,EOB,XINTPO,XINT,ZXINT,IFMCUT,&100)	1330
IF(RTITEL(1).NE.RT) GOTO 802	790	IF(ANZ.NE.0) GOTO 200	1340
C-----PRINT RT CARD AND PERFORM RT-INITIALIZATION.	800	IF(NR.NE.0) GOTO 132	1350
WRITE(KOUT,600) RTITEL	810	WRITE(KOUT,620)	1360
1003 CALL INIT2	820	GOTO 100	1370
DC 1004 I=1,1000	830	132 WRITE(KOUT,621)	1380
1004 PS(I)=1.	840	GOTO 100	1390
100 ANZ=0	850	C-----PLOT SIZE DEFINITION.	1400
SMOQ=NEIN	860	C-----COMPLETE PREVIOUS PLOT, IF IT WAS CALCOMP.	1410
KDK=.FALSE.	870	145 XPH(1)=XMILJ	1420
C-----TEST COMMAND (=NAMLIST NAME.)	880	XPH(2)=XPH(1)	1430
READ(KIN,500,END=902,ERR=803) STEUER	890	YPH(1)=YMILJ	1440
DC 101 J=1,MAXOPT	900	YPH(2)=YPH(1)	1450
CALL ASSIGN(STEUER,TA,J,NR,MAXOPT)	910	CALL PLOTA(XPH,YPH,-2,2,9,1,1,1,10,XMAX,XMIN,SX,YMAX,YMIN,SY,1,	1460
IF(NR.EQ.0) GOTO 101	920	1)DPL0T,0,0,0)	1470
BACKSPACE KIN	930	WRITE(KOUT,645)	1480
GOTO (110,120,130,140,150,160,170),J	940	STCP	1490
101 CONTINUE	950	140 CALL PREPLO	1500
GOTO 804	960	GOTO 100	1510
C-----FORMATTED INPUT IS REQUESTED.	970	C-----FORMATTED OUTPUT IS TO BE PRODUCED AND OUTPUT	1520
110 READ(KIN,KARTEN)	980	C-----PARAMETERS MUST BE DEFINED.	1530
WRITE(KOUT,636)	990	150 READ(KIN,STAFGR)	1540
WRITE(KOUT,612) KARTEH,FMT,ANZ,DATADD,DRUCK,STANZ,PLO,PLOZEI,	1000	WRITE(KOUT,636)	1550
1)KURVE,PLOKU,DRUKU,STAKU,DISP,EINHXY,EINHXY,EINHXY	1010	WRITE(KOUT,622) STAFMT	1560
WRITE(KOUT,653) E1,E2,REWD	1020	WRITE(KOUT,636)	1570
IF(SMOQ.NE.NEIN) GOTO 111	1030	1STAFQ=1	1580
WRITE(KOUT,613) SMOQ,PLOSM,DRUSM,STASM,SMGEW,FISS	1040	GOTO 100	1590
GOTO 112	1050	C-----INTERPOLATION REQUESTED. SAVE CURRENT INTERPOLATION	1600
111 WRITE(KOUT,614) SMOQ,PLOSM,DRUSM,STASM,SMGEW,FISS	1060	C-----POINTS.	1610
1SMO=1	1070	160 CALL STRING(Z3(1),XINT(1),4*ZXINT)	1620
112 CALL LESENI(KARTEH,FMT,SMGEW,DELTA,E1,E2,REWD,&100)	1080	NZXINT=ZXINT	1630
C-----CONVERT DATA TO INTERNAL PHYSICAL UNITS, IF REQUIRED.	1090	READ(KIN,OUTINT)	1640
CALL ECONV(ANZ,XP,EINHXY)	1100	WRITE(KOUT,636)	1650

C-----BRANCH TO CORRECT READING ROUTINE FOR INTERPOL. POINTS.	1660	DO 2001 I=1,ANZ	2210
IF(ZXINT.EQ.NDAT) GOTO 161	1670	XP(I)=Z1(I)	2220
IF(ZXINT.EQ.KEDA) GOTO 167	1680	YP(I)=Z2(I)	2230
IF(ZXINT.EQ.NEUD) GOTO 169	1690	2001 DELTA(I)=Z6(I)	2240
IF(ZXINT.EQ.KART) GOTO 1670	1700	IF(.NOT.DRUCK) GOTO 201	2250
CALL ECONV(ZXINT,XINT,EINH TX)	1710	C-----PRINT DATA POINTS.	2260
WRITE(KOUT,623) FMTOUT,ZXINT,EINH TX,ADD,(XINT(I),I=1,ZXINT)	1720	WRITE(KOUT,630)	2270
GCTO 162	1730	IF(KDK) GOTO 2004	2280
161 WRITE(KOUT,641) FMTOUT,ZXINT,EINH TX,ADD	1740	IF(SMGEW) GOTO 2002	2290
GCTO 165	1750	2004 CALL DRUCKX(XP,YP,ANZ)	2300
C-----APPEND DATA, IF REQUESTED.	1760	GCTO 201	2310
162 IF(.NOT.ADD) GOTO 165	1770	2002 CALL DRUCK3(ANZ,XP,YP,DELTA)	2320
IF(IFMOUT.NE.1) GOTO 166	1780	201 IF(.NOT.STANZ) GOTO 203	2330
WRITE(KOUT,642)	1790	C-----FORMATTED OUTPUT OF DATA POINTS.	2340
IF(ZXINT+NZXINT.LE.1000) GOTO 163	1800	IF(ISTAFO.EQ.1) GOTO 202	2350
WRITE(KOUT,643)	1810	WRITE(KOUT,602)	2360
NZXINT=1000-ZXINT	1820	202 IF(DISP.EQ.MOD) GOTO 2020	2370
163 I=ZXINT+1	1830	WRITE(KPU,500) RTITEL	2380
DO 164 J=1,NZXINT	1840	2020 IF(SMGEW) GOTO 2021	2390
XINT(I)=Z3(J)	1850	WRITE(KPU,STAFMT) (XP(I),YP(I),I=1,ANZ)	2400
164 I=I+1	1860	GOTO 203	2410
ZXINT=ZXINT+NZXINT	1870	2021 WRITE(KPU,STAFMT) (XP(I),YP(I),DELTA(I),I=1,ANZ)	2420
165 IFMOUT=1	1880	203 IF(.NOT.PLO) GOTO 206	2430
WRITE(KOUT,636)	1890	C-----PLOT DATA POINTS.	2440
GCTO 100	1900	IF(IPLOGR.EQ.1) GOTO 204	2450
166 WRITE(KOUT,644)	1910	WRITE(KOUT,603)	2460
GCTO 165	1920	GCTO 206	2470
C-----RETRIEVE INTERPOLATION ENERGIES FROM KEDAK.	1930	204 IZ=0	2480
167 WRITE(KOUT,647) FMTOUT,ZXINT,PLEH,NAMZ,(NAMEN(I),I=1,NAMZ),VON,BIS	1940	IF(ANZ.NE.1) GOTO 205	2490
CALL LDFOPN(PLEH,IDATUM,8903)	1950	IZ=1	2500
168 CALL KEDEN(NAMZ,NAMEN,VON,BIS,XINT,ZXINT,81697)	1960	CALL XCORR	2510
WRITE(KOUT,652) ZXINT,(XINT(I),I=1,ZXINT)	1970	C-----A SINGLE DATA POINTS CANNOT BE HANDLED BY PLOTA.	2520
GCTO 162	1980	205 JP=JP+1	2530
C-----RETRIEVE INTERPOLATION ENERGIES FROM NEUDACA TAPE.	1990	NP=NNP(JP)	2540
169 CALL NEUINT(XINT,BDEH,ACCN,8100)	2000	IF(PLOZEI.NE.-1) NP=PLOZEI	2550
GCTO 162	2010	C-----FOR PLOTS WITH ONE AXIS LOGARITHMIC, USE PLOGXY.	2560
1697 IFMOUT=0	2020	IF(RAST(1).EQ.XLOG.OR.RAST(2).EQ.YLOG) GOTO 2003	2570
GCTO 100	2030	CALL PLOTA(XP,YP,ANZ,1,NP,1,1,1,INDZ,XMAX,XMIN, SX, YMAX, YMIN, SY, 1, I	2580
1670 WRITE(KOUT,654) ZXINT,KARTEH,ANZ,VON,BIS,RWD,FMT,FMTOUT	2040	IDPLOT,0,0,0)	2590
CALL LESEN(KARTEH,FMT,REWD,ZXINT,ANZ,VON,BIS,XINT,81697)	2050	WRITE(KOUT,604) NP	2600
CALL ECONV(ZXINT,XINT,EINH TX)	2060	CALL SUB1(INDZ,RAST,NEIN,XLOG,YLOG,KREX,KREY,8211)	2610
WRITE(KOUT,651) EINH TX,ZXINT,(XINT(I),I=1,ZXINT)	2070	C-----PLOT SCALE ALSO, IF REQUESTED.	2620
GCTO 162	2080	CALL RASTER(KREX,KREY,XMAX,XMIN, SX, YMAX, YMIN, SY, INDZ, XSK, YSK)	2630
C-----RELATIVE WEIGHTS FOR SMOOTH ARE REQUESTED.	2090	GOTO 211	2640
170 READ(KIN,RELGEW)	2100	2003 CALL PLOGXY(XP,YP,ANZ,1,NP,1,INDZ,1,	2650
CALL ECONV(ZAHLXR,XR,EINH TX)	2110	1 XL,YL,RAST,XMAL,XMIL,XMAX,XMIN, SX, YMAL, YMIL, YMAX, YMIN, SY,	2660
WRITE(KOUT,636)	2120	2 IDPLOT,NLGGX,NLGGY,NX,NY,XSK,YSK,BSK)	2670
WRITE(KOUT,624) ZAHLXR,EINH TX,(XR(I),PR(I),I=1,ZAHLXR)	2130	WRITE(KOUT,604) NP	2680
WRITE(KOUT,636)	2140	CALL SUB1(INDZ,RAST,NEIN,XLOG,YLOG,KREX,KREY,8211)	2690
GCTO 100	2150	CALL RASTER(KREX,KREY,XMAL,XMIL, SX, YMAL, YMIL, SY, INDZ, XSK, YSK)	2700
C-----	2160	GCTO 211	2710
C-----	2170	211 IF(IZ.NE.1) GOTO 206	2720
C-----DO THE REQUESTED WORK.-----S	2180	ANZ=1	2730
C-----SORT DATA.	2190	C-----IS INTERPOLATION OF DATA POINTS TO PRODUCE A	2740
200 CALL ORDNEI(ANZ,XP,YP,DELTA,Z1,Z2,Z6)	2200	C-----CURVE REQUESTED?	2750

C-----TWO MODES OF INTERPOLATION ARE SUPPLIED:	2760	215 XINT(I)=Z1(I)	3310
C-----LINEAR AND SPLINE.	2770	CALL KURVE4(ANZ,XP,YP,ZXINT,XINT,S,S2,A,B,C,D,NA1,NA2,IKU,&220)	3320
206 IF(KURVE.EQ.NEIN) GOTO 220	2780	IF(DISP.EQ.MOD) GOTO 2150	3330
IF(KURVE.EQ.LINE) IKU=1	2790	WRITE(KPU,500) RTITEL	3340
IF(KURVE.EQ.SPLI) IKU=2	2800	2150 WRITE(KPU,STAFMT) (XINT(I),S(I),I=NA1,NA2)	3350
IF(.NOT.PLOKU) GOTO 208	2810	GOTO 220	3360
IF(IPLOGR.EQ.1) GOTO 207	2820	C-----SAVE DATA FOR SMOOTH FIT IF REQUESTED.	3370
WRITE(KOUT,603)	2830	220 IF(SMOO.EQ.NEIN) GOTO 230	3380
GOTO 208	2840	IF((ANZ+ISMAX).LE.1000) GOTO 224	3390
207 CALL KURVE1(ANZ,XP,YP,M,T,S,S2,A,B,C,D,IKU,&243)	2850	ANZ=1000-ISMAX	3400
IF(.NOT.TST) GOTO 241	2860	WRITE(KOUT,658) ANZ	3410
WRITE(KOUT,650) (T(I),S(I),I=1,M)	2870	658 FORMAT(/5X,'FEHLERNACHRICHT'/8X,'ZAHL DER PUNKTE FUER SMOOTH-FIT'	3420
241 IF(.NOT.FISS) GOTO 240	2880	*,' WUERDE GROESSER ALS 1000. ES WERDEN VON DEN SOBEN GELESENEN ',	3430
C-----FISSION SPECTRUM AVERAGE.	2890	*'DATEN NUR DIE ERSTEN',I5,' PUNKTE GENOMMEN.')	3440
CALL QFISS(M,T,S,A)	2900	IF(ANZ.LE.0) GOTO 230	3450
C-----PLOT CURVE WITH REQUESTED INTERPOLATION.	2910	224 DO 221 I=1,ANZ	3460
240 IF(RAST(1).EQ.XLOG.OR.RAST(2).EQ.YLOG) GOTO 2041	2920	XS(I+ISMAX)=XP(I)	3470
CALL PLOTA(T,S,M,2,0,1,1,1,INDZ,XMAX,XMIN,SY,YMAX,YMIN,SY,1,1DPL0T	2930	221 YS(I+ISMAX)=YP(I)	3480
1,0,0,0)	2940	FQU=FQU+SMO0	3490
CALL SUB1(INDZ,RAST,NEIN,XLOG,YLCG,KREX,KREY,&243)	2950	IF(.NOT.SMGEW) GOTO 223	3500
CALL RASTER(KREX,KREY,XMAX,XMIN,SY,YMAX,YMIN,SY,INDZ,XSK,YSK)	2960	DO 222 I=1,ANZ	3510
GOTO 243	2970	222 PS(I+ISMAX)=DELTA(I)	3520
2041 CALL PLOGXY(T,S,M,2,0,1,INDZ,1,	2980	223 ISMAX=ISMAX+ANZ	3530
1 XL,YL,RAST,XMAL,XMIL,XMAX,XMIN,SY,YMAL,YMIL,YMAX,YMIN,SY,	2990	GOTO 230	3540
2 IDPLOT,NLGGX,NLGGY,NX,NY,XSK,YSK,BSK)	3000	230 GOTO 100	3550
CALL SUB1(INDZ,RAST,NEIN,XLOG,YLCG,KREX,KREY,&243)	3010	C-XX	3560
CALL RASTER(KREX,KREY,XMAL,XMIL,SY,INDZ,XSK,YSK)	3020	C-XX	3570
GOTO 243	3030	C-----SMOOTH FIT SECTION.	3580
208 IF(.NOT.FISS) GOTO 243	3040	300 IF(ISMAX.LT.3) GOTO 809	3590
CALL KURVE1(ANZ,XP,YP,M,T,S,S2,A,B,C,D,IKU,&243)	3050	C-----SORT DATA.	3600
CALL QFISS(M,T,S,A)	3060	CALL ORDNEI(ISMAX,XS,YS,PS,Z1,Z2,Z3)	3610
C-----PRINT CURVE ACCORDING TO USER SPECIFICATICNS.	3070	DC 301 I=1,ISMAX	3620
243 IF(.NOT.DRUKU) GOTO 209	3080	XS(I)=Z1(I)	3630
IF(IFMOUT.NE.0) GOTO 212	3090	YS(I)=Z2(I)	3640
WRITE(KOUT,627)	3100	301 PS(I)=Z3(I)	3650
GOTO 220	3110	CALL EQUEN(ISMAX,XS,YS,PS)	3660
212 IF(ZXINT.NE.NDAT) GOTO 216	3120	IF(ZAHLXR.EQ.0) GOTO 310	3670
ZXINT=ANZ	3130	C-----SORT WEIGHTS.	3680
DO 217 I=1,ZXINT	3140	309 CALL ORDNEI(ZAHLXR,XR,PR,Z3,Z1,Z2,Z3)	3690
217 XINT(I)=XP(I)	3150	C-----SUPPLY RELATIFV WEIGHTING.	3700
216 CALL ORDNEI(ZXINT,XINT,Z2,Z2,Z1,Z2,Z2)	3160	DC 302 I=1,ZAHLXR	3710
DO 213 I=1,ZXINT	3170	XR(I)=Z1(I)	3720
213 XINT(I)=Z1(I)	3180	302 PR(I)=Z2(I)	3730
CALL KURVE4(ANZ,XP,YP,ZXINT,XINT,S,S2,A,B,C,D,NA1,NA2,IKU,&209)	3190	ASSIGN 305 TO NST	3740
WRITE(KOUT,628)	3200	K=1	3750
WRITE(KOUT,FMTOUT) (XINT(I),S(I),I=NA1,NA2)	3210	DC 327 I=1,ISMAX	3760
C-----FORMATTED UOTPUT OF CURVE ACCORDING USER SPECIFICATIONS.	3220	303 IF(XS(I).LT.XR(K)) GOTO NST,(305,304)	3770
209 IF(.NOT.STAKU) GOTO 220	3230	IF(K.EQ.ZAHLXR) GOTO 306	3780
IF(ISTAFO.EQ.1) GOTO 210	3240	K=K+1	3790
WRITE(KOUT,602)	3250	ASSIGN 304 TO NST	3800
210 IF(IFMOUT.NE.0) GOTO 214	3260	GOTO 303	3810
WRITE(KOUT,627)	3270	306 IANF=I	3820
GOTO 220	3280	GOTO 307	3830
214 CALL ORDNEI(ZXINT,XINT,Z2,Z2,Z1,Z2,Z2)	3290	304 P(I)=PS(I)*PR(K-1)	3840
DC 215 I=1,ZXINT	3300	GOTO 327	3850



```

305 P(I)=PS(I)
327 CONTINUE
GCTO 310
307 DO 308 I=IANF,ISMAX
308 P(I)=PS(I)*PR(ZAHLXR)
GOTO 3102
310 DO 3100 I=1,ISMAX
3100 P(I)=PS(I)
C-----PRINT IDENTIFICATION AND PROVIDE PRINTOUT OF DATA.
3102 WRITE(KOUT,631) FQU
CALL DRUCK3(ISMAX,XS,YS,P)
C-----SMOOTH. USE PROGRAM OF H. SPAETH. (SEE INR-REPORT)
CALL SMOOTH(ISMAX,XS,YS,P,FQU,A,B,C,D,Z1,Z2,Z3,Z4,Z5,Z6,Z7)
C-----PLOT SMOOTHED CURVE, IF REQUESTED.
IF(.NOT.PLOSM) GOTO 313
IF(IPLOGR.EQ.1) GOTO 312
WRITE(KOUT,603)
GOTO 313
312 CALL KURVE2(ISMAX,XS,YS,M,T,S,A,B,C,D,8343)
C-----PROVIDE TESTPRINTOUT OF SMOOTHED CURVE, IF REQUESTED.
IF(TSTS) CALL TSTPRO(ISMAX,XS,YS,P,A,FQU)
IF(.NOT.TST) GOTO 3124
WRITE(KOUT,655)
CALL DRUCKX(T,S,M)
3124 IF(.NOT.FISS) GOTO 340
C-----FISSION SPECTRUM AVERAGE.
CALL QFISS(M,T,S,Z1)
C-----PLOT INTO NEW OR OLD PLOT?
C-----OLD PLOT.
340 IF(NEUZEI) GOTO 320
C-----INCREMENT PLOT IDENTIFIER.
ISMNP=ISMNP+1
NP=NNP(ISMNP)
IF(ISMNP.EQ.1) GOTO 321
NPA=M-1
IF(RAST(1).EQ.XLOG.OR.RAST(2).EQ.YLOG) GOTO 304C
CALL PLOTA(T,S,M,3,NP,1,1,NPA,INDZ,XMAX,XMIN,SX,YMAX,YMIN,SY,1,
1IDPLOT,0,0,0)
WRITE(KOUT,633) NP
322 CALL SUB1(INDZ,RAST,NEIN,XLOG,YLOG,KREX,KREY,8343)
319 CALL RASTER(KREX,KREY,XMAX,XMIN,SX,YMAX,YMIN,SY,INDZ,XSK,YSK)
GCTO 343
3040 CALL PLOGXY(T,S,M,3,NP,NPA,INDZ,1,
1 XL,YL,RAST,XMAL,XMIL,XMAX,XMIN,SX,YMAL,YMIL,YMAX,YMIN,SY,
2 IDPLOT,NLGGX,NLGGY,NX,NY,XSK,YSK,BSK)
WRITE(KOUT,633) NP
3220 CALL SUB1(INDZ,RAST,NEIN,XLOG,YLOG,KREX,KREY,8343)
3190 CALL RASTER(KREX,KREY,XMAL,XMIL,SX,YMAL,YMIL,SY,INDZ,XSK,YSK)
GOTO 343
313 IF(.NOT.FISS) GOTO 343
CALL KURVE2(ISMAX,XS,YS,M,T,S,A,B,C,D,8343)
IF(.NOT.TST) GOTO 3134
WRITE(KOUT,655)
CALL DRUCKX(T,S,M)
3134 CALL QFISS(M,T,S,Z1)

```

```

3860 343 IF(.NOT.DRUSM) GOTO 316
3870 C-----PRINT SMOOTHED CURVE ACCORDING USFR SPECIFICATIONS.
3880 IF(IFMOUT.NE.0) GOTO 314
3890 WRITE(KOUT,610)
3900 GCTO 330
3910 314 IF(ZXINT.NE.NDAT) GOTO 324
3920 ZXINT=ISMAX
3930 DO 323 I=1,ZXINT
3940 323 XINT(I)=XS(I)
3950 324 CALL ORDNEI(ZXINT,XINT,Z2,Z2,Z1,Z2,Z2)
3960 DO 315 I=1,ZXINT
3970 315 XINT(I)=Z1(I)
3980 CALL KURVE3(ISMAX,XS,YS,ZXINT,XINT,S,A,B,C,D,NA1,NA2,8330)
3990 WRITE(KOUT,629)
4000 WRITE(KOUT,FMTOUT) (XINT(I),S(I),I=NA1,NA2)
4010 316 IF(.NOT.STASM) GOTO 330
4020 C-----PROVIDE FORMATTED OUTPUT, IF REQUESTED.
4030 IF(ISTAFO.EQ.1) GOTO 311
4040 WRITE(KOUT,602)
4050 311 IF(IFMOUT.NE.0) GOTO 317
4060 WRITE(KOUT,610)
4070 GCTO 330
4080 317 CALL ORDNEI(ZXINT,XINT,Z2,Z2,Z1,Z2,Z2)
4090 DO 318 I=1,ZXINT
4100 318 XINT(I)=Z1(I)
4110 CALL KURVE3(ISMAX,XS,YS,ZXINT,XINT,S,A,B,C,D,NA1,NA2,8330)
4120 IF(DISP.EQ.MOD) GOTO 3180
4130 WRITE(KPU,500) RTITEL
4140 3180 WRITE(KPU,STAFMT) (XINT(I),S(I),I=NA1,NA2)
4150 GCTO 330
4160 C-----NEW PLOT WAS REQUESTED FOR SMOOTH.
4170 320 IDPLOT=IDPLOT+1
4180 IF(PLOFQU) GOTO 325
4190 IF(RAST(1).EQ.XLOG.OR.RAST(2).EQ.YLOG) GOTO 32CC
4200 CALL PLOTA(XS,YS,ISMAX,1,4,1,1,1,JNDZ,XMAX,XMIN,SX,YMAX,YMIN,SY,1,
1IDPLOT,0,0,0)
4220 GCTO 326
4230 3200 CALL PLOGXY(XS,YS,ISMAX,1,4,1,JNDZ,1,
1 XL,YL,RAST,XMAL,XMIL,XMAX,XMIN,SX,YMAL,YMIL,YMAX,YMIN,SY,
2 IDPLOT,NLGGX,NLGGY,NX,NY,XSK,YSK,BSK)
4250 GOTO 3260
4260 325 CALL CONVX(FQU,NTEXT(2),5HE12.4)
4280 IF(RAST(1).EQ.XLOG.OR.RAST(2).EQ.YLOG) GOTO 3250
4290 CALL PLOTA(XS,YS,ISMAX,1,4,1,2,1,JNDZ,XMAX,XMIN,SX,YMAX,YMIN,SY,
1NTEXT,1IDPLOT,0,0,0)
4310 GCTO 326
4320 3250 CALL PLOGXY(XS,YS,ISMAX,1,4,1,JNDZ,NTEXT,
1 XL,YL,RAST,XMAL,XMIL,XMAX,XMIN,SX,YMAL,YMIL,YMAX,YMIN,SY,
2 IDPLOT,NLGGX,NLGGY,NX,NY,XSK,YSK,BSK)
4330 GCTO 3260
4340 326 JJ=0
4350 IF(JNDZ.GE.10) JJ=10
4360 CALL PLOTA(T,S,M,2,0,1,1,1,1,JJ,XMAX,XMIN,SX,YMAX,YMIN,SY,1,
1IDPLOT,
10,0,0)
4370 JJ=1
4380
4390
4400

```

```

4410
4420
4430
4440
4450
4460
4470
4480
4490
4500
4510
4520
4530
4540
4550
4560
4570
4580
4590
4600
4610
4620
4630
4640
4650
4660
4670
4680
4690
4700
4710
4720
4730
4740
4750
4760
4770
4780
4790
4800
4810
4820
4830
4840
4850
4860
4870
4880
4890
4900
4910
4920
4930
4940
4950

```

```

CALL SUB1(J,RAST,NEIN,XLOG,YLOG,KREX,KREY,8313)
WRITE(KOUT,634)
GOTO 319
3260 JJ=0
IF(JNDZ.GE.10) JJ=10
CALL PLOGXY(T,S,M,2,0,1,JJ,1,
1 XL,YL,RAST,XMAL,XMIL,XMAX,XMIN,SX,YMAL,YMIL,YMAX,YMIN,SY,
2 IDPLOT,NLGGX,NLGGY,NX,NY,XSK,YSK,BSK)
J=1
CALL SUB1(J,RAST,NEIN,XLOG,YLOG,KREX,KREY,8313)
GOTO 3190
321 IF(RAST(1).EQ.XLOG.OR.RAST(2).EQ.YLOG) GOTO 321C
CALL PLOTA(T,S,M,2,NP,1,1,1,INDZ,XMAX,XMIN,SX,YMAX,YMIN,SY,1,
1 ICPL0T,0,0,0)
GOTO 322
3210 CALL PLOGXY(T,S,M,2,NP,1,INDZ,1,
1 XL,YL,RAST,XMAL,XMIL,XMAX,XMIN,SX,YMAL,YMIL,YMAX,YMIN,SY,
2 IDPLOT,NLGGX,NLGGY,NX,NY,XSK,YSK,BSK)
GOTO 3220
C-----END OF SMOOTH. IS ANOTHER ONE REQUIRED?
330 ISMO=0
ZHLXR=0
GOTO 100
C*****
C***** FEHLERROUTINEN *****
801 WRITE(KOUT,605)
GOTO 100
802 WRITE(KOUT,606) RTITEL
BACKSPACE KIN
GOTO 1003
803 WRITE(KOUT,607)
GOTO 100
804 IF(ISMO.NE.1) GOTO 805
BACKSPACE KIN
WRITE(KOUT,636)
GOTO 300
805 IF(STEUER(1).NE.TB(1)) GOTO 807
IF(STEUER(2).NE.TB(2)) GOTO 807
BACKSPACE KIN
ZHLXR=0
READ(KIN,SMOWDH)
WRITE(KOUT,636)
WRITE(KOUT,639) ZHLXR,NEUZEI,PLOFQU
IF(ISMAX.LT.3) GOTO 809
IF(ZHLXR.EQ.0) GOTO 310
WRITE(KOUT,640) EINHTX
CALL ECONV(ZHLXR,XR,EINHTX)
WRITE(KOUT,626) (XR(I),PR(I),I=1,ZHLXR)
GOTO 309
807 IF(STEUER(1).EQ.TEND) GOTO 810
IF(STEUER(1).NE.RT) GOTO 808
BACKSPACE KIN
WRITE(KOUT,638)
GOTO 1000

```

```

4960 808 WRITE(KOUT,601) STEUER 5510
4970 GOTO 100 5520
4980 809 WRITE(KOUT,632) 5530
4990 ISMO=0 5540
5000 GOTO 100 5550
5010 810 WRITE(KOUT,637) 5560
5020 IF(CALPJ) GOTO 145 5570
5030 STOP 5580
5040 C ***** ENDRROUTINEN ***** 5590
5050 901 WRITE(KOUT,611) 5600
5060 STOP 5610
5070 902 WRITE(KOUT,611) 5620
5080 IF(ISMO.EQ.1) GOTO 300 5630
5090 STOP 5640
5100 903 WRITE(KOUT,608) 5650
5110 GOTO 100 5660
5120 C ***** FORMATE ***** 5670
5130 500 FORMAT(20A4) 5680
5140 600 FORMAT(1H0,7X,84(' ')/8X,'*',82X,'**/8X,'*',1X,20A4,1X,'**/8X,'*', 5690
5150 182X,'**/8X,84(' ')/8X) 5700
5160 601 FORMAT(/5X,'FEHLERNACHRICHT'/8X,'EINGELESENE KARTE:'/8X,'***',20A 5710
5170 14,'**'/8X,'WURDE NICHT ALS STEUERKARTE ERKANNT,OBWOHL SOLCHE ERWA 5720
5180 2RTET WURDE'/5X,'AKTION:UEBERLESEN BIS ERKENNEN EINER STEUER-,ENDE- 5730
5190 3CDE RT-KARTE'//) 5740
5200 602 FORMAT(/5X,'NACHRICHT:'/8X,'STANZEN WURDE GEWUENSCHT.DA STAFORKAR 5750
5210 1TE FEHLT,WIRD (2E12.4) ALS AUSGABEFORMAT ANGENOMMEN') 5760
5220 603 FORMAT(/5X,'FEHLERNACHRICHT:'/8X,'PLOT WURDE GEWUENSCHT.PLOTGR-KA 5770
5230 1RTE FEHLT'//) 5780
5240 604 FORMAT(' PLOTNACHRICHT: NP =' ,I3) 5790
5250 605 FORMAT(/5X,'FEHLERNACHRICHT'/8X,'LESEFEHLER BEIM VERSUCH,TITEL-KA 5800
5260 1RTE ZU LESEN'/5X,'AKTION:KARTE WIRD UEBERGANGEN.ES WIRD VERSUCHT,N 5810
5270 2AECHESTE KARTE ALS STEUERKARTE ZU LESEN'//) 5820
5280 606 FORMAT(/5X,'FEHLERNACHRICHT'/8X,'ALS RT-KARTE EINGELESENES RECORD 5830
5290 1 TRUG KEINE RT-KENNUNG'/8X,'RECORDTEXT:*****',20A4,'*****'/5X,'AKT 5840
5300 2ION:RECORD WIRD UEBERLESEN.NAECHESTES RECORD WIRD ALS STEUERKARTE E 5850
5310 3INGELESEN'//) 5860
5320 607 FORMAT(/5X,'FEHLERNACHRICHT'/8X,'LESEFEHLER BEIM LESEN EINER STEU 5870
5330 1ERKARTE'/5X,'AKTION:KARTE WIRD UEBERLESEN.ES WIRD VERSUCHT NAECHEST 5880
5340 2E KARTE ALS STEUERKARTE ZU LESEN'//) 5890
5350 608 FORMAT(/5X,'FEHLERNACHRICHT'/8X,'LDFOPN-AUFRUF BLIEB ERFOLGLOS'/5 5900
5360 1X,'AKTION:KEDAK-OPERATIONEN UNTERBLEIBEN.NAECHESTE STEUERKARTE WIRD 5910
5370 2 GESUCHT'//) 5920
5380 610 FORMAT(/5X,'FEHLERNACHRICHT'/8X,'AUSGABE VON SMOOTH ANDERS ALS DU 5930
5390 1RCH PLOT WURDE GEFORDERT.ES WURDE JEDOCH KEIN SCUTINT /XINTPO GEFU 5940
5400 2NDEN'/5X,'AKTION:DIESE AUSGABE ERFOLGT NICHT'//) 5950
5410 611 FORMAT(/5X,'FEHLERNACHRICHT'/8X,'BEIM VERSUCH,EINE KARTE EINZULES 5960
5420 1EN WURDE DAS ENDE DER KARTENEINGABE ERREICHT'/5X,'AKTION:GEWUENSCH 5970
5430 2TE UND NOCH NICHT AUSGEFUEHRTE SMOOTHOPERATIONEN WERDEN DURCHGEFUE 5980
5440 3HRT'//) 5990
5450 612 FORMAT(/5X,'KARTEN KARTEH=' ,I3,' , FMT=' ,20A4/ 6000
5460 15X,'ANZ=' ,I5,' , DATADD=' ,L2,' , CRUCK=' ,L2,' , STANZ=' ,L2,' , PL 6010
5470 XO=' ,L2,' , PLOZEI=' ,I2/ 6020
5480 25X,'KURVE=' ,A4,' , PLOKU=' ,L2,' , CRUKU=' ,L2,' , STAKU=' ,L2,' , DI 6030
5490 XSP=' ,A4/ 6040
5500 35X,'EINHTX=' ,A4,' , EINHTY=' ,A4,' , EINHTD=' ,A4) 6050

```

```

613 FORMAT(5X,'SMOO=' ,A4,' , PLOSM=' ,L2,' , DRUSM=' ,L2,' , STASM=' ,L2,' , 6060
    1SMGEW=' ,L2,' , FISS=' ,L2) 6070
614 FORMAT(5X,'SMOO=' ,E12.3,' , PLOSM=' ,L2,' , DRUSM=' ,L2,' , STASM=' ,L2,' , 6080
    1' , SMGEW=' ,L2,' , FISS=' ,L2) 6090
619 FORMAT(///' &KEDAK PLEH=' ,I3,' , NAMZ=' ,I3,' , NAMEN=' ,A8,2X,A8) 6100
620 FORMAT(//5X,'FEHLERNACHRICHT'/8X,'ZAHL DER KEDAKENERGIEN ZWISCHEN 6110
    1EU UND ED = 0'/5X,'AKTION:KEIN PLOT VON KEDAKWERTEN'//) 6120
621 FORMAT(//5X,'FEHLERNACHRICHT'/8X,'ZAHL DER IM KEDAKAUFRUF ERHALTEN 6130
    1EN WERTE = 0'/5X,'AKTION:KEIN PLOT VON KEDAKWERTEN'//) 6140
622 FORMAT(//' &STAFOR STAFMT=' ,20A4) 6150
623 FCORMAT(//' &OUTINT FMTOUT=' ,20A4/5X,'ZXINT=' ,I5,' , EINHTX=' ,A4,' , A 6160
    1DD=' ,L2,' , XINT=' /(10(1X,E12.5))) 6170
624 FORMAT(//' &RELGEW ZAHLXR=' ,I3,' , EINHTX=' ,A4,' , XR,PR=' /(10(1X,E12 6180
    1.5))) 6190
625 FORMAT(//' &SMOWDH FQU=' ,E13.5,' , ZAHLXR=' ,I3,' , EINHTX=' ,A4,' , NEU 6200
    1ZEI=' ,L2) 6210
626 FCORMAT(5X,'XR,PR=' /(10(1X,E12.5))) 6220
627 FORMAT(//5X,'FEHLERNACHRICHT'/8X,'AUSGABE DER KURVE ANDERS ALS DU 6230
    1RCH PLOT WURDE GEFORDERT.ES WURDE JEDOCH KEIN &OUTINT /XINTPO GEFU 6240
    2NDEN'/5X,'AKTION:DIESE AUSGABE ERFOLGT NICHT'//) 6250
628 FCORMAT(5X,'AUSDRUCK DER AUF XINT INTERPOLIERTEN KURVE') 6260
629 FORMAT(5X,'AUSDRUCK DES AUF XINT INTERPOLIERTEN SMOOTH') 6270
630 FORMAT(5X,'EINGABEDATEN X/Y BZW X/Y/P :') 6280
631 FCORMAT(5X,'SMOOTH WIRD MIT FOLGENDEN WERTEN ANGELAUFEN:' /8X,'FQU= 6290
    1' ,E12.5/8X,'X/Y/P ='//) 6300
632 FCORMAT(//5X,'NACHRICHT'/8X,'DA DIE ANZAHL DER PLNKTE FUER SMOOTH < 6310
    13 IST,WIRD SMOOTH NICHT AUSGEFUEHRT'//) 6320
633 FCORMAT(5X,'NACHRICHT'/8X,'SMOOTH WIRD MIT DEM ZL NP=' ,I2,' GEHOERE 6330
    1NDFM ZEICHEN GEPLOTTET.DIESE ZEICHEN BEFINDEN SICH AM ANFANG UND A 634C
    2M ENDE DER KURVE ') 6350
634 FCORMAT(5X,'NACHRICHT'/8X,'SMOOTH WIRD AUF NEUES BLATT GEPLOTTET.DI 6360
    1E DEM FIT ZUGRUNDE LIEGENDEN PUNKTE WERDEN MIT NP=4 GEZEICHNET') 6370
636 FCORMAT('-----' 6380
    1-----' 6390
    2-----') 6400
637 FORMAT(5X,'NACHRICHT'/8X,'*ENDE*-RECORD WURDE ERREICHT.PROGRAMMAUS 641C
    1FUEHRUNG WIRD BEENDET') 6420
638 FCORMAT(1H1) 6430
639 FCORMAT(///' &SMOWDH ZAHLXR=' ,I3,' , NEUZEI=' ,L2,' , PLOFQU=' ,L2) 6440
640 FORMAT(5X,'EINHTX=' ,A4) 6450
641 FORMAT(//' &OUTINT FMTOUT=' ,20A4/5X,'ZXINT=' ,A4,' , EINHTX=' ,A4,' 6460
    1' , ADD=' ,L2) 6470
642 FORMAT(8X,'DIESE PUNKTE WERDEN ZU DEN SCHON BESTEHENDEN INTERPOLAT 6480
    1IONSPUNKTEN HINZUGEFUEGT') 6490
643 FORMAT(8X,'MIT DEN NEU HINZUGEKOMMENEN PUNKTEN WAERE DEREN ZAHL > 6500
    11000.DIE P(N) MIT N>1000 WERDEN WEGGELASSEN') 6510
644 FORMAT(//5X,'FEHLERNACHRICHT'/8X,'ADD WURDE IN &OUTINT SPEZIFIZIER 6520
    1T.ES LIEGEN BIS JETZT JEDOCH NOCH KEINE INTERPOLATIONSPUNKTE VOR'/ 6530
    28X,'AKTION:DIE PUNKTE WERDEN OHNE AENDERUNG UEBERNOMMEN') 6540
645 FORMAT(5X,'CALCOMPLOT WURDE ABGESCHLOSSEN') 6550
647 FORMAT(//' &OUTINT FMTOUT=' ,20A4/ 6560
    1 5X,'ZXINT=' ,A4,' , PLEH=' ,I2,' , NAMZ=' ,I2,' , NAMEN=' ,2A8/ 6570
    15X,'VON=' ,E14.6,' ,EV' ,5X,'BIS=' ,E14.6,' ,EV'//) 6580
650 FORMAT(//' KONTROLLAUSDRUCK DER GEPLOTTETEN KURVE:'// 6590
    1(10E13.6)) 6600

```

```

651 FORMAT(//5X,'EINHTX=' ,A4,' , ZXINT=' ,I5,' , XINT=' /(10E13.6)) 6610
652 FORMAT(//5X,'ZXINT=' ,I5,' , XINT=' /(10E13.6)) 6620
653 FCORMAT(8X,'FRCM=' ,E13.6,' , TO=' ,E13.6,' IN UNITS OF EINHTX.' ,3 663C
    XX,'REWD=' ,L2) 6640
654 FORMAT(//5X,'&OUTINT ZXINT=' ,A4,' , KARTEH=' ,I2,' , ANZ=' ,I6,' F 6650
    XRCM=' ,E13.6,' , TO=' ,E13.6,' , REWD=' ,L2/ 6660
    1 8X,'FMT=' ,20A4/ 6670
    1 8X,'FMTOUT=' ,20A4/) 6680
655 FCORMAT(///' TESTAUSDRUCK DER SMOOTHKURVE FUER PLOT ODER SIGMA( 6690
    1FISS).'//) 6700
656 FCORMAT('+' ,54X,E14.6) 6710
657 FCORMAT(8X,'FROM=' ,E14.6,' , TO=' ,E14.6/ 6720
    15X,'DRUCK=' ,L2,' , STANZ=' ,L2,' , PLC=' ,L2,' , PLCZEI=' ,I2/ 6730
    25X,'KURVE=' ,A4,' , PLOKU=' ,L2,' , DRUKU=' ,L2,' , STAKU=' ,L2,' , DI 6740
    XSP=' ,A4/ 6750
    35X,'XINTPO=' ,L2,' , FISS=' ,L2) 6760
    END 6770

```

```

SUBROUTINE NEUDA(XINT,DELTA,*) 10
C-----READ DATA POINTS FROM NEUDADA TRANSMISSION FILE. 20
    INTEGER PLEH,ZXINT,WNR,ZAHL,PLOZEI 30
    COMMON/INOUT/KOUT,KIN 40
    INTEGER ACCNRV,BDEH,ACCNR,ANZ 50
    LOGICAL DRUCK,STANZ,PLC,PLOKU,DRUKU,STAKU,PLOSM,DRUSM,STASM,SMGEW, 60
    1 ADD,PLOFQU,REWD,SEQN,CALPJ,DATADD,NEUZEI,TST,XINTPC,FISS,TSTS, 70
    2 FODATA 80
    REAL KURVE,MEDIUM,MEV,NULL,LTR,NEW,NEIN/'NEIN'/ 90
    DIMENSION XINT(1),DELTA(1) 100
    DIMENSION ARR(6),ACONV(10),ACNRV(32) 110
    COMMON XP(1000),YP(1000),ANZ 120
    COMMON/PARM1/ MAXOPT,PLEH,PLCZEI,ZXINT,DPLOT,CALPJ 130
    COMMON/PARM2/ I PLOGR,I STAF0,I SMO,I FMOUT,I SMNP,DISP,EINHTX,EINHTY, 140
    1 EINHTD,XEINHT,YEINHT,DEINHT,LTR,KARTEH,FMTOUT(20),STAFMT(20), 150
    2 FMT(20),DRUCK,STANZ,PLC,KURVE,PLOKU,DRUKU,STAKU,PLOSM,DRUSM, 160
    3 STASM,SMGEW,NEUZEI,XINTPO,PLOFQU,FISS,SEQN,ADD,DATADD,TST, 170
    4 REWD,NANZ,NACD,ZAHL,MEDIUM,WNR, 180
    5 RAST(2),ISMAX,FQU,NAMZ,ZAHLXR,TSTS,FODATA 190
    COMMON/PARM3/SMOO 200
    COMMON/PARM4/FROM,TO 210
    EQUIVALENCE (FROM,EMIN,VON,E1,EUNT),(TO,EMAX,BIS,E2,E0B) 220
    NAMELIST/NEUDAD/BDEH,ACCNR,DRUCK,STANZ,PLC,KURVE,PLOKU,DRUKU,STAKU 230
    1,SMOO,PLOSM,DRUSM,STASM,SMGEW,MEDIUM,KARTEH,FMT,ZAHL,WNR,XEINHT,YE 240
    2INHT,DEINHT,FISS,LTR,SEQN,FROM,TO,XINTPO,PLOZEI,DATADD,TST,DISP 250
    3,TSTS,EMIN,VON,E1,EUNT,EMAX,BIS,E2,ECB,FODATA 260
    DATA ACONV/'0 ' ,1 ' ,2 ' ,3 ' ,4 ' ,5 ' ,6 ' ,7 ' 270
    1,'8 ' ,9 ' / 280
    DATA ACCNRV/32*0/,PERC/'PERC'/ 290
C-----READ CONTROL INPUT LIST. 300
    READ(KIN,NEUDAD) 310
    WRITE(KOUT,636) 320
    WRITE(KOUT,616) BDEH,ACCNR,LTR,FROM,TO,SEQN,DATADD,DRUCK,STANZ,PLC 33C
    1,PLOZEI,KURVE,PLOKU,DRUKU,STAKU,DISP,XEINHT,YEINHT,DEINHT, 340

```

```

2 FISS,XINTPO,FODATA
IF(SMOO.NE.NEIN) GOTO 125
WRITE(KOUT,617) SMOO,PLOSM,DRUSM,STASM,SMGEW,MEDIUM,KARTEH,FMT,
1 ZAHL,WNR
GOTO 126
125 ISMO=1
WRITE(KOUT,618) SMOO,PLOSM,DRUSM,STASM,SMGEW,MEDIUM,KARTEH,FMT,
1ZAHL,WNR
C-----CONVERT DATA SET ID TO HOLLERITH.
C-----AND CHECK CURRENT TAPE POSITION.
126 CALL ACONVE(ACCNR,BRR,ACONV)
BRR(1)=LTR
IF(.NOT.SEQN) GOTO 122
IF(ACCNRV(BDEH).EQ.0) GOTO 122
IF(ACCNRV(BDEH)-ACCNR) 121,122,122
121 BACKSPACE BDEH
GOTO 123
122 REWIND BDEH
C-----FETCH DATA FROM TAPE.
123 CALL BANDDT(BDEH,BRR,NR,SMGEW,DELTA,MEDIUM,ZAHL,FMT,KARTEH,
1 WNR,FROM,TO,FODATA)
IF(ANZ.EQ.0) GOTO 124
ACCNRV(BDEH)=ACCNR
IF(NR.EQ.1) ACCNRV(BDEH)=0
C-----CONVERT TO INTERNAL PHYSICAL UNITS.
CALL ECONV(ANZ,XP,XEINH)
CALL ECONV(ANZ,YP,YEINH)
C-----INTERPOLATION POINTS ARE TO BE READ.
IF(.NOT.XINTPO) GOTO 127
ZXINT=ANZ
CALL STRING(XINT(1),XP(1),4*ZXINT)
IF(IFMOUT.EQ.1) WRITE(KOUT,646)
IFMOUT=1
C-----HANDLE ERROR BARS.
127 IF(.NOT.SMGEW) RETURN
IF(DEINH.EQ.PERC) GOTO 128
CALL ECONV(ANZ,DELTA,DEINH)
RETURN
128 CALL PERCNT(ANZ,YP,DELTA)
RETURN
124 ACCNRV(BDEH)=0
RETURN
C
C-----INTERPOLATION POINTS ARE REQUESTED ONLY.
ENTRY NEUINT(XINT,BDEH,ACCNR,*)
WRITE(KOUT,648) FMTOUT,ZXINT,BDEH,LTR,ACCNR,SEQN,VON,BIS
CALL ACONVE(ACCNR,BRR,ACONV)
BRR(1)=LTR
IF(.NOT.SEQN) GOTO 1695
IF(ACCNRV(BDEH).EQ.0) GOTO 1695
IF(ACCNRV(BDEH)-ACCNR) 1694,1695,1695
1694 BACKSPACE BDEH
GOTO 1696
1695 REWIND BDEH
1696 CALL NEUDEN(BDEH,BRR,VON,BIS,XINT,ZXINT,FODATA,61697)

```

```

350 CALL ECONV(ZXINT,XINT,EINH)
360 WRITE(KOUT,651) EINH,XZINT,(XINT(I),I=1,ZXINT)
370 RETURN
380 1697 IFMOUT=0
390 RETURN
400 616 FORMAT(///' ANEUDAD BDEH=',I3,', ACCNR=',I6,', LTR=',A2,', FROM='
410 X,E12.4,', TO=',E12.4,', SEQN=',L2,', DATADD=',L2/
420 15X,', DRUCK=',L2,', STANZ=',L2,', PLO=',L2,', PLOZEI=',I2/
430 25X,'KURVE=',A4,', PLOKU=',L2,', DRUKU=',L2,', STAKU=',L2,', DI
440 XSP=',A4/
450 35X,'XEINH=',A4,', YEINH=',A4,', DEINH=',A4,', FISS=',L2,
460 X', XINTPO=',L2,', FODATA=',L2)
470 617 FORMAT(5X,'SMOO=',A4,', PLOSM=',L2,', DRUSM=',L2,', STASM=',L2/5X,
480 1'SMGEW=',L2,', MEDIUM=',A4,', KARTEH=',I3,', FMT=',20A4/5X,'ZAHL='
490 2,I5,', WNR=',I3)
500 618 FORMAT(5X,'SMOO=',E12.5,', PLOSM=',L2,', DRUSM=',L2,', STASM=',L2/
510 15X,'SMGEW=',L2,', MEDIUM=',A4,', KARTEH=',I3,', FMT=',20A4/5X,'ZAH
520 2L=',I5,', WNR=',I3)
530 636 FCRMAT(2X,130('-'))
540 646 FORMAT(///' ***** NACHRICHT'/8X,'DIE BSHERIGEN INTERPOLATIONSPUNKT
550 1E WERDEN DURCH DIE NEUDADAPUNKTE ERSETZT'//)
560 648 FCRMAT(/' &OUTINT FMTOUT=',20A4/
570 1 5X,'ZXINT=',A4,', BDEH=',I2,', LTR=',A2,', ACCNR=',I6,', SEQN
580 1=',L2/5X,'VON=',E14.6,5X,'BIS=',E14.6/)
590 651 FORMAT(//5X,'EINH=',A4,' XZINT=',I5,' XINT='/(10E13.6))
600 END
610
620
630
640
650
660 *PROCESS OPTIMIZE(1)
670 SUBROUTINE LESENI(KE,FMT,SMGEW,DELTA,FROM,TO,REWD,*)
680 C-----FORMATTED INPUT READING ROUTINE.
690 COMMON/INOUT/KOUT,KIN
700 DIMENSION XP(1000),YP(1000),TEXT(20),FMT(20),DELTA(1000)
710 INTEGER ANZ
720 LOGICAL SMGEW,REWD
730 COMMON XP,YP,ANZ
740 NPEC=0
750 C-----MAXIMUM IS 1000 DATA POINTS.
760 IF(ANZ.LE.1000) GOTO 1
770 WRITE(KOUT,607)
780 ANZ=1000
790 C-----REWIND TO BE DONE?
800 1 IF(.NOT.REWD) GOTO 6
810 C-----REJECT REWIND ON CONTROL INPUT UNIT.
820 IF(KE.NE.KIN) GOTO 3
830 WRITE(KOUT,601) KIN
840 RETURN
850 3 REWIND KE
860 C-----SEARCH FOR FROM,TO AS SPECIFIED.
870 6 READ(KE,500,END=11,ERR=21) TEXT
880 WRITE(KOUT,600) TEXT
890 FF=-1.E-6*FROM+FROM
FT=1.E-6*TO+TO
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

```

CC 7 I=1,ANZ	260	IF(XP(I).LT.FF) GOTC 46	81C
7 XP(I)=-1.E+10	270	J I=-1	820
4 READ(KE,FMT,END=13,ERR=22) XX	280	GCTO 48	830
NREC=NREC+1	290	46 CONTINUE	840
IF(XX.GT.FT) GOTO 14	300	48 IF(JI.EQ.0) GOTO 50	850
IF(XX.GE.FF) GOTO 5	310	ANZ=ANZ-JI	860
GCTO 4	320	CALL STRING(XP(I),XP(JI+1),4*ANZ)	870
C-----FOUND. BACKSPACE TO BE ON SAFE SIDE.	330	CALL STRING(YP(I),YP(JI+1),4*ANZ)	880
5 BACKSPACE KE	340	CALL STRING(DELTA(I),DELTA(JI+1),4*ANZ)	890
IF(NREC.EQ.1) GOTO 2	350	50 WRITE(KOUT,610) ANZ	900
BACKSPACE KE	360	RETURN	910
GCTO 2	370	500 FORMAT(20A4)	920
C-----READ DATA.	380	600 FORMAT(10X,20A4//)	930
C-----REMEMBER THAT AN EVEN NUMBER OF DATA ITEMS	390	601 FORMAT(//5X,'REWIND AUF',I2,'WURDE VERLANGT. AKTION: REWIND WIRD	940
C-----IS PRESCRIBED FOR EACH RECORD.	400	XNICHT AUSGEFUEHRT.DATENSATZ WIRD NICHT VERARBEITET.//)	950
2 IF(SMGEW) GOTO 30	410	602 FORMAT(//5X,'DATENFILE ENDE AUF EINHEIT',I2,' .KEIN WERT >FROM WUR	960
READ(KE,FMT,END=12,ERR=22) (XP(I),YP(I),I=1,ANZ)	420	XDE GEFUNDEN' /	970
GCTO 40	430	18X,'AKTION: KEINE VERARBEITUNG//)	980
11 WRITE(KOUT,603)	440	603 FORMAT(//5X,'FEHLERNACHRICHT'/8X,'BEIM VERSUCH,DEN DATENTITEL FINZ	990
RETURN 1	450	1ULESEN WURDE DAS DATENFILEENDE ERREICHT'/5X,'AKTION:DATENSATZ WIRD	1000
C-----SEARCH FOR END OF DATA.	460	X NICHT VERARBEITET'//)	1010
12 WRITE(KOUT,604) TEXT	470	604 FORMAT(//5X,'FEHLERNACHRICHT'/8X,'BEIM VERSUCH, INFORMATION ZU'/8X,	1020
WRITE(KOUT,611) (XP(I),YP(I),I=1,ANZ)	480	1'***',20A4,'***'/8X,'EINZULESEN,WURDE DAS ENDE DES DATENFILES ERRE	1030
611 FORMAT(' CONTENTS OF ARRAYS XP AND YP: '(2E15.5))	490	ZICHT'/5X,'AKTION:BISHER ERHALTENE WERTE WERDEN VERARBEITET'//)	1040
DO 15 I=1,ANZ	500	605 FORMAT(//5X,'FEHLERNACHRICHT'/8X,'BEIM VERSUCH,DEN DATENTITEL FINZ	1050
C-----READ IN CASE ERROR BARS ARE TO BE READ ALSO.	510	1ULESEN TRAT EIN LESEFEHLER AUF'/5X,'AKTION:DATENTITEL WIRD UEBERGA	1060
IF(XP(I).EQ.-1.E+10) GOTO 17	520	2NGFN')	1070
IF(XP(I).EQ.0..AND.I.NE.1.AND.XP(I-1).GT.XP(I)) GCTO 17	530	606 FORMAT(//5X,'FEHLERNACHRICHT'/8X,'LESEFEHLER BEIM EINLESEN VON INF	1080
GCTO 15	540	ORMATION ZU'/8X,'***',20A4,'***'/5X,'AKTION:AUSFUEHRUNG DES LETZTE	1090
17 J=I-1	550	2N STEUERBEFFHLS WIRD ABGEBROCHEN.NAECHESTER STEUERBEFFHL WIRD GESUC	1100
GCTO 16	560	3HT')	1110
15 CCNTINUE	570	607 FORMAT(//5X,' ANZ > 1000 NICHT ERLAUBT.DATENSATZ WIRD GEKUEZT'//)	1120
J=ANZ	580	608 FORMAT(//5X,'KEINE DATEN IM ANGEgebenEN INTERVALL'//)	1130
16 ANZ=J	590	609 FORMAT(10X,'SINCE NO DATA WERE READ BEFORE END OF DATA WAS ENCOUNT	1140
IF(ANZ.NE.0) GOTO 40	600	1ERED,NO PROCESSING IS ATTEMPTED.//)	1150
WRITE(KOUT,609)	610	610 FORMAT(//5X,'NUMBER OF DATAPPOINTS TRANSMITTED=',I4)	1160
RETURN 1	620	END	1170
13 WRITE(KOUT,602) KE	630		
RETURN 1	640		
14 WRITE(KOUT,608)	650		
RETURN 1	660		
21 WRITE(KOUT,605)	670		
GCTO 2	680	SUBROUTINE PLOGXY(XP,YP,ANZ,NCUR,NP,NPA,INDZ,NTEXT,	10
22 BACKSPACE KE	690	1XL,YL,RAST,XMAL,XMIL,XMAX,XMIN, SX, YMAL, YMIL, YMAX, YMIN, SY,	20
WRITE(KOUT,606)	700	2ICPLOT,NLGGX,NLGGY, NX, NY, XSKA, YSKA, BSK)	30
RETURN 1	710	C-----PLOTTING ROUTINE FOR LOG. PLOTS.	40
30 READ(KE,FMT,END=12,ERR=22) (XP(I),YP(I),DELTA(I),I=1,ANZ)	720	DIMENSION XDUM(2),YDUM(2),NDIR(10),NSC(10),XSKA(100),YSKA(100),	50
GOTO 40	730	1XP(2),YP(2),XL(2),YL(2),RAST(2),NTEXT(1)	60
C-----SKIP ALL DATA OUTSIDE SPECIFIED INTERVAL (FROM,TO)	740	DATA NDIR/10*2/,NSC/10*1/	70
40 DO 42 I=1,ANZ	750	REAL*8 BSK(200)	80
IF(XP(I).LE.FT) GOTO 42	760	INTEGER ANZ	90
ANZ=I-1	770	DATA XLOG/'XLOG',YLOG/'YLOG'//	100
GCTO 44	780	C-----CONVERT DATA TO LOG.	110
42 CCNTINUE	790	CALL CONLOG(XP,XL,ANZ,RAST(1),XLOG)	120
44 DO 46 I=1,ANZ	800	CALL CONLOG(YP,YL,ANZ,RAST(2),YLOG)	130
		IF(INDZ.EQ.0.OR.INDZ.EQ.10) GOTO 50	140

```

MC=1
DX=80.*SX
DY=80.*SY
IF(INDZ.GT.10) DX=DX+DX
IF(INDZ.GT.10) DY=DY+DY
IF(NX.EQ.0) DX=XMAX-XMIN
IF(NY.EQ.0) DY=YMAX-YMIN
C-----PLOT AXIS.
CALL PLOTA(XL,YL,ANZ,NCUR,NP,1,1,NPA,INDZ,XMAL,XMIL,SX,
1YMAL,YMIL,SY,NTEXT,1DPL0T,
2NLGGX,XMIN,DX,XMAX,4HE9.2,1,-1,+1,NX,
3NLGGY,YMIN,DY,YMAX,4HE9.2,1,+1,-1,NY,0)
IF(INDZ.GT.10) L=10
IF(INDZ.LT.10) L=0
XDUM(1)=XMIL
XDUM(2)=XDUM(1)
YDUM(1)=YMIL
YDUM(2)=YDUM(1)
IF(NX.NE.0) GOTO 20
C-----PLOT SCALE.
CALL SKALA(XMIN,XMIL,XMAL,SX,YMIL,SY,NSKA,NPLOS,LASKA,
1XSKA,YSKA,BSK,RAST(1),XLOG,1)
IF(INDZ.LT.10.OR.RAST(1).EQ.XLOG) GOTO 10
CALL CALSKA(BSK,XSKA,YSKA,NPLOS,LASKA)
C M=10*NPLOS+LASKA
C WRITE(KOUT,600) M,(XSKA(1),YSKA(1),BSK(2*I-1),BSK(2*I),I=1,M)
C 600 FORMAT(I6,' SKALENPUNKTE'/(2E12.3,5X,2A8))
10 IF(NPLOS.EQ.0) GOTO 12
C-----IDENTIFIERS TO BE PLOTTED ALONG AXIS.
DO 11 N=1,NPLOS
11 CALL PLOTA(XDUM,YDUM,2,2,0,1,1,1,L,XMAL,XMIL,SX,YMAL,YMIL,SY,1,
1IDPLOT,0,0,
210,XSKA(10*N-9),YSKA(10*N-9),NDR,NSC,
3BSK(20*N-19),BSK(20*N-17),BSK(20*N-15),BSK(20*N-13),BSK(20*N-11),
4BSK(20*N-9),BSK(20*N-7),BSK(20*N-5),BSK(20*N-3),BSK(20*N-1))
12 IF(LASKA.EQ.0) GOTO 20
CALL PLOTA(XDUM,YDUM,2,2,0,1,1,1,L,XMAL,XMIL,SX,YMAL,YMIL,SY,1,
1IDPLOT,0,0,
2LASKA,XSKA(10*NPL0S+1),YSKA(10*NPL0S+1),NDR,NSC,
3BSK(20*NPL0S+1),BSK(20*NPL0S+3),BSK(20*NPL0S+5),
4BSK(20*NPL0S+7),BSK(20*NPL0S+9),BSK(20*NPL0S+11),
5BSK(20*NPL0S+13),BSK(20*NPL0S+15),BSK(NPLOS*20+17))
20 IF(MC.EQ.2) GOTO 60
IF(NY.NE.0) GOTO 60
MC=MC+1
CALL SKALA(YMIN,YMIL,YMAL,SY,XMIL,SX,NSKA,NPLOS,LASKA,YSKA,XSKA,
1BSK,RAST(2),YLOG,2)
C M=10*NPL0S+LASKA
C WRITE(KOUT,600) M,(XSKA(1),YSKA(1),BSK(2*I-1),BSK(2*I),I=1,M)
GOTO 10
50 CALL PLOTA(XL,YL,ANZ,NCUR,NP,1,1,NPA,INDZ,XMAL,XMIL,SX,
1YMAL,YMIL,SY,1,1DPL0T,0,0,0)
GOTO 60
60 RETURN
END

```

```

150
160
170
180
190
200
210
220
230
240
250
260
270
280
290
300
310
320
330
340
350
360
370
380
390
400
410
420
430
440
450
460
470
480
490
500
510
520
530
540
550
560
570
580
590
600
610
620
630
640
650
660
670
680
690
SUBROUTINE KURVE1(IM,X,Y,M,T,S,S2,A,B,C,D,KENN,*)
C-----CONTROL INTERPOLATION OF DATA POINTS TO CURVE.
C-----KURVE1,KURVE2 FOR AUTOMATIC SELECTION OF
C-----INTERPOLATION POINTS (FOR PLOTS ONLY).
COMMON/INOUT/KOUT
DIMENSION X(IM),Y(IM),T(1000),S(1000),S2(1000),A(1000),B(1000),C(1
1000),D(1000)
GOTO 10
ENTRY KURVE2(IM,X,Y,M,T,S,A,B,C,D,*)
KENN=0
10 NUMP=999/(IM-1)
M=(IM-1)*NUMP+1
FNUMP=FLOAT(NUMP)
J=0
MAX=M-1
DC 6 I=1,MAX,NUMP
J=J+1
T(I)=X(J)
IF(NUMP.EQ.1) GOTO 6
DT=(X(J+1)-X(J))/FNUMP
DC 7 L=2,NUMP
7 T(I+L-1)=T(I+L-2)+DT
6 CONTINUE
T(IM)=X(IM)
GOTO 8
8 IF(KENN.NE.0) GOTO 20
C-----KURVE 3,KURVE4 FOR GIVEN INTERPOLATION POINTS.
ENTRY KURVE3(IM,X,Y,M,T,S,A,B,C,D,NA,NB,*)
C-----CHECK RANGE OF INTERPOLATION POINTS.
CALL PCHECK(IM,X,M,T,NA,NB)
IF(NA.GT.NB) RETURN
K=1
DC 4 I=NA,NB
2 IF(T(I).LT.X(K+1)) GOTO 3
K=K+1
IF(K.NE.IM) GOTO 2
IF(I.FQ.NB) GOTO 5
WRITE(KOUT,604)
RETURN
5 K=K-1
3 DT=T(I)-X(K)
S(I)=A(K)+B(K)*DT+C(K)*DT**2+D(K)*DT**3
4 CONTINUE
RETURN
ENTRY KURVE4(IM,X,Y,M,T,S,S2,A,B,C,D,NA,NB,KENN,*)
20 CALL PCHECK(IM,X,M,T,NA,NB)
IF(NA.GT.NB) RETURN
MM=NB-NA+1
IF(KENN.EQ.1) GOTO 30
IF(IM.LT.4) GOTO 41
CALL SPLINE(IM,X,Y,S2,0,0,MM,T(NA),S(NA),NR,A,E,C,D)

```

```

WRITE(KOUT,600) NR
RETURN
30 IF(IM.LT.2) GOTO 40
CALL LINEAR(IM,X,Y,MM,T(NA),S(NA),NR)
WRITE(KOUT,601) NR
RETURN
40 WRITE(KOUT,602)
RETURN
41 WRITE(KOUT,603)
RETURN
600 FORMAT(//2X,'SPLINENACHRICHT,NR = ',I3)
601 FORMAT(//2X,'NACHRICHT DER LINEAREN INTERPOLATION,NR = ',I3)
602 FORMAT(//5X,'FEHLERACHRICHT'/8X,'KURVE KANN NICHT GELEGT WERDEN.F
INTWEDER IST DIE ANZAHL DER PUNKTE=1 ODER X(ANZ)-X(1) KLEINER ALS E
2 IN PLOTTERSCHRITT'/5X,'AKTION:ES WIRD KEINE KURVE GELEGT UND AUCH
3 NICHT AUSGEGEBEN'//)
603 FORMAT(//5X,'FEHLERACHRICHT'/8X,'SPLINE WURDE GEWUNSCHT.ANZAHL D
IER ANGELIEFERTEN PUNKTE IST < 4'/5X,'AKTION:ES WIRD KEINE SPLINE I
NTERPOLATION VORGENOMMEN'//)
604 FORMAT(//' ***** ERROR IN SUBROUTINE KURVE1.'//)
END

```

```

SUBROUTINE INIT
INTEGER PLEH,ZXINT,WNR,ZAHL,PLOZEI,ZAHLXR
COMMON/INOUT/KOUT,KIN,KPUN,KINC
LOGICAL DRUCK,STANZ,PLC,PLOKU,DRUKU,STAKU,PLOSM,DRUSM,STASM,SMGEW,
1 NEUZEI,XINTPO,PLOFQU,FISS,SEQN,ADD,CALP,CALPJ,DATAADD,TST,REWD
2 ,TSTS,FODATA
REAL NEIN,KURVE,MEDIUM,MEV,NULL,LTR,NFW
DIMENSION FMTA(2)
COMMON/PARM1/ MAXOPT,PLEH,PLOZEI,ZXINT,IDPLOT,CALPJ
COMMON/CLP/ CALP
COMMON/PARM2/ IPLOGR,IStaFO,ISMO,IFMOUT,ISMNP,DISP,EINH TX,EINH TY,
1 EINH TD,XEINH T,YEINH T,DEINH T,LTR,KARTEH,FMTOUT(20),STAFMT(20),
2 FMT(20),DRUCK,STANZ,PLC,KURVE,PLOKU,DRUKU,STAKU,PLOSM,DRUSM,
3 STASM,SMGEW,NEUZEI,XINTPO,PLOFQU,FISS,SEQN,ADD,CATADD,TST,
4 REWD,NANZ,NADD,ZAHL,MEDIUM,WNR,
5 RAST(2),ISMAX,FQU,NAMZ,ZAHLXR,TSTS,FODATA
COMMON/PARM4/ FROM,TD
DATA NEW/'NEW'/,EV/'EV'/,BARN/'BARN'/,MEV/'MEV'/,NULL/'0'/,
1 NEIN/'NEIN'/,BLANK/' '/,FMTA/'(2E1','4.6)'/

```

```

C-----PROGRAM INITIALISATION.
ENTRY INIT1
KCUT=6
KIN=9
KINC=5
KPUN=8
MAXOPT=7
PLEH=1
PLOZEI=-1
ZXINT=0

```

```

520 IDPLOT=0
530 CALPJ=.FALSE.
540 RETURN
550 C
560 C-----RT INITIALISATION.
570 ENTRY INIT2
580 IPLOGR=0
590 IStaFO=0
600 ISMO=0
610 IFMOUT=0
620 ISMNP=0
630 DISP=NEW
640 EINH TX=EV
650 EINH TY=BARN
660 EINH TD=BARN
670 XEINH T=MEV
680 YEINH T=BARN
690 DEINH T=BARN
700 LTR=NULL
710 KARTEH=KIN
720 CC 1001 I=1,2
FMTOUT(I)=FMTA(I)
STAFMT(I)=FMTA(I)
1001 FMT(I)=FMTA(I)
DO 1002 I=3,20
FMTOUT(I)=BLANK
STAFMT(I)=BLANK
1002 FMT(I)=BLANK
DRUCK=.FALSE.
STANZ=.FALSE.
PLC=.TRUE.
KURVE=NEIN
PLOKU=.TRUE.
DRUKU=.FALSE.
STAKU=.FALSE.
PLOSM=.TRUE.
DRUSM=.FALSE.
STASM=.FALSE.
SMGEW=.FALSE.
NEUZEI=.FALSE.
XINTPO=.FALSE.
PLCFQU=.TRUE.
FISS=.FALSE.
SEQN=.TRUE.
CALP=.FALSE.
ADD=.FALSE.
DATAADD=.FALSE.
TST=.FALSE.
REWD=.FALSE.
NANZ=0
NADD=0
FROM=0.
TC=1.E+20
ZAHL=0
MEDIUM=BLANK

```

310  
320  
330  
340  
350  
360  
370  
380  
390  
400  
410  
420  
430  
440  
450  
460  
470  
480  
490  
500  
510  
520  
530  
540  
550  
560  
570  
580  
590  
600  
610  
620  
630  
640  
650  
660  
670  
680  
690  
700  
710  
720  
730  
740  
750  
760  
770  
780  
790  
800  
810  
820  
830  
840  
850

```

WNR=54
RAST(1)=NEIN
RAST(2)=BLANK
ISMAX=0
FCU=0.
NAMZ=2
ZAHXR=0
TSTS=.FALSE.
FODATA=.TRUE.
RETURN
END

SUBROUTINE BANDDT(KB,BLL,NR,SMGEW,DELTA,MEDIUM,ZAHL,FMT,KE,WNR,
1 FROM,TO,FODATA)
C-----READ NEUDADA TAPE.
COMMON/INOUT/KOUT,KIN
DIMENSION BLL(6),BAND(54),XP(1000),YP(1000),DELTA(1000),FMT(20)
REAL MEDIUM
INTEGER ZAHL,WNR
LOGICAL SMGEW,FODATA
COMMON XP,YP,IM
DATA CARD/'CARD',TAPE/'TAPE'/
CALL ERRSET(214,200,-10,1,1)
IERRN=0
NR=0
IM=0
IP=1
C-----SEARCH FOR CORRECT ID.
19 IF(FODATA) READ(KB,101,END=30,ERR=28) (BAND(I),I=1,54)
IF(.NOT.FODATA) READ(KB,120,END=30,ERR=28) (BANC(I),I=1,54)
IF(IERRN.EQ.2) GOTO 81
20 CALL BDVGL(BAND,BLL,&19)
C-----IF FOUND, VERIFY.
WRITE(KOUT,113) (BAND(I),I=1,31)
C-----DETERMINE PHYSICAL UNITS OF ERROR BARS AND RESOLUTION.
WRITE(KOUT,117)
CALL ERRDET(BAND(33),IRTY)
WRITE(KOUT,112)
CALL ERRDET(BAND(39),IRTY)
IF(.NOT.SMGEW) GOTO 18
IF(MEDIUM.EQ.CARD) GOTO 18
IF(MEDIUM.EQ.TAPE) GOTO 40
WRITE(KOUT,102)
SMGEW=.FALSE.
GOTO 18
40 XX=BAND(48)
C-----SEARCH FOR FROM,TO OR END OF THIS ID.
IF(XX.LT.FROM) GOTO 21
IF(XX.GT.TO) GOTO 31
IM=IM+1
IF(IM.GT.1000) GOTO 92
XP(IM)=BAND(48)

```

```

860 YP(IM)=BAND(53) 410
870 DELTA(IM)=BAND(WNR) 420
880 21 IF(FODATA) READ(KB,101,END=27,ERR=25) (BAND(I),I=1,54) 430
890 IF(.NOT.FODATA) READ(KB,120,END=27,ERR=25) (BANC(I),I=1,54) 440
900 IP=2 450
910 IF(IERRN.EQ.2) GOTO 81 460
920 22 CALL BDVGL(BAND,BLL,&27) 470
930 GOTO 40 480
940 18 XX=BAND(48) 490
950 IRTYP=0 500
960 IF(XX.LT.FROM) GOTO 23 510
IF(XX.GT.TO) GOTO 31 520
IM=IM+1 530
IF(IM.GT.1000) GOTO 92 540
XP(IM)=BAND(48) 550
YP(IM)=BAND(53) 560
23 IF(FODATA) READ(KB,101,END=27,ERR=29) (BAND(I),I=1,54) 570
IF(.NOT.FODATA) READ(KB,120,END=27,ERR=29) (BANC(I),I=1,54) 580
IP=3 590
IF(IERRN.EQ.2) GOTO 81 600
24 CALL BDVGL(BAND,BLL,&27) 610
GOTO 18 620
27 WRITE(KOUT,107) KB,BLL,IM 630
IF(IRTY.NE.2) GOTO 26 640
IF(IM.EQ.0) GOTO 26 650
DC 32 I=1,IM 660
32 DELTA(I)=DELTA(I)*YP(I)*.01 670
GOTO 26 680
C-----ERROR EXITS. 690
28 WRITE(KOUT,108) BLL 700
IP=1 710
GOTO 80 720
25 WRITE(KOUT,116) BLL 730
IP=2 740
GOTO 80 750
29 WRITE(KOUT,109) BLL 760
IP=3 770
GOTO 80 780
31 IF(IM.NE.0) GOTO 27 790
NR=1 800
WRITE(KOUT,106) KB,BLL 810
GOTO 26 820
30 WRITE(KOUT,110) BLL 830
NR=1 840
C-----IF ERROR BARS ARE DESIRED, CONTINUE. ELSE RETURN. 850
26 IF(.NOT.SMGEW) RETURN 860
IF(MEDIUM.EQ.TAPE) RETURN 870
C-----ERROR BARS ARE TO BE READ FROM CARDS. 880
IF(ZAHL.EQ.IM) GOTO 50 890
WRITE(KOUT,103) IM,ZAHL 900
50 READ(KE,FMT,END=90,ERR=91) (DELTA(I),I=1,ZAHL) 910
C-----CHECK IF NUMBER OF DATA AGREE. 920
IF(ZAHL.GT.IM.OR.ZAHL.EQ.IM) RETURN 930
IANF=ZAHL+1 940
DO 51 I=IANF,IM 950

```



51 DELTA(I)=DELTA(ZAHL)	960		
RETURN	970		
80 WRITE(KOUT,114)	980		
READ(KB)	990		
IERRN=2	1000		
GCTO(19,21,23),IP	1010		
81 WRITE(KOUT,115) BAND	1020	C-----	SLROUTINE TSTPRO(ISMAX,XS,YS,P,A,FQU) 10
IERRN=0	1030	-----TEST PRINTOUT.	20
GCTO(20,22,24),IP	1040	DIMENSION P(1),A(1),XS(1),YS(1)	30
90 WRITE(KOUT,104)	1050	COMMON/INOUT/KOUT	40
SMGEW=.FALSE.	1060	WRITE(KOUT,604)	50
RETURN	1070	604 FCRMAT('1')	60
91 WRITE(KOUT,105)	1080	WRITE(KOUT,600)	70
SMGEW=.FALSE.	1090	600 FCRMAT( 50X,'TESTPRINTOUT OF SMOOTH RESULTS'////)	80
BACKSPACE KE	1100	WRITE(KOUT,602)	90
RETURN	1110	DDX=0	100
92 WRITE(KOUT,111)	1120	DATA LLM/50/	110
IM=IM-1	1130	L=0	120
GOTO 27	1140	DC 100 I=1,ISMAX	130
101 FCRMAT(47A1,7E12.5,1X)	1150	IF(L.LT.LLM) GOTO 10	140
120 FCRMAT(47A1,7A4)	1160	WRITE(KOUT,604)	150
102 FORMAT(/5X,'FEHLERNACHRICHT'/8X,'SMGEW=T GEFUNDEN.MEDIUMANWEISUNG	1170	WRITE(KOUT,602)	160
1 FEHLT'/5X,'AKTION:SMGEW=F GESETZT.DADURCH WERDEN DIE GEWICHTE AUT	1180	L=0	170
2OMATISCH = 1'////)	1190	10 DX=ABS(YS(I)-A(I))	180
103 FCRMAT(/5X,'FEHLERNACHRICHT'/8X,'DIE ANZAHL DER AUF NEUDADA GEFUN	1200	DDP=DX/P(I)	190
1DENEN DATENPUNKTE IM=',I5/8X,'STIMMT NICHT MIT DER IN DER NEUDAD -	1210	DCX=DDX+DDP**2	200
2 KARTE ANGEGEBENEN ANZAHL ZAHL=',I5,'UEBEREIN'/5X,'AKTION:FALLS IM	1220	601 FCRMAT(1P6E15.5)	210
3>ZAHL ERHALTEN DIE LETZTEN IM-ZAHL DATENPUNKTE DAS GEWICHT DES ZAH	1230	602 FORMAT( T8,'ENERGY',T19,'EXPERIMENTAL',T36,'FITTED',T49,'ABS(FI	220
4L-TEN PUNKTES'/8X,'FALLS IM<ZAHL WERDEN DIE LETZTEN ZAHL-IM GEWICH	1240	IT-EXP)',T65,'EXP.ERROR',T79,'D(FIT)/C(EXP)'/)	230
5TE IGNORIERT'////)	1250	WRITE(KOUT,601) XS(I),YS(I),A(I),DX,P(I),DDP	240
104 FCRMAT(/5X,'FEHLERNACHRICHT'/8X,'BEIM EINLESEN DER SMOOTHGEWICHTE	1260	100 L=L+1	250
1 WURDE DAS ENDE DES DATENFILES ERREICHT'/5X,'AKTION:SMGEW=FALSE GE	1270	WRITE(KOUT,603) FQU,DDX	260
2SETZT'////)	1280	603 FCRMAT(/10X,'FQU=',1PE13.5,5X,'RESIDUAL OF FIT=',1PE13.5'////)	270
105 FORMAT(/5X,'FEHLERNACHRICHT'/8X,'BEIM EINLESEN DER SMOOTHGEWICHTE	1290	RETURN	280
1 TRAT EIN LESEFEHLER AUF'/5X,'AKTION:SMGEW=FALSE GESETZT'////)	1300	END	290
106 FORMAT(' *****FEHLERNACHRICHT'/8X,'ANZAHL DER DATENPUNKTE AUF NEUD	1310		
1ADABAND',I3,' MIT ACCNR=',6A1/	1320		
28X,' = 0 ZWISCHEN DEN ANGEGEBENEN ENFRGIEGRENZEN')	1330		
107 FORMAT(5X,'ANZAHL DER AUF DEM NEUDADABAND ',I3,' GEFUNDENEN DATEN	1340		
1 MIT ACCESSNUMMER ',6A1,' = ',I5)	1350	C-----DATA POINTS WITH EQUAL ENERGIES ARE NOT PERMITTED FOR	10
108 FORMAT(2X,6A1,' FEHLER IN 19 READ')	1360	C-----SMOOTH. THESE DATA POINTS ARE COMBINED TO A WEIGHTED AVERAGE	20
109 FCRMAT(2X,6A1,' FEHLER IN 23 READ')	1370	C-----IN THIS ROUTINE.	30
110 FCRMAT(2X,6A1,'NICHT GEFUNDEN,BANDENDE ERREICHT')	1380	SLROUTINE EQUFN(MAX,X,Y,P)	40
111 FORMAT(/' ANZAHL DER UEBERTRAGENEN DATENPUNKTE HAT DIE MAXIMALZA	1390	COMMON/INOUT/KOUT	50
1HL 1000 ERREICHT'/' REST WIRD ABGESCHNITTEN.'////)	1400	DIMENSION X(1),Y(1),P(1)	60
112 FORMAT(/' ERRORTYPE OF Y-VALUES:')	1410	DATA VH/1.00001/	70
113 FCRMAT(/5X,31A1/)	1420	IF(MAX.LT.2) GOTO 200	80
114 FCRMAT(/5X,'NACHRICHT'/8X,' RECORDS WURDEN AUSGELASSEN WEGEN IHC	1430	MT=0	90
1218-ERROR'/8X,'ES IST ZU UEBERPRUEFFEN,OB DARUNTER KEINES DER BEIM	1440	MP=0	100
12AUFRUF GEFUENSCHTEN IST'/8X,'ZU DIESEM ZWECK WIRD VERSUCHT,DAS NAC	1450	DC 100 I=2,MAX	110
13HFOLGENDE RECORD AUSZUDRUCKEN')	1460	IF(X(I).GT.VH*X(I-1)) GOTO 90	120
115 FCRMAT(5X,'RECORD LAUTET: '/1X,47A1,7E12.5'////)	1470	MF=MP+1	130
116 FORMAT(2X,6A1,' FEHLER IN 21 READ')	1480	GCTC 99	140
117 FORMAT(/' ERRORTYPE OF X-VALUES:')	1490	80 IF(MT.EQ.0) WRITE(KOUT,600)	150
END	1500	600 FCRMAT(/20X,'DATAPOINTS WITH EQUAL ENERGIES ARE NOT ALLOWED FOR	160
		1SMOOTH'//T20,'FOR THE FOLLOWING ENERGIES DATAPCINTS WERE COMBINED	170
		1FCR THIS REASON.'////)	180

```

WRITE(KOUT,601) X(I-MP)
601 FCRMAT(10X,1PE13.5)
MT=MT+MP
MP=MP+1
YY=0
PP=0
PS=0
DO 81 J=1,MP
YY=YY+Y(I-J)/P(I-J)
PP=PP+1./P(I-J)
81 PS=PS+P(I-J)**2
Y(I-MT-1)=YY/PP
P(I-MT-1)=SQRT(PS)/MP
MP=0
GOTO 92
90 IF(MP.NE.0) GOTO 80
52 Y(I-MT)=Y(I)
X(I-MT)=X(I)
P(I-MT)=P(I)
GOTO 100
99 IF(I.NE.MAX) GOTO 100
GOTO 80
100 CONTINUE
MAX=MAX-MT
200 RETURN
END

```

```

SUBROUTINE LESEN(KE,FMT,RWD,N1,N2,XMIN,XMAX,X,*)
C-----READ INTERPOLATION ENERGIES FROM FORMATTED INPUT, E.G. CARDS.
DIMENSION X(1),FMT(1),TEXT(20)
COMMON/INOUT/KOUT,KIN
LCGICAL RWD
NREC=0
IF(N2.LE.1000) GOTO 1
WRITE(KOUT,607)
607 FORMAT(//'* ANZ > 1000 NICHT ERLAUBT.DATENSATZ WIRD GEKUEZT'//)
N2=1000
C-----REWIND? IF ON CONTROL INPUT, REJECT.
1 IF(.NOT.RWD) GOTC 6
IF(KE.NE.KIN) GOTO 3
WRITE(KOUT,601) KIN
601 FORMAT(//5X,'REWIND AUF',I2,'WURDE VERLANGT. AKTION: REWIND WIRD
NICHT AUSGEFUEHRT.DATENSATZ WIRD NICHT VERARBEITET.'//)
RETURN 1
3 REWIND KE
6 READ(KE,500,END=11) TEXT
C-----REMEMBER, THAT COMMENT MUST PRECEDE.
500 FCRMAT(20A4)
WRITE(KOUT,600) TEXT
600 FORMAT(//'* COMMENT=',2CA4)
FF=XMIN*0.999999
FT=XMAX*1.000001

```

```

190 DC 7 I=1,N2
200 7 X(I)=-1.E+10
210 C-----SEARCH FOR REQUESTED INTERVAL.
220 C-----ALLOWANCE FOR MACHINE INTERNAL INACCURACIES
230 C-----WAS PROVIDED.
240 4 READ(KE,FMT,END=11) XX
250 NREC=NREC+1
260 IF(XX.GT.FT) GOTO 14
270 IF(XX.GE.FF) GOTO 5
280 GOTO 4
290 5 BACKSPACE KE
300 IF(NREC.EQ.1) GOTO 2
310 BACKSPACE KE
320 GOTO 2
330 2 READ(KE,FMT,END=12) (X(I),I=1,N2)
340 GOTO 40
350 11 WRITE(KOUT,603) KE
360 603 FCRMAT(//'* END OF DATA ON UNIT=',I3,'. DATASET IS NOT TRANSMI
370 1TTED'//)
380 RETURN 1
390 C-----END OF DATA. TRANSMIT REST.
400 12 WRITE(KOUT,604) KE
410 604 FCRMAT(//5X,'END OF DATA ON UNIT ',I3,' ENCOUNTERED.DATA ALREADY R
420 LEAD ARE USED.'//)
430 DC 15 I=1,N2
440 IF(X(I).NE.-1.E+10) GOTO 15
J=I-1
GOTO 16
15 CONTINUE
J=N2
16 N2=J
10 IF(N2.NE.0) GOTO 40
20 WRITE(KOUT,609)
30 609 FORMAT(5X,'SINCE NO DATA WERE READ ,NO PROCESSING IS ATTEMPTED.'//
40 1)
50 RETURN 1
60 C-----ERROR: NO DATA FOUND.
70 14 WRITE(KOUT,608)
80 608 FORMAT(//5X,'KEINE DATEN IM ANGEGEBENEN INTERVALL'//)
90 RETURN 1
100 DO 42 I=1,N2
110 IF(X(I).LE.FT) GOTO 42
120 N2=I-1
130 GOTO 44
140 42 CONTINUE
150 44 DO 46 I=1,N2
160 IF(X(I).LE.FF) GOTO 46
170 JI=I-1
180 GOTO 48
190 46 CONTINUE
200 48 IF(JI.EQ.0) GOTO 50
210 N2=N2-JI
220 CALL STRING(X(1),X(JI+1),4*N2)
230 50 N1=N2
240 WRITE(KOUT,602) ANZ
250
260
270
280
290
300
310
320
330
340
350
360
370
380
390
400
410
420
430
440
450
460
470
480
490
500
510
520
530
540
550
560
570
580
590
600
610
620
630
640
650
660
670
680
690
700
710
720
730
740
750
760
770
780
790
800

```

```

602 FORMAT(/ /5X,'NUMBER OF DATAPPOINTS TRANSMITTED=',I4)
      RETURN
      END

      SUBROUTINE SKALA(XMTN,XMIL,XMAL,SX,YMIL,SY,NSKA,NPLOS,LASKA,
1XSKA,YSKA,BSK,RAST,SLOG,KONTR)
C-----PROVIDE SCALING FOR PLOTS.
      DIMENSION XSKA(2),YSKA(2),BSK(2),DX(2),DY(2),DSKA(9),DXL(2),DYL(2),
1,CYZW(2)
      REAL*8 POINT1/'./',POINT2/'..',POINT3/'...'/,
1BLANK/'./',BSK,ZERO/'0.',BZERO/'0./'
      DATA DX/30.,8./,DY/24.,56./,DXL/12.,8./,DYL/28.,56./,DYZW/12.,36./
      DATA DSKA/.0458,.3010,.1761,.1250,.0969,.0792,.0669,.0580,.0511/
      LOGICAL*1 BV(10),BW(7),PLUS/'+'/,PT/'./',MINUS/'-'/'
      INTEGER*2 TST,TSTB/'E'/
      LOGICAL*1 BL/'./'
      INTEGER*2 INTNUM(11)/'0','1','2','3','4','5','6','7','8',
1'9','0'/,
      2NUMEX,NUMAX,ST/'**'/
      EQUIVALENCE (NUMEX,BV(9)),(NUMAX,BV(7))
      IF(RAST.EQ.SLOG) GOTO 3
C-----SCALE FOR LIN. AXIS.
      XSKA(1)=XMIL
      NSKA=0
1 NSKA=NSKA+1
      IF(NSKA.GE.100) GOTO 11
      XSKA(NSKA+1)=XSKA(NSKA)+80.*SX
      IF(XSKA(NSKA+1).LE.XMAL) GOTO 1
11 NPLOS=NSKA/10
      LASKA=NSKA-NPLOS*10
      DO 2 I=1,NSKA
      J=2*I-1
      BSK(J)=BLANK
      BSK(J+1)=POINT1
      CALL CONVX(XSKA(I),BV(1),5HE10.3)
      BV(1)=BV(4)
      BV(2)=BV(3)
      DC 18 K=3,6
18 BV(K)=BV(K+2)
      CALL STRING(TST,BV(7),2)
      IF(TST.EQ.TSTB) BV(6)=PLUS
      BV(9)=BL
      IF(TST.EQ.TSTB) GOTO 21
      DO 19 K=1,10
      IF(NUMEX.NE.INTNUM(K)) GOTO 19
      NUMEX=INTNUM(K+1)
      GOTO 20
19 CONTINUE
24 NUMEX=ST
      GOTO 20
21 DC 22 K=1,9
      IF(NUMEX.NE.INTNUM(K)) GOTO 22

```

```

810 IF(K.EQ.1) GOTO 23
820 NUMEX=INTNUM(K-1)
830 GOTO 20
23 NUMEX=INTNUM(2)
      BV(6)=MINUS
      GOTO 20
22 CONTINUE
      GOTO 24
20 BV(7)=BV(10)
      BV(8)=PT
      CALL STRING(BSK(J),BV(1),8)
      IF(BSK(J).EQ.ZERO) BSK(J)=BZERO
      XSKA(I)=XSKA(I)-DX(KONTR)*SX
2 YSKA(I)=YMIL-DY(KONTR)*SY
      GOTO 10
C-----SCALE IS LOG. SCALING DETERMINED BY LENGTH CF DECADE.
3 XSKA(1)=XMIL-DXL(KONTR)*SX
      YSKA(1)=YMIL-DYL(KONTR)*SY
      NSKA=0
4 NSKA=NSKA+1
      IF(NSKA.GE.100) GOTO 41
      XSKA(NSKA+1)=XSKA(NSKA)+DSKA(MOD(NSKA,9)+1)
      YSKA(NSKA+1)=YSKA(1)
      IF(XSKA(NSKA+1).LE.XMAL) GOTO 4
41 NPLOS=NSKA/10
      LASKA=NSKA-NPLOS*10
      JEXP=0
      DO 6 I=1,NSKA,9
      J=2*I-1
      BSK(J)=POINT2
      BSK(J+1)=BLANK
      CALL CONVX(XMIN*(10.**JEXP),BV(1),4HE8.1)
      BV(1)=BV(4)
      BV(2)=BV(3)
      BV(3)=BV(5)
      BV(4)=BV(6)
      CALL STRING(TST,BV(5),2)
      IF(TST.EQ.TSTB) BV(4)=PLUS
      BV(7)=BL
      IF(TST.EQ.TSTB) GOTO 45
      DO 43 K=1,10
      IF(NUMAX.NE.INTNUM(K)) GOTO 43
      NUMAX=INTNUM(K+1)
      GOTO 44
43 CONTINUE
48 NUMAX=ST
      GOTO 44
45 DC 46 K=1,9
      IF(NUMAX.NE.INTNUM(K)) GOTO 46
      IF(K.EQ.1) GOTO 47
      NUMAX=INTNUM(K-1)
      GOTO 44
47 NUMAX=INTNUM(2)
      BV(4)=MINUS
      GOTO 44
490
500
510
520
530
540
550
560
570
580
590
600
610
620
630
640
650
660
670
680
690
700
710
720
730
740
750
760
770
780
790
800
810
820
830
840
850
860
870
880
890
900
910
920
930
940
950
960
970
980
990
1000
1010
1020
1030

```

```

46 CONTINUE
GCTO 48
44 BV(5)=BV(8)
CALL STRING(BSK(J),BV(1),5)
DC 5 N=1,8
BSK(J+2*N)=POINT3
BSK(J+2*N+1)=BLANK
YSKA(I+N)=YSKA(I+N)+DYZH(KONTR)*SY
5 CALL CONVX(N+1,BSK(J+2*N),2HI2)
JEXP=JEXP+1
6 CONTINUE
GCTO 10
10 RETURN
END

SUBROUTINE KEDSUB(NR,NAMZ,NAME,EUNT,EOB,XINTPC,XINT,ZXINT,IFMOUT,
1*)
C-----READ KEDAK DATA.
REAL*8 NAME(1),NAMEN(3)
COMMON/INOUT/KOUT
INTEGER ANZ,ZXINT
LOGICAL XINTPC
DIMENSION NARG(3),ARG(2),XP(1000),YP(1000),XINT(ZXINT)
COMMON XP,YP,ANZ
DC 1 I=1,NAMZ
1 NAMEN(I)=NAMF(I)
NR=0
ANZ=0
NARG(1)=NAMZ
C-----LOCATE DATA.
CALL LDFLOC(NR,NARG,NAMEN,ARG)
IF(NR.EQ.0) GOTO 5
3 IF(ARG(1).LT.EUNT) GOTO 2
IF(ARG(1).GT.EOB) GOTO 4
ANZ=ANZ+1
XP(ANZ)=ARG(1)
YP(ANZ)=ARG(2)
IF(ARG(1).EQ.EOB) GCTO 4
IF(ANZ.LT.1000) GOTO 2
WRITE(KOUT,604)
GOTO 4
C-----LOOP TO SEARCH FOR ENERGY INTERVAL.
2 CALL LDFNXT(NR,NARG,NAMEN,ARG)
IF(NR.EQ.0) GOTO 6
GCTO 3
4 WRITE(KOUT,603) ANZ
IF(XINTPC) GOTO 10
RETURN
5 WRITE(KOUT,601) NAMZ,(NAME(I),I=1,2)
IF(NAMZ.GE.3) WRITE(KOUT,606) NAME(3)
WRITE(KOUT,602)
RETURN 1

```

```

1040 6 WRITE(KOUT,601) NAMZ,(NAME(I),I=1,2) 380
1050 IF(NAMZ.GE.3) WRITE(KOUT,606) NAME(3) 390
1060 WRITE(KOUT,608) ANZ 400
1070 RETURN 410
1080 C-----INTERPOLATION ENERGIES WERE REQUESTED ALSO. 420
1090 10 IF(IFMOUT.EQ.1) GOTO 12 430
1100 9 IFMOUT=1 440
1110 ZXINT=ANZ 450
1120 DO 11 I=1,ZXINT 460
1130 11 XINT(I)=XP(I) 470
1140 RETURN 480
1150 12 WRITE(KOUT,605) 490
1160 GCTO 9 500
1170 601 FORMAT(/5X,'MESSAGE FROM DATARETRIEVAL FROM KEDAK FOR:/' 510
112X,'NUMBER OF NAMES=',I2/ 520
1 12X,'NAMES=',A8,2X,A8) 530
602 FCRMAT(8X,'NO ENTRY ON KEDAK FOR THIS DATA TYPE'//) 540
603 FORMAT(5X,'ANZAHL DER VCN KEDAK GEHOLTEN DATENPUNKTE = ',I5) 550
604 FORMAT(/5X,'NACHRICHT'/8X,'BEIM KEDAKAUFRUF WURDEN 1000 DATENPUNK 560
1TE EINGELESEN.DAMIT IST DAS ZUR VERFUEGUNG STEHENDE FELD GEFUELLT' 570
2/5X,'AKTION:KEDAK EINLESEN WIRD ABGEBROCHEN.BISHER ERHALTENE WERTE 580
3 WERDEN VERARBEITET'//) 590
605 FCRMAT(/5X,'NACHRICHT'/8X,'DIE UEBER 8OUTINT EINGELESENEN INTERPO 600
1LATIONSPUNKTE VERLIEREN IHRE GUELTIGKEIT.SIE WERDEN DURCH DIE KEDA 610
2KENERGIEEN ERSETZT') 620
606 FORMAT('+',36X,E14.6) 630
607 FCRMAT(8X,'EO OR EU > HIGHEST ENERGY FOR THIS TYPE') 640
608 FORMAT(8X,'NUMBER OF DATAPOINTS FOUND =',I5//) 650
END 660

SUBROUTINE DRUCK3(IM,X,Y,P) 10
C-----PRINT DATA POINTS PLUS ERROR BARS. 20
C-----60 LINES PER PAGE, 3 TIMES 3 COLUMNS. 30
DIMENSION X(IM),Y(IM),P(IM) 40
COMMON/INOUT/KCUT 50
NUMSPA=3 60
GCTO 10 70
ENTRY DRUCKX(X,Y,IM) 80
NUMSPA=4 90
10 JM=IM/NUMSPA 100
JMS=JM*NUMSPA 110
IF(JMS.LT.IM) JM=JM+1 120
IMS=0 130
1 IF(JM.GT.60) GOTO 3 140
JAN=IMS+1 150
JMS=IMS+JM 160
IMS=IM 170
IF(NUMSPA.EQ.3) GOTO 31 180
DO 2 I=JAN,JMS 190
2 WRITE(KOUT,601) (X(J),Y(J),J=I,IMS,JM) 200
RETURN 210
31 DC 32 I=JAN,JMS 220

```

```

32 WRITE(KOUT,602) (X(J),Y(J),P(J),J=I,IMS,JM)
RETURN
3 JAN=IMS+1
JMS=IMS+60
IMS=IMS+60*NUMSPA
IF(NUMSPA.EQ.3) GOTO 33
DC 4 I=JAN,JMS
4 WRITE(KOUT,601) (X(J),Y(J),J=I,IMS,60)
5 JM=JM-60
WRITE(KOUT,600)
GOTO 1
33 DO 34 I=JAN,JMS
34 WRITE(KOUT,602) (X(J),Y(J),P(J),J=I,IMS,60)
GOTO 5
600 FCRMAT(1H1)
601 FORMAT (4(5X,2(1PE13.5)))
602 FCRMAT (3(5X,3(1PE13.5)))
END

SUBROUTINE ADF(XL,YL,Z6,NADD,NANZ,ANZ,XP,YP,DELTA,DATAADD,*)
C-----APPEND DATA TO EXISTING SET.
COMMON/INOUT/KOUT
DIMENSION XL(1),YL(1),Z6(1),XP(1),YP(1),DELTA(1)
LOGICAL DATAADD
INTEGER ANZ
IF(NADD.NE.1) GOTO 1290
IF(ANZ+NANZ.LE.1000) GOTO 129
WRITE(KOUT,649)
NANZ=1000-ANZ
129 CALL STRING(XP(ANZ+1),XL(1),4*NANZ)
CALL STRING(YP(ANZ+1),YL(1),4*NANZ)
CALL STRING(DELTA(ANZ+1),Z6(1),4*NANZ)
ANZ=ANZ+NANZ
NADD=0
1290 IF(.NOT.DATAADD) RETURN
NADD=1
DC 1291 J=1,ANZ
XL(J)=XP(J)
YL(J)=YP(J)
1291 Z6(J)=DELTA(J)
NANZ=ANZ
RETURN 1
649 FORMAT(//' ADD=T WAR ANGEGEBEN.GESAMTZAHL DER PUNKTE WIRD DADURCH
1 > 1000.//' ERSTER DATENSATZ WIRD GEKUEZT'//)
END

```

```

230 SUBROUTINE ASSIGN(S,T,J,NR,M) 10
240 C-----COMPARE HOLLERITH DATA. 20
250 DIMENSION S(2),T(2,M) 30
260 NR=0 40
270 DO 1 I=1,2 50
280 IF(S(I).NE.T(I,J)) GOTO 2 60
290 1 CONTINUE 70
300 NR=1 80
310 2 RETURN 90
320 END 100
330
340
350
360
370

```

```

380 SUBROUTINE BCVGL(B,T,*) 10
390 C-----COMPARE NEUDADA ID. 20
400 DIMENSION B(54),T(6) 30
DC 1 J=1,6 40
JB=41+J 50
IF(B(JB).NE.T(J)) GOTO 2 60
1 CONTINUE 70
RETURN 80
2 RETURN 90
END 100

```

```

SUBROUTINE ACCNVE(ACCNR,BRR,ACCNV) 10
C-----CONVERT INTEGER ID TO HOLLERITH. 20
DIMENSION BRR(6),ACCNV(10),IC(5) 30
INTEGER ACCNR 40
IZ=ACCNR 50
ID(1)=1 60
DO 1 K=2,5 70
1 ID(K)=ID(K-1)*10 80
DO 2 I=1,5 90
K=6-I 100
JZ=IZ/ID(K) 110
BRR(I+1)=ACCNV(JZ+1) 120
2 IZ=IZ-JZ*ID(K) 130
RETURN 140
END 150

```

```

SUBROUTINE XCORR 10
C-----PLOTS DOES NOT ALLOW PLOTTING OF A SINGLE DATA POINT. 20
C-----HERE A POINT IS ADDED FOR PLOT ONLY WHICH IS REMOVED LATER. 30
DIMENSION XP(1000),YP(1000) 40
COMMON XP,YP,IM 50
COMMON/PAR/XMAX,SX,YMAX,SY 60
XP(2)=XP(1) 70
YP(2)=YP(1) 80

```

```

IM=2
RETURN
END

SUBROUTINE SUB1(INDZ,RAST,NEIN,XLOG,YLOG,KREX,KREY,*)
C-----SET FLAGS FOR PLOT SCALES.
COMMON/CLP/CALP
LOGICAL CALP
REAL NEIN
DIMENSION RAST(2)
IF(INDZ.EQ.0.OR.INDZ.EQ.10) RETURN1
INDZ=0
IF(CALP) INDZ=10
IF(RAST(1).EQ.NEIN) RETURN1
KREX=1
IF(RAST(1).EQ.XLOG) KREX=-1
KREY=1
IF(RAST(2).EQ.YLOG) KREY=-1
RETURN
END

SUBROUTINE QFISS(MAX,T,S,F)
C-----CALCULATE FISSION SPECTRUM AVERAGE.
C-----AN ANALYTICAL FISSION SPECTRUM IS USED BY DEFAULT.
C-----BUT BY REPLACING CHIF THE USER MAY SUPPLY HIS OWN
C-----SPECTRUM.
DIMENSION T(MAX),S(MAX),F(MAX)
COMMON/INOUT/KOUT
CHIF(F)=.4527*EXP(-F/.965E+6)*SINH(SQRT(2.29E-6*E))
VSUM=0.
SLM=0.
SUMCHI=0.
WRITE(KOUT,600)
600 FORMAT(/5X,' ES BEDEUTEN IM NACHFOLGENDEN AUSDRUCK:/'
A10X,' E=ENERGIE'/
11X,' QF=MIT SPALTSPEKTRUM GEMITTELTER QUERSCHNITT ZWISCHEN DIESER
2 ENERGIE UND DER VORHERGEHENDEN'//
36X,'E',11X,'QF'/
45X,'0. ')
IF(T(1).LE.0.) GOTO 9
DT=T(1)/100.
X1=0.
X2=DT
DO 5 I=1,100
SUMCHI=SUMCHI+DT*(CHIF(X2)+CHIF(X1))
X1=X2
5 X2=X1+DT
9 F(1)=S(1)*CHIF(T(1))
DO 10 I=2,MAX

```

```

90 F(I)=S(I)*CHIF(T(I))
100 SUMCHI=SUMCHI+(CHIF(T(I))+CHIF(T(I-1)))*(T(I)-T(I-1))
110 SUM=SUM+(F(I)+F(I-1))*(T(I)-T(I-1))
IF((I/16)*16.NE.I) GOTO 10
DSUM=(SUM-VSUM)*.5E-3
VSUM=SUM
CHIE=CHIF(T(I))
WRITE(KOUT,601) T(I),DSUM,CHIE
10 CONTINUE
601 FORMAT(1X,E12.3,1X,E12.3,' CHIF=',E10.3)
SUM=SUM*.5E-3
WRITE(KOUT,602) SUM
602 FORMAT(//' QFISS(CHIF(U235))=',E12.4,'MILLIBARN')
SUMCHI=SUMCHI*.5E-6
WRITE(KOUT,603) SUMCHI
603 FORMAT(//' SUMCHI=',5X,F12.6)
RETURN
END

SUBROUTINE RASTER(KREX,KREY,XMAX,XMIN,SX,YMAX,YMIN,SY,INDZ,X,Y)
C-----RASTER CONTROLS PLOTTING OF SCALES.
DIMENSION X(2),Y(2)
DATA NRAST/9999/
IF(INDZ.NE.0.AND.INDZ.NE.10) GOTO 50
IF(KREY.EQ.-1) GOTO 10
CALL RALIN(YMIN,YMAX,SY,XMIN,XMAX,M,X,Y)
GOTO 20
10 CALL RALOG(YMIN,YMAX,SY,XMIN,XMAX,M,X,Y)
20 CALL PLOTA(X,Y,M,2,0,1,1,1,INDZ,XMAX,XMIN,SX,YMAX,YMIN,SY,
11,NRAST,0,0,0)
IF(KREX.EQ.-1) GOTO 30
CALL RALIN(XMIN,XMAX,SX,YMIN,YMAX,M,Y,X)
GOTO 40
30 CALL RALOG(XMIN,XMAX,SX,YMIN,YMAX,M,Y,X)
40 CALL PLOTA(X,Y,M,2,0,1,1,1,INDZ,XMAX,XMIN,SX,YMAX,YMIN,SY,
11,NRAST,0,0,0)
50 RETURN
END

SUBROUTINE CALSKA(BSK,XSKA,YSKA,NPLOS,LASKA)
DIMENSION XSKA(1),YSKA(1)
REAL*8 BSK(1)
C CALSKA BEWIRKT,DASS NUR JEDE ZWEITE SKALENTEILUNG BESCHRIFTET WIRD.
M=10*NPLOS+LASKA
L=1
DO 10 I=1,M,2
BSK(I)=BSK(2*I-1)
BSK(I+1)=BSK(2*I)

```

XSKA(L)=XSKA(I)			
YSKA(L)=YSKA(I)			
10 L=L+1			
M=L-1			
NPLOS=M/10			
LASKA=M-10*NPLOS			
RETURN			
END			
SUBROUTINE CONLOG(X,XL,MAX,RAST,SLOG)			
C-----CONVERT DATA TO LOG, IF REQUIRED.			
DIMENSION X(2),XL(2)			
IF(RAST.NE.SLOG) GOTO 10			
DC 5 M=1,MAX			
5 XL(M)=ALOG10(X(M))			
GOTO 20			
10 CALL STRING(XL(1),X(1),4*MAX)			
20 RETURN			
END			
SUBROUTINE NEUDEN(KB,BRR,FROM,TO,X,NX,FODATA,*)			
C-----FETCH INTERPOLATION POINTS FROM NEUDAN.			
C-----SAME TECHNIQUE IS USED AS IN NEUDA-ROUTINE.			
C-----ERRSET IS USED TO HANDLE BAD TAPES.			
DIMENSION X(2),BAND(54),BRR(6)			
LOGICAL*4 FODATA			
COMMON/INOUT/KOUT			
CALL ERRSET(214,200,-10,1,1)			
NR=0			
NX=0			
10 IF(FODATA) READ(KB,100,END=80,ERR=90) (BAND(L),L=1,54)			
IF(.NOT.FODATA) READ(KB,120,END=80,ERR=90) (BAND(L),L=1,54)			
100 FORMAT(47A1,7E12.5,1X)			
120 FORMAT(47A1,7A4)			
CALL BDVGL(BAND,BRR,810)			
WRITE(KOUT,101) (BAND(I),I=1,31)			
101 FORMAT(/' GEFUNDEN:',31A1)			
C-----CONTROL SCALING OF PLOTS.			
15 XX=BAND(48)			
IF(XX.LT.FROM) GOTO 20			
IF(XX.GT.TO) GOTO 93			
GOTO 25			
20 IF(FODATA) READ(KB,100,END=81,FRR=91) (BAND(I),I=1,54)			
IF(.NOT.FODATA) READ(KB,120,END=81,ERR=91) (BAND(I),I=1,54)			
CALL BDVGL(BAND,BRR,494)			
GOTO 15			
25 NX=NX+1			
X(NX)=XX			
27 IF(FODATA) READ(KB,100,END=82,ERR=92) (BAND(I),I=1,54)			
	110	IF(.NOT.FODATA) READ(KB,120,FND=82,FRR=92) (BAND(I),I=1,54)	300
	120	CALL BDVGL(BAND,BRR,870)	310
	130	XX=BAND(48)	320
	140	IF(XX.GT.TO) GOTO 72	330
	150	IF(NX.FQ.1000) GOTO 74	340
	160	GOTO 25	350
	170	70 WRITE(KOUT,102) NX	360
	180	102 FORMAT(' ENDE DIESERS DATENSATZES.ES WURDEN ',I5,' DATENPUNKTE UEB	370
		1ERTRAGEN'/)	380
		RETURN	390
		72 WRITE(KOUT,103) NX	400
	103	FORMAT(' IM ANGFGERENEN ENERGIEBEREICH WURDEN ',I5,' DATENPUNKTE	410
		1GEFUNDEN'/)	420
		RETURN	430
		74 WRITE(KOUT,104) NX	440
	104	FORMAT(' ERLAUBTE MAXIMALZAHL VON DATEN =',I5,' UEBERTRAGEN.EINLE	450
		1SEN WIRD ABGEBROCHEN'/)	460
		RETURN	470
		80 WRITE(KOUT,105) BRR,KB	480
	105	FORMAT(/' FUER:',6A1,' WURDEN KEINE DATEN AUF BAND ',I2,' GEFUND	490
		1EN'/)	500
		RETURN 1	510
		81 WRITE(KOUT,106) FROM,TO,BRR,KB	520
	106	FORMAT(/' ZWISCHEN EMIN=',E14.6,' UND EMAX=',F14.6,' WURDEN FUER	530
		1',6A1,' KEINE DATEN AUF BAND ',I2,' GEFUNDEN'/)	540
		RETURN 1	550
		82 WRITE(KOUT,102) NX	560
		RETURN	570
		90 WRITE(KOUT,107)	580
	107	FORMAT(/' *** ERROR ON TAPE. RECORDS HAD TO BE SKIPPED.'/)	590
		READ(KB)	600
		GOTO 10	610
		91 WRITE(KOUT,107)	620
		READ(KB)	630
		GOTO 20	640
		92 WRITE(KOUT,107)	650
		WRITE(KOUT,108)	660
	108	FORMAT(' UNTER DEN AUSGELASSENEN RECORDS BEFINDEN SICH WAHRSCHE	670
	110	1NLICH AUCH RECORDS,DIE DEM GESUCHTEN DATENSATZ ANGEHOREN'/)	680
		GOTO 27	690
		93 WRITE(KOUT,106) FROM,TO,BRR,KB	700
		RETURN 1	710
		94 WRITE(KOUT,109) BRR,KB,FROM	720
	109	FORMAT(/' FUER ',6A1,' WURDEN AUF BAND ',I2,' KEINE DATENPUNKTE	730
		10BERHALB EMIN=',E14.6,'EV GEFUNDEN'/)	740
		RETURN 1	750
		END	760

```

SUBROUTINE LINEAR(IM,XP,YP,M,T,S,NR)
C-----LINEAR INTERPOLATION.
DIMENSION XP(IM),YP(IM),T(M),S(M)
COMMON/INOUT/KOUT
NR=0
IKU=1
IF(IM-1) 1701,1701,1702
1701 NR=1
GOTO 1703
1702 DO 1704 K=1,M
DO 1705 I=IKU,IM
IF(XP(I)-T(K)) 1705,1705,1706
1706 IK=I
GOTO 1707
1705 CCNTINUE
IF(T(K).NE.XP(IM)) GOTO 1710
IK=IM
GOTO 1707
1707 IF(IK-1) 1708,1708,1709
1708 NR=2
WRITE(KOUT,600) K,T(K),IK,XP(IK)
600 FORMAT(/5X,'DER ',I5,'. INTERPOLATIONSPUNKT ',E13.6,' LIEGT LINKS
X VOM ',I3,'. DATENPUNKT: ',E13.6/
18X,'AKTION: DIESER INTERPOLATIONSPUNKT WIRD UEBERGANGEN'//)
GOTO 1704
1709 D=(YP(IK)-YP(IK-1))/(XP(IK)-XP(IK-1))
S(K)=D*(T(K)-XP(IK))+YP(IK)
IKU=IK
1704 CONTINUE
1703 RETURN
1710 WRITE(KOUT,601) K
601 FORMAT(/5X,'AB DEM ',I5,'. INTERPOLATIONSPUNKT LIEGEN DIE INTERPO
LATIONSPUNKTE RECHTS VOM LETZTEN DATENPUNKT'/
18X,'AKTION: DIESE INTERPOLATIONSPUNKTE WERDEN LEBERGANGEN'//)
GOTO 1703
END

```

```

SUBROUTINE PREPLO
C-----INITIALIZE PLOT ACCORDING CONTROL INPUT FOR
C-----PLOT SIZE.
INTEGER PLEH,ZXINT,WNR,ZAHL,PLOZEI
COMMON/INOUT/KOUT,KIN
LCGICAL DRUCK,STANZ,PLO,PLOKU,DRUKU,STAKU,PLOSM,DRUSM,STASM,SMGEW,
1 NEUZEI,XINTPC,PLOFQU,FISS,SEQN,ADD,CALP,CALPJ,CATADD,TST,REWD
REAL NEIN,KURVE,MEDIUM,MEV,NULL,LTR,NEW
COMMON/CLP/CALP
COMMON/PARM1/ MAXOPT,PLEH,PLCZEI,ZXINT,IDPLOT,CALPJ
COMMON/PARM2/ IPLOGR,ISTAFO,ISMO,IFMOUT,ISMNP,DISP,EINHXTX,EINHXY,
1 EINHDT,XEINHDT,YEINHDT,DEINHDT,LTR,KARTEH,FMTOUT(20),STAFMT(20),
2 FMT(20),DRUCK,STANZ,PLO,KURVE,PLOKU,DRUKU,STAKL,PLOSM,DRUSM,
3 STASM,SMGEW,NEUZEI,XINTPO,PLOFQU,FISS,SEQN,ADD,CATADD,TST,
4 REWD,NANZ,NADD,ZAHL,MEDIUM,WNR,

```

```

10 5 RAST(2),ISMAX,FQU,NAMZ 160
20 COMMON/PAR/XMAX,SX,YMAX,SY,XMIN,YMIN,XMAL,XMIL,YMAL,YMIL,NLGGX, 170
30 1 ALGGY,NX,NY,INDZ,JNDZ,XMILJ,YMILJ,JP 180
40 DATA XLOG/'XLOG'/,YLOG/'YLOG'/,NEIN/'NEIN'/,XSTA/-1.E-60/ 190
50 DATA XMINO,XMAXO,YMINO,YMAXO/4*0./ 200
60 NAMELIST/PLOTGR/XMIN,XMAX,XLG,YMIN,YMAX,YLG,RAST,EINHXTX,EINHXY, 210
70 ICALP 220
80 COMMON/PARM4/FROM,TO 230
90 XMIN=XSTA 240
100 C-----TEST WHETHER DATA FOR PLOT LIMITS WERE ACTUALLY IN INPUT. 250
110 C-----IF YES, PERFORM CONVERSION TO INTERNAL PHYSICAL UNITS. 260
120 XMAX=XSTA 270
130 YMIN=XSTA 280
140 YMAX=XSTA 290
150 READ(KIN,PLOTGR) 300
160 IF(XMIN.EQ.XSTA) GOTO 1410 310
170 CALL ECONVL(XMIN,EINHXTX) 320
180 XMINO=XMIN 330
190 GOTO 1415 340
200 1410 XMIN=XMINO 350
210 1415 IF(XMAX.EQ.XSTA) GOTO 1420 360
220 CALL ECONVL(XMAX,EINHXTX) 370
230 XMAXO=XMAX 380
240 GOTO 1425 390
250 1420 XMAX=XMAXO 400
260 1425 IF(YMIN.EQ.XSTA) GOTO 1430 410
270 CALL ECONVL(YMIN,EINHXY) 420
280 YMINO=YMIN 430
290 GOTO 1435 440
300 1430 YMIN=YMINO 450
310 1435 IF(YMAX.EQ.XSTA) GOTO 1440 460
320 CALL ECONVL(YMAX,EINHXY) 470
330 YMAXO=YMAX 480
340 GOTO 1450 490
350 1440 YMAX=YMAXO 500
360 C-----INCREMENT PLOTID AND SET PLOT SIGN SELECTICN FLAG. 510
1450 ISMNP=0 520
ICFLOT=IDPLOT+1 530
PLOZEI=-1 540
JP=0 550
IF(RAST(1).EQ.NEIN) RAST(2)=NEIN 56C
C-----CONFIRM INPUT. 570
WRITE(KOUT,636) 580
WRITE(KOUT,615) XMIN,XMAX,XLG,YMIN,YMAX,YLG,RAST,EINHXTX,EINHXY, 590
*CALP 600
WRITE(KOUT,636) 610
IPLOGR=1 620
FROM=XMIN 630
TO=XMAX 64C
SX=(XMAX-XMIN)/(40.*XLG) 650
SY=(YMAX-YMIN)/(40.*YLG) 660
XMAL=XMAX 670
XMIL=XMIN 680
YMAL=YMAX 690
YMIL=YMIN 700

```



```

NLGGX=+1
NLGGY=+1
NX=+1
NY=+1
C-----LOG. SCALE.
IF(RAST(1).NE.XLOG) GOTO 146
X*AL=ALOG10(X*MAX)
XMIL=ALOG10(XMIN)
SX=(X*AL-XMIL)/(40.*XLG)
NLGGX=-1
IF(XMIL.LT.0.) NY=0
146 IF(RAST(2).NE.YLOG) GOTO 147
Y*AL=ALOG10(Y*MAX)
YMIL=ALOG10(YMIN)
SY=(Y*AL-YMIL)/(40.*YLG)
NLGGY=-1
IF(YMIL.LT.0.) NX=0
C-----ESTABLISH PLOT SIZE INDICATOR FOR PLOTA.
147 F=XLG/25.
IF(CALP) F=XLG/60.
IF(F.LT.4.) INDZ=INT(F)+1
IF(F.EQ.4.) INDZ=4
IF(F.GT.4.) GOTO 806
IF(.NOT.CALP) GOTO 148
INDZ=INDZ+10
CALPJ=.TRUE.
XMILJ=XMIL
YMILJ=YMIL
148 JNDZ=INDZ
RETURN
806 WRITE(KOUT,609) F
INDZ=4
IF(CALP) INDZ=INDZ+10
JNDZ=INDZ
RETURN
609 FORMAT(/ /5X,'FEHLERNACHRICHT'/8X,'F=XLG/YLG>4 NICHT ERLAUBT:F =',E
112.5/5X,'AKTIGN: PLOTFORMAT - F=4 GESETZT'//)
615 FORMAT(' PLOTGR XMIN=',E12.3,' , XMAX=',E12.3,' , XLG=',E12.3/9X,
1'YMIN=',E12.3,' , YMAX=',E12.3,' , YLG=',E12.3/8X,' RAST=',A4,' ,',
2A4/9X,' EINHTX=',A4,' , EINHTY=',A4,' , CALP=',L2)
636 FCRRAT(2X,130(' - '))
END

SUBROUTINE CDINPT(A,TEND)
C-----COPY CONTROL INPUT TO FT09 TO ENABLE BACKSFACING.
DIMENSION A(20),TEND(20)
COMMON/INOUT/KOUT,KIN,DUMMY,KINC
DATA TTEND/'ENDE'/,BLANK/' '/
TEND(1)=TTEND
DO 100 I=2,20
100 TEND(I)=BLANK
WRITE(KOUT,600)

```

```

710 600 FCRRAT('1',10X,'NACHRICHT DER SUBROUTINE *CDINPT*'/11X,'CDINPT LEG 100
720 IT DIE KARTENEINGABE AUF G.FTC9'/ 110
730 211X,'DIE MAXIMALZAHL VON EINGABEKARTEN IST DURCH DEN SPACEPARAMETE 120
740 3R FUER FT09 GEGEBEN') 130
750 N=1 140
760 1000 READ(KINC,500,END=99,ERR=9) A 150
770 WRITE(KIN,500) A 160
780 500 FCRRAT(20A4) 170
790 N=N+1 180
800 GOTO 1000 190
810 9 WRITE(KOUT,601) N 200
820 601 FORMAT('/ FEHLER BEIM LESEN DER',I5,'-TEN KARTE'/) 210
830 GOTO 1000 220
840 95 WRITE(KIN,500) TEND 230
850 REWIND KIN 240
860 WRITE(KOUT,602) N 250
870 602 FCRRAT('/ DER DATASET AUF G.FT09 BESTEHT AUS',I5,' RECORDS,'/5X,' 260
880 IALS LETZTES RECORD IST EIN *ENDE*-RECORD HINZUGEUEGT WORDEN') 270
890 RETURN 280
900 END 290
910 920 930 940 950 960 970 980 990 1000 1010 1020 1030 1040 1050 1060 1070 1080 1090 1100 1110 1120
SUBROUTINE ORDNRN(JMAX,X,Y,Z,U,V,W) 10
C-----SORT DATA. ROUTINE IS OPTIMIZED ASSUMING THAT DATA 20
C-----WILL BE SORTED USUALLY. 30
DIMENSION X(JMAX),Y(JMAX),Z(JMAX),U(JMAX),V(JMAX),W(JMAX) 40
U(1)=X(1) 50
V(1)=Y(1) 60
W(1)=Z(1) 70
IF(JMAX.LT.2) GOTO 20 80
DO 10 J=2,JMAX 90
M=J-1 100
1 IF(X(J).LT.U(M)) GOTO 2 110
IF(J.NE.(M+1)) GOTO 3 120
U(J)=X(J) 130
V(J)=Y(J) 140
W(J)=Z(J) 150
GOTO 10 160
2 M=M-1 170
IF(M.NE.0) GOTO 1 180
3 IA=M+1 190
IF=J-1 200
CC 4 K=IA,IE 210
I=IA+IE-K 220
U(I+1)=U(I) 230
V(I+1)=V(I) 240
4 W(I+1)=W(I) 250
U(IA)=X(J) 260
V(IA)=Y(J) 270
W(IA)=Z(J) 280
10 CCNTINUE 290
20 RETURN 300
END 310

```

```

SUBROUTINE KEDEN(NARG,NAME,EUNT,EOB,XINT,ZXINT,*)
C-----RETRIEVE INTERPOLATION ENERGIES FROM KEDAK.
C-----SAME TECHNIQUE IS USED AS IN ROUTINE KEDSUB.
  REAL*8 NAME(1),NAMEN(3)
  COMMON/INOUT/KOUT
  INTEGER ZXINT
  DIMENSION XINT(2),NAMZ(3),VAL(2)
  DO 10 I=1,NARG
10  NAMEN(I)=NAME(I)
  NR=0
  ZXINT=0
  NAMZ(1)=NARG
  CALL LDFLOC(NERR,NAMZ,NAMEN,VAL)
  IF(NERR.EQ.0) GOTO 100
  IF(VAL(1).GT.EOB) GOTO 105
  IF(VAL(1).LT.EUNT) GOTO 15
  GOTO 20
15  CALL LDFNXT(NERR,NAMZ,NAMEN,VAL)
  IF(NERR.EQ.0) GOTO 110
  IF(VAL(1).LT.EUNT) GOTO 15
  IF(VAL(1).GT.EOB) GOTO 115
  GOTO 20
25  CALL LDFNXT(NERR,NAMZ,NAMEN,VAL)
  IF(NERR.EQ.0) GOTO 50
  IF(VAL(1).GT.EOB) GOTO 60
  IF(ZXINT.EQ.1000) GOTO 120
20  ZXINT=ZXINT+1
  XINT(ZXINT)=VAL(1)
  GOTO 25
50  WRITE(KOUT,601) (NAME(I),I=1,2)
  IF(NARG.GE.3) WRITE(KOUT,608) NAME(3)
  WRITE(KOUT,609) XINT(ZXINT),ZXINT
601  FORMAT(/5X,'MESSAGE FROM DATARETRIEVAL FROM KEDAK FOR:/'
1 12X,A8,2X,A8)
608  FORMAT('+',32X,E14.6)
609  FORMAT(8X,'LAST KEDAKVALUE ENCOUNTERED BEFORE SPECIFIED EOB WAS RE
  LACHED'/
2 8X,'LAST ENERGY TRANSMITTED:',E14.6/
3 8X,'NUMBER OF ENERGIES TRANSMITTED:',I6//)
  RETURN
60  WRITE(KOUT,601) (NAME(I),I=1,2)
  IF(NARG.GE.3) WRITE(KOUT,608) NAME(3)
  WRITE(KOUT,602) XINT(ZXINT),ZXINT
602  FORMAT(8X,'LAST ENERGY IN SPECIFIED INTERVAL FOUND'/
1 8X,'ITS VALUE IS ',E14.6/
2 8X,'AND THE NUMBER OF ENERGIES TRANSMITTED IS ',I6//)
  RETURN
100  WRITE(KOUT,601) (NAME(I),I=1,2)
  IF(NARG.GE.3) WRITE(KOUT,608) NAME(3)
  WRITE(KOUT,603)
603  FORMAT(8X,'NO ENTRY FOR THIS DATA TYPE ON KEDAK'//)

```

```

  RETURN 1
105  WRITE(KOUT,601) (NAME(I),I=1,2)
  IF(NARG.GE.3) WRITE(KOUT,608) NAME(3)
  WRITE(KOUT,604)
604  FORMAT(8X,'FIRST ENERGY OF THIS DATA TYPE IS > SPECIFIED EUNT'//)
  RETURN 1
110  WRITE(KOUT,601) (NAME(I),I=1,2)
  IF(NARG.GE.3) WRITE(KOUT,608) NAME(3)
  WRITE(KOUT,605)
605  FORMAT(8X,'ALL ENERGIES OF THIS DATA TYPE ARE < SPECIFIED EUNT'//)
  RETURN 1
115  WRITE(KOUT,601) (NAME(I),I=1,2)
  IF(NARG.GE.3) WRITE(KOUT,608) NAME(3)
  WRITE(KOUT,606)
606  FORMAT(8X,'NO DATAPPOINTS IN SPECIFIED ENERGY INTERVAL'//)
  RETURN 1
120  WRITE(KOUT,607) ZXINT
607  FORMAT(//' ERLAUBTE MAXIMALZAHL VON DATEN=',I5,' WURDE UEBERTRAGE
  IN.EINLESEN WIRD ABGEBROCHEN.'//)
  RETURN
  END

```

```

SUBROUTINE RALOG(YMIN,YMAX,SY,XMIN,XMAX,NPOINT,X,Y)
C-----PROVIDE LOG. SCALE LINES.
  DIMENSION X(2),Y(2),DYL(9)
  COMMON/INOUT/KOUT
  DATA DYL/.0458,.3010,.1761,.1250,.0969,.0792,.0669,.0580,.0511/
  YL=YMAX-YMIN
  NDEK=INT(YL)
  IF((YL-FLOAT(NDEK)).GT.C.8) NDEK=NDEK+1
  IF(NDEK.GT.11) GOTO 25
  IF(SY.GT..005) GOTO 25
  NLIN=9*NDEK+1
  NPOINT=2*NLIN
  J=0
  Y1=YMIN
  Y2=YMIN+DYL(2)
  DO 20 I=1,NPOINT,4
  X(I)=XMIN
  X(I+1)=XMAX
  X(I+2)=XMAX
  X(I+3)=XMIN
  Y(I)=Y1
  Y(I+1)=Y1
  Y(I+2)=Y2
  Y(I+3)=Y2
  J=J+2
  Y1=Y2+DYL(MOD(J,9)+1)
  Y2=Y1+DYL(MOD(J+1,9)+1)
20  CONTINUE
  Y(NPOINT-1)=YMAX
  Y(NPOINT)=YMAX

```

```

GCTO 30
25 NPOINT=0
WRITE(KOUT,601)
601 FCRMAT(' ***** ERRORMESSAGE FROM RASTER - ROUTINE. . .MINIMUM LENG
1TH OF A DECADE IS 5 CM. '/
220X,'DECREASE NUMBER OF DECADES OR USE CALCOMP - PLOTTER. ')
30 RETURN
END

```

```

SUBROUTINE RALIN(YMIN,YMAX,SY,XMIN,XMAX,NPOINT,X,Y)
C-----PROVIDE LIN SCALE LINES.
DIMENSION X(2),Y(2)
DY=80.*SY
YL=YMAX-YMIN
NLIN=INT(YL/DY)
IF((YL-FLOAT(NLIN)*DY).GT.0.99*DY) NLIN=NLIN+1
NLIN=NLIN+1
NPOINT=2*NLIN
IF(NPOINT.GT.200) NPOINT=200
Y1=YMIN
Y2=YMIN+DY
DC 10 I=1,NPOINT,4
X(I)=XMIN
X(I+1)=XMAX
X(I+2)=XMAX
X(I+3)=XMIN
Y(I)=Y1
Y(I+1)=Y1
Y(I+2)=Y2
Y(I+3)=Y2
Y1=Y2+DY
Y2=Y1+DY
10 CONTINUE
Y(NPOINT)=YMAX
Y(NPOINT-1)=YMAX
RETURN
END

```

```

SUBROUTINE ECCNV(M,X,E)
C-----CONVERT DATA TO INTERNAL PHYSICAL UNITS (EV USUALLY.)
DIMENSION X(M),D(M)
REAL MEV,KEV,MIBA
KCNTN=0
GCTO 10
ENTRY ECONV1(Y,E)
KCNTN=1
10 K=C
DATA MEV/'MEV ',KEV/'KEV ',MIBA/'MIBA ',PERC/'PERC' /
IF(E.EQ.PERC) GOTO 10

```

```

310 IF(E.EQ.MEV) K=6
320 IF(E.EQ.KEV) K=3
330 IF(E.EQ.MIBA) K=-3
340 IF(KONTR.EQ.1) GOTO 2
350 DO 1 I=1,M
360 1 X(I)=X(I)*(10.**K)
370 RETURN
380 2 Y=Y*(10.**K)
RETURN
ENTRY PERCNT(M,X,D)
DO 3 I=1,M
3 D(I)=D(I)*X(I)*.01
RETURN
END

```

```

SUBROUTINE ERRDET(ET,IRTP)
C-----DETERMINE PHYSICAL UNITS OF ERROR BARS
C-----READ FROM NEUDADA TAPE ACCORDING NEUDADA CONVENTIONS.
INTEGER*2 ETS,IRA,IRP,IRS,ERA(6),ERP(6),ERS(2)
COMMON/INOUT/KOUT
INTEGER*2 ITS
DIMENSION ET(2)
DATA IRA/6/,IRP/6/,IRS/2/
DATA ERA/'LE','LR','LL','L<','L>','LS' /
DATA ERP/'PE','PR','PP','P<','P>','PS' /
DATA ERS/'MR','NR' /
DATA ATEXT /'ABSO','LUTE','ERR','OR-U','NITS','ABSO','LUTE',
'ERR','OR-U','NITS','ABSO','LUTE','ERR','OR-U',
2'NITS','AVER','AGE','ABS.','ERR','OR<','AVER','AGE','ABS.',
3'ERR','OR>','ABS.','STAT','TIST','ERR','OR' /
DATA PTEXT /'ERRO','R IN','PER','CENT','ERRO','R IN',
1'PER','CENT','ERRO','R IN','PER','CENT',
2'','AVER','PE','RC','ERRO','R<','AVER','PE','RC',
3'ERRO','R>','STAT','IST','ERR','OR I','N %' /
DATA STEXT/'ERRO','R IN','MS','METE','R','ERRO','R IN',
1'NS','METE','R' /
LOGICAL*1 LETS(2)
EQUIVALENCE(LETS(1),ETS)
CALL STRING(LETS(1),ET(1),1)
CALL STRING(LETS(2),ET(2),1)
GCTO 5
ENTRY ERRT(ITS,IRTP)
ETS=ITS
5 DC 10 J=1,IRA
IF(ETS.NE.ERA(J)) GOTO 10
IRTP=1
WRITE(KOUT,601) ETS,(ATEXT(I,J),I=1,5)
601 FCRMAT(8X,' ERRORTYPE ',A2,' MEANS ',5A4//)
GOTO 90
10 CONTINUE
DC 20 J=1,IRP

```

```

10
20
30
40
50
60
70
80
90
100
110

```

```

IF(ETS.NE.ERP(J)) GOTO 20
IRTP=2
WRITE(KOUT,601) ETS,(PTEXT(I,J),I=1,5)
GOTO 90
20 CCNTINUE
DO 30 J=1,IRS
IF(ETS.NE.ERS(J)) GOTO 30
IRTP=2+J
WRITE(KOUT,601) ETS,(STEXT(I,J),I=1,5)
GOTO 90
30 CONTINUE
WRITE(KOUT,600) ETS
600 FORMAT(' FEHLERTYP=',A2,' NICHT ERKANNT'//)
IRTP=0
GOTO 90
90 RETURN
END

```

```

380
390
400
410
420
430
440
450
460
470
480
490
500
510
520
530
540

```

```

11N THE RANGE,WHERE THE CURVE IS DEFINED.*/
2 10X,'ACTION TAKEN : CURVE IS NOT INTERPOLATED.'//)
130 RETURN
END

```

```

350
360
370
380

```

```

SUBROUTINE PCHECK(N1,X1,N2,X2,NA,NB)
DIMENSION X1(1),X2(1)
COMMON/INOUT/KOUT
C
C CHECK WHETHER ALL INTERPOLATION POINTS LY WITHIN THE BOUNDARY
C VALUES X1(1),X1(N1)
DO 10 I=1,N2
IF(X2(I).LT.X1(1)) GOTO 10
NA=I
GOTO 20
10 CCNTINUE
NA=NA+1
NB=NA
GOTO 100
20 J=N2
DO 30 I=NA,N2
IF(X2(J).GT.X1(N1)) GOTO 30
NB=J
GOTO 100
30 J=J-1
NB=NB-1
100 IF(NA.EQ.1) GOTO 110
WRITE(KOUT,601) X2(NA),X1(1)
601 FORMAT(/5X,'MESSAGE:ALL INTERPOLATION POINTS <=',1PE13.5/
1 10X,'ARE BELOW FIRST POINT FOR WHICH THE CURVE IS DEFINED =' ,1PE13
23.5/10X,'ACTION TAKEN: THESE POINTS ARE SKIPPED.'//)
110 IF(NB.EQ.N2) GOTO 120
WRITE(KOUT,602) X2(NB),X1(N1)
602 FORMAT(/5X,'MESSAGE: ALL POINTS >= ',1PE13.5/
1 10X,'ARE ABOVE LAST POINT FOR WHICH THE CURVE IS DEFINED =' ,1PE13
2.5/10X,'ACTION TAKEN:THESE POINTS ARE SKIPPED.'//)
120 IF(NB.GE.NA) GOTO 130
WRITE(KOUT,603)
603 FORMAT(/5X,'ERRORMESSAGE:NCNE OF THE INTERPOLATION POINTS IS WITH

```

```

10
20
30
40
50
60
70
80
90
100
110
120
130
140
150
160
170
180
190
200
210
220
230
240
250
260
270
280
290
300
310
320
330
340

```

#### 4.2 Karlsruhe Version of SCORE

The SCORE program developed by C.L. Dunford for interactive graphic cross section evaluation was adapted to the file formats used at Karlsruhe (KEDAK for evaluated data, NEUDADA for experimental data) by I. Langner and R. Meyer. The following user guide is essentially a translation of unpublished internal reports written by I. Langner.

Adaptation of SCORE to KEDAK and NEUDADA Data  
SCORE Version III-K-172

I. Langner

## Contents

### Introduction

#### 4.2.1 Information for SCORE users

- 4.2.1.1 General Introduction to SCORE
- 4.2.1.2 Job control cards for SCORE
- 4.2.1.3 Input for SCORE at the IBM-2250 terminal
- 4.2.1.4 Job control cards and input for NAP

#### 4.2.2 Modifications in SCORE

- 4.2.2.1 Changes in the overlay structure
- 4.2.2.2 SCORE routines
  - 4.2.2.2.1 MAIN PROGRAM SCORE
  - 4.2.2.2.2 Subroutines ERRCOR, SPDATA, CHOSZA
- 4.2.2.3 Program NAP - Compilation of SCORE library from experimental (NEUDADA) data

#### 4.2.3 New subroutines for SCORE

- 4.2.3.1 General subroutines
  - 4.2.3.1.1 SELEC - Choice between NEUDADA and KEDAK
  - 4.2.3.1.2 IDENT - User and case identification
- 4.2.3.2 Subroutines for the organization of KEDAK data
  - 4.2.3.2.1 KNDF - KEDAK data for curve display
  - 4.2.3.2.2 TAPKND - Calculation of numeric KEDAK names
  - 4.2.3.2.3 RKNDF - Reading of KEDAK data
  - 4.2.3.2.4 SKIPØI - Deletion of data points
  - 4.2.3.2.5 SELK - Calculation of the differential elastic-scattering cross section
  - 4.2.3.2.6 KEDAK - KEDAK data without experimental (NEUDADA) data
  - 4.2.3.2.7 ANERG - Listing of energies on the screen

### References

### Appendices

- A1 List of changes in SCORE routines
- A2 Source program lists of the new subroutines
- A3 Input summary

## Introduction

SCØRE - a program system for cross section evaluation - was developed by Atomics International and IBM Palo Alto for interactive graphic data evaluation at special terminal stations, the IBM 2250 terminal /9/.

The SCØRE system /1/, /12/ has recently been adapted to experimental data in NEUDADA format /7/ and to evaluated data in KEDAK format /3/, /4/.

Originally SCØRE was designed for processing of experimental data in SCISRS format (SCISRS: Sigma Center Information Storage and Retrieval System) and evaluated data in ENDF/B format /2/.

The Karlsruhe Version III-K-1972 is based on SCØRE Version III-1970. The changes and additions permit reading of KEDAK data from the KNDF library (Karlsruhe Nuclear Data File) and their display on screen. Curves of KEDAK data can be displayed either alone or together with experimental data. The SCØRE Version III-K-1972 was initiated by Dr. R. Meyer who also participated in its development.

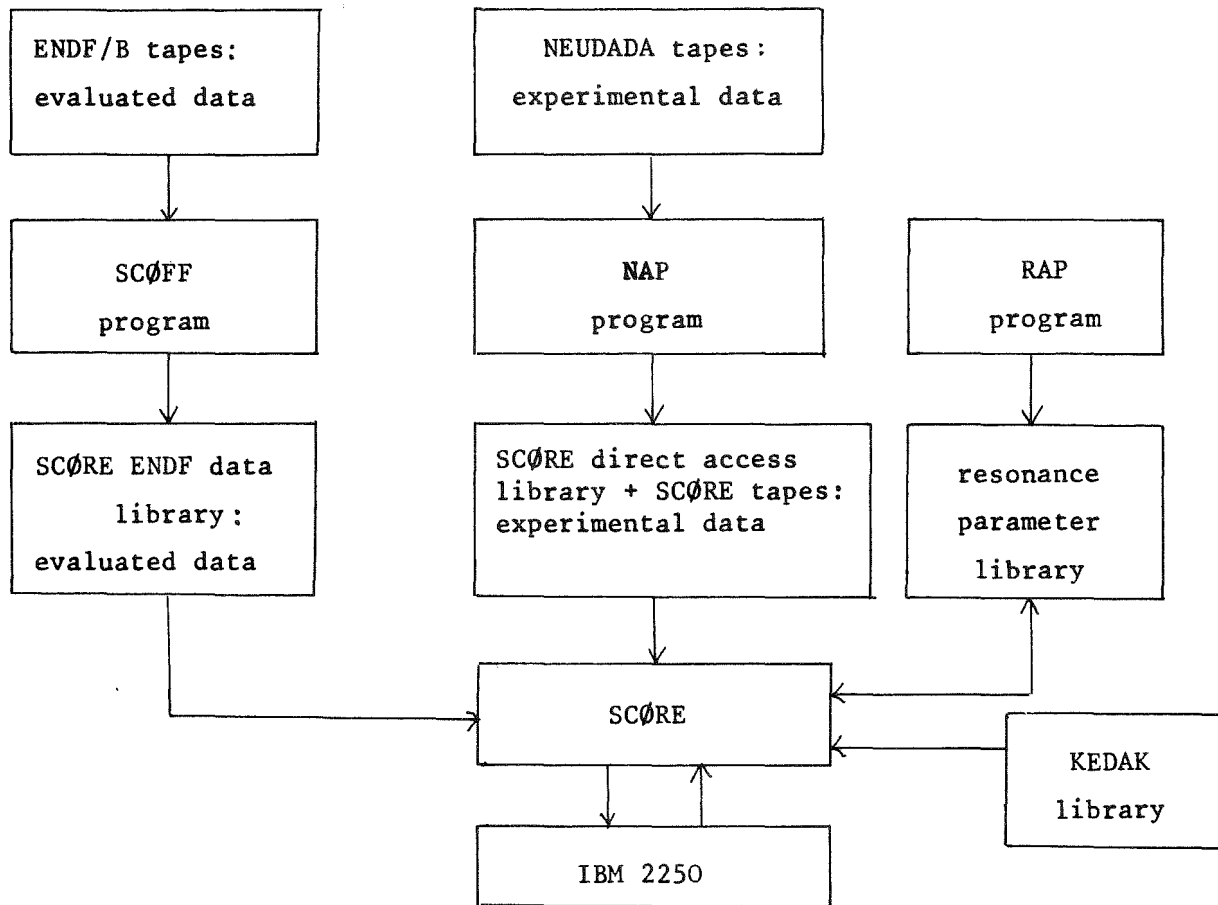
A very informative general description of SCØRE with reproductions of the various Figures displayed on the screen can be found in Ref. /12/.

### 4.2.1 Information for SCØRE users

#### 4.2.1.1 General Introduction to SCØRE

SCØRE consists of the SCØRE program proper and three administration programs for the associated data libraries: SCØFF, RAP and NAP.





The structure of the data files, as they are distributed on tape by the neutron data centers, does not allow rapid processing. Therefore the data are reorganized in special formats that were developed for SCØRE:

1. The SCØFF program treats ENDF/B tapes.
2. The RAP program establishes the resonance parameter library.
3. The NAP program (NEUDADA Adaptation Program) writes NEUDADA files /7/ of experimental data in a SCØRE format and produces, apart from SCØRE tapes, a direct-access library on disc (see 2.3).

SCØRE can process data from tape, disc (direct access) and from the IBM-2250 terminal - via keyboard, light pen and programmed-function keys.

4.2.1.2 The job control cards for SCORE

The following listing shows all possible job control cards for SCORE:

```
//INR048SC JOB (0048,101,P6M1A),LANGNER,CLASS=G,REGION=160K,TIME=3
/*SETUP DEVICE=2250,ID=RESERV
/*SETUP DEVICE=TAPE9,ID=(SCORE1,NORING,,NL)
/*SETUP DEVICE=TAPE9,ID=(ENDFB1,NORING,,NL)
/*SETUP DEVICE=2314,ID=GFKC16
//EXEC FTG,LIB=NUSYS,NAME=SCORE,REGION.G=160K,TIME.G=5
//G.SCOPE DD UNIT=2250
//TAPE01 DD UNIT=2400,VOL=SER=SCORE1,DISP=CLD,LABEL=(1,NL)
//TAPE02 DD UNIT=2400,VOL=SER=ENDFB1,DISP=OLD,LABEL=(1,NL)
//DISK01 DD UNIT=2314,VOL=SER=GFK016,DISP=SHR
//G.FT09F001 DD VOL=REF=*.DISK01,DISP=(NEW,DELETE,DELETE),
//   DCB=(RECFM=V,LRECL=3524,BLKSIZE=3528,BUFNO=1),
//   SPACE=(3524,(5,1)),DSNAME=RESTEMP
//G.FT16F001 DD VOL=REF=*.DISK01,DISP=OLD,DSNAME=ENDFB,
//   DCB=(RECFM=F,BLKSIZE=3600,LRECL=3600,BUFNO=1)
//G.FT17F001 DD VOL=REF=*.DISK01,DISP=OLD,DSNAME=NEUDAD,
//   DCB=(RECFM=F,BLKSIZE=3520,LRECL=3520,BUFNO=1)
//G.FT19F001 DD VOL=REF=*.DISK01,DISP=OLD,DSNAME=RESLIB
//   DCB=(RECFM=U,BLKSIZE=400,BUFNO=1)
//G.FT20F001 DD VOL=REF=*.TAPE01,DISP=OLD,LABEL=(1,NL,,IN),
//   DSN=NAP.LABEL,DCB=(RECFM=VBS,LRECL=68,BLKSIZE=3744)
//G.FT21F001 DD VOL=REF=*.TAPE01,DISP=OLD,LABEL=(2,NL,,IN),
//   DSN=NA023,DCB=*.FT20F001
//G.FT22F001 DD VOL=REF=*.TAPE01,DISP=OLD,LABEL=(3,NL,,IN),
//   DSN=FE000,DCB=*.FT20F001
//G.FT23F001 DD VOL=REF=*.TAPE01,DISP=OLD,LABEL=(4,NL,,IN),
//   DSN=FE054,DCB=*.FT20F001
//G.FT24F001 DD VOL=REF=*.TAPE01,DISP=CLD,LABEL=(5,NL,,IN),
//   DSN=LI000,DCB=*.FT20F001
//G.FT25F001 DD VOL=REF=*.TAPE01,DISP=CLD,LABEL=(6,NL,,IN),
//   DSN=LI007,DCB=*.FT20F001
//G.FT26F001 DD VOL=REF=*.TAPE01,DISP=OLD,LABEL=(7,NL,,IN),
//   DSN=PU241,DCB=*.FT20F001
//G.FT27F001 DD VOL=REF=*.TAPE02,DISP=CLD,LABEL=(1,NL,,IN),
//   DSN=SCOFF.LABEL,DCB=(RECFM=V,LRECL=360,BLKSIZE=3608,BUFNO=1)
//G.FT28F001 DD VOL=REF=*.TAPE02,DISP=CLD,LABEL=(2,NL,,IN),
//   DSN=LI007B,DCB=*.FT27F001
//G.FT29F001 DD VOL=REF=*.TAPE02,DISP=OLD,LABEL=(3,NL,,IN),
//   DSN=NA023B,DCB=*.FT27F001
//G.FT30F001 DD VOL=REF=*.TAPE02,DISP=OLD,LABEL=(4,NL,,IN),
//   DSN=FE000B,DCB=*.FT27F001
//G.FT31F001 DD VOL=REF=*.TAPE02,DISP=OLD,LABEL=(5,NL,,IN),
//   DSN=PU241B,DCB=*.FT27F001
//G.SYSIN DD DUMMY
```

The following example of job control cards requires only KEDAK and NEUDADA data.

```
//INR048SC JOB (0048,101,P6M1A),LANGNER,CLASS=G,REGION=160K,TIME=3
// EXEC FHG,LIB=NUSYS,NAME=SCORE,REGION.G=160K,TIME.G=5
//G.SCOPE DD UNIT=250
//TAPE01 DD UNIT=2400,VOL=SER=SCORE1,DISP=OLD,LABEL=(1,NL)
//G.FTC1F001 DD UNIT=2314,VOL=SER=NUSYS0,DISP=SHR,DSN=KNDF
//G.FT17F001 DD UNIT=2314,VOL=SER=GFK016,DISP=SHR,DSN=NEUDADA.DA.LAN,
//   DCB=(RECFM=F,LRECL=3520,BLKSIZE=3520,BUFNO=1)
//G.FT20F001 DD VOL=REF=*.TAPE01,DISP=OLD,LABEL=(1,NL,,IN),
//   DCB=(RECFM=VBS,LRECL=68,BLKSIZE=3744,BUFNO=1),DSN=NAP.LABEL
//G.FT21F001 DD VOL=REF=*.TAPE01,DISP=OLD,LABEL=(2,NL,,IN),
//   DSN=NA023,DCB=*.FT20F001
//G.FT22F001 DD VOL=REF=*.TAPE01,DISP=OLD,LABEL=(3,NL,,IN),
//   DSN=FE000,DCB=*.FT20F001
//G.FT23F001 DD VOL=REF=*.TAPE01,DISP=OLD,LABEL=(4,NL,,IN),
//   DSN=FE054,DCB=*.FT20F001
//G.FT24F001 DD VOL=REF=*.TAPE01,DISP=OLD,LABEL=(5,NL,,IN),
//   DSN=LI000,DCB=*.FT20F001
//G.FT25F001 DD VOL=REF=*.TAPE01,DISP=OLD,LABEL=(6,NL,,IN),
//   DSN=LI007,DCB=*.FT20F001
//G.FT26F001 DD VOL=REF=*.TAPE01,DISP=OLD,LABEL=(7,NL,,IN),
//   DSN=PU241,DCB=*.FT20F001
//G.SYSIN DD DLMYY
```

As shown by the EXEC card SCORE is part of the library LOAD.NUSYS.

#### 4.2.1.3 Input for SCORE at the IBM 2250 terminal

A summary of the SCORE input is appended under A3. below.

##### 4.2.1.2.1 Input modes

Three different input modes exist:

1. light pen display on the screen,
2. input via keyboard as alphameric text, fixed- or floating-point numbers,
3. input via the programmed-function keyboard controlling program execution.

Note: Key no. 1 of the programmed-function keyboard has the same function as the END key of the keyboard.

#### 4.2.1.3.2 Input for SCØRE, mainly for KEDAK data

Reproductions of many of the Figures displayed on the screen can be found in Ref. /12/.

First figure: The user can choose between two options: KEDAK and NEUDADA (cf. 4.2.3.1). The option selected by light pen is displayed in large size /9/ on the screen. The information is transferred to the main program when the END key is pressed.

On the second figure SCØRE requests user and case identification (cf. /1/, Vol. I, p. 16).

#### KEDAK option

If the KEDAK option was selected on the first figure a list of the alphameric names /5/ of the isotopes available in the KNDF library /3/, /4/ is displayed in the third figure. Fifteen names are shown per figure. Selecting the option PAGE with the light pen the user can "turn the page", i.e. get the next 15 names displayed. He chooses the isotope with the light pen. The chosen isotope is shown in large size. If the KNDF library does not contain any reaction types for the chosen isotope that can be processed by SCØRE, pressing of the END key causes display of the first 15 isotopes again. Otherwise a list of reaction types available for SCØRE processing follows in the next figure. Again the choice is made by light pen. The chosen reaction type is displayed in large size in the list, and also a request for the energy limits. The energy limits must be typed in. Pressing of the END key causes display of the cross section curve in "Basic SCØRE Data Display" (/1/, Vol. I, p. 16). Exceptions are the reaction types SGNC (differential elastic-scattering cross section) and SGIZ (partial inelastic-scattering cross section) for which a list of incident (SGNC) or excitation (SGIZ) energies is shown first, with PAGE having the same function as described above. After selecting the appropriate energy with the light pen the user can get the cross section curve displayed by pressing the END key.

The following selection of the original options (cf. /1/, Vol. I, p. 27) are valid, i.e. can be chosen with the light pen:

MAIN 1/3                    option list 1 -  
RESTART  
PRINT  
EXPAND  
QUIT  
MAIN 2/3                    option list 2 -  
KEDAK

Meaning:

MAIN 1/3                    display option list MAIN 2/3,  
  
RESTART                    jump back to the start of the program,  
                          i.e. display first picture again,  
  
PRINT                      print the data displayed on the screen,  
  
EXPAND                     Expand the graph portion of the display to  
                          full screen size with omission of comment and  
                          option lists, (EXPAND can be turned off again  
                          by pressing the END key),  
  
QUIT                        end terminal session,  
  
MAIN 2/3                    display option list MAIN 3/3 ,  
  
REPLOT                     change co-ordinate system according to input  
                          furnished by user ,  
  
MAIN 3/3                    display option list MAIN 1/3 ,  
  
KEDAK                      display KEDAK (evaluated) data in curve  
                          form together with experimental data points.

NEUDADA option

If the NEUDADA option is chosen in the first figure the description in Ref. /1/  
is valid for subsequent input. Selecting KEDAK from option list 3/3 one obtains  
the (evaluated) KEDAK data in curve form together with the (experimental)  
NEUDADA data /7/.

#### 4.2.1.3.3 Summary of light-pen input

- Figure 1 (1) select KEDAK or NEUDADA option by light pen,  
(2) press END key.
- Figure 2 user identification etc., cf. Ref. /1/, Vol. I, pp. 16, 17
- Figure 2 list of isotope names  
(1) turn page, if necessary, by touching PAGE with the light pen,  
(2) select isotope,  
(3) press END key.
- Figure 4 list of reaction types  
(1) select reaction type with the light pen,  
(2) supply energy limits using the key board,  
(3) press END key.
- Figure 5 list of energies (for SGNZ, SGIZ only)  
(1) turn page, if necessary, by touching PAGE with the light pen,  
(2) select energy,  
(3) press END key.

#### 4.2.1.4 Job control cards and input for NAP

For NAP input see also Ref. /1/, Vol. II, pp. 19-21. The following example shows the job control cards for reformatting of a NEUDADA tape (to be requested from CCDN, Saclay, in "internal expanded format") in SCORE format and creation of a SCORE library of experimental data in direct-access form on disc (cf.4.2.2.3).

```
//INRC48NA JOB (0048,101,P6M10),LANGNER,CLASS=A,TIME=3
/*FORMAT PU,DDNAME=FT04F001,FORMS=STANZ
/*SETUP DEVICE=TAPE9,ID=(SCORE1,,NL)
/*SETUP DEVICE=TAPE9,ID=(DVO645,NORING,,SL)
/*SETUP DEVICE=2314,ID=GFK016
// EXEC FHG,LIB=NUSYS,NAME=NAP
//G.FTC7F001 DD SYSOUT=B
//G.FTC8F001 DD UNIT=TAPE9,VOL=SER=DVO645,DSN=NEUDA1,
// LABEL=1,DISP=OLD
//G.FTC9F001 DD UNIT=TAPE9,VOL=SER=SCORE1,LABEL=(8,NL),DISP=(,PASS),
// DSN=NAP.LABEL,DCB=(RECFM=VBS,LRECL=68,BLKSIZE=3744)
//G.FTC9F002 DD UNIT=TAPE9,VOL=SER=SCORE1,LABEL=(9,NL),DISP=(,PASS),
// DSN=FE000,DCB=*.FT09F001
//G.FTC9F003 DD UNIT=TAPE9,VOL=SER=SCORE1,LABEL=(10,NL),DISP=(,PASS),
// DSN=FE056,DCB=*.FT09F001
//G.FT17F001 DD UNIT=2314,VOL=SER=GFK016,DSN=NEUDAD.DA,
// SPACE=(3520,520),DISP=(NEW,KEEP)
//G.SYSIN DD *
16/11/71 1 1 1 -1 0
SCORE02 SECOND SCORE TAPE
FE000 NEUDADA CCDN
26000 55.847 2 1
FE056 NEUDADA CCDN
26056 55.935 2 2
```

The second example shows the job control cards needed to print a table of content of a SCORE tape and to set up a direct-access file (if no such file exists):

```
//INRO48NA JOB (0048,101,P6M10),LANGNER,CLASS=A,TIME=3
/*FORMAT PU,DDNAME=FT04F001,FORMS=STANZ
/*SETUP DEVICE=2314,ID=GFK016
/*SETUP DEVICE=TAPE9,ID=(SCORE1,NORING,,NL)
// EXEC FHG,LIB=NUSYS,NAME=NAP
//G.FT07F001 DD SYSOUT=B
//G.FT10F001 DD UNIT=TAPE9,VOL=SER=SCORE1,LABEL=(8,NL,,IN),
// DSN=NAP.LABEL,DISP=OLD,DCB=(RECFM=VBS,LRECL=68,BLKSIZE=3744)
//G.FT10F002 DD UNIT=TAPE9,VOL=SER=SCORE1,LABEL=(9,NL,,IN),
// DSN=FE000,DISP=OLD,DCB=*.FT10F001
//G.FT10F003 DD UNIT=TAPE9,VOL=SER=SCORE1,LABEL=(10,NL,,IN),
// DSN=FE056,DISP=OLD,DCB=*.FT10F001
//G.FT17F001 DD UNIT=2314,VOL=SER=GFK016,DSN=NEUDAD.LAN,
// SPACE=(3520,520),DISP=(NEW,KEEP)
//G.SYSIN DD *
18/11/71 0 0 1 1 3
8 2
SCORE001 SCORE002
11023 1C10000 7/19/67 2230
26000 1C20000 7/19/67 7485
26054 1C30000 7/20/67 238
3000 1C40000 7/20/67 1123
3007 1050000 7/20/67 1037
94241 1C6000011/27/67 16544
26000 201000018/10/71 17373
26056 2C2000018/10/71 128
```

Short input description

The structure of the experimental-data SCORE library and the related terminology are explained in Ref. /1/, Vol. II, chap. II.

first card, format (2A4,1X,I3,5I12)

date                    2 words

NF                      number of files to be converted on the NEUDADA tape,  
= 0 if no conversion is requested

IFILE                   = 0: do not write an identification file.  
= 1: write an identification file for a new SCORE tape

IDACS                   = 0: do not create a direct-access library  
= 1: create a direct-access library

IUPD                   = -1: create table of tape identifications and isotope  
                         table (these tables are neither printed nor  
                         punched if NF=0)  
= 0: the tables are to be transferred from the  
                         direct-access library disc  
= 1: the tables are to be read in

NPRINT                 number of files on the SCORE tape for which the tables  
                         are to be printed.

tape identification card, format (16A4)

(missing if IFILE=0)

TAPEX                 2 words, tape identification,  
HEAD                   14 words, arbitrary information.

tape identification and isotope table card, format (2I12)

(missing if IFILE=1)

NIS                    number of isotopes  
NTS                    number of tapes



tape identification card, format (2A4,4X,2A4,4X etc.)

TAPE(J,I)        NTS tape identifications are to be read, six per card  
                  (J= 1,2, I=1,NTS). I is the tape number.

isotope table card, format (2I12,2A4,4X,I12)

ISOTOP(I)        isotopic identification in the form Z\*1000+A,

ISNAM(I)        tape and file number of the actual and preceding  
                  data sets in the form (NT\*100+NF)<sub>act.</sub> +10000+(NT\*100+NF)<sub>prec.</sub>

DATEIS(J,I)     date, 2 words (J=1,2)

NRECS(I)        total number of sets on the file

The last two cards must be read in for all NIS isotopes.

Title card, format (A2,A3,7X,E12.8,2I12)

ZNUM            2 alphanumeric characters, proton number Z

ANUM            3 alphanumeric characters, nucleon number A

AMASS           atomic weight (<sup>12</sup>C system)

NT              tape number

NF              file number

Title card and isotope card are to be given for each isotope.

Sorting of NEUDADA data (cf. 4.2.2.3) with SORT/MERGE Utility /9/

Job control cards and input:

```
//INR048SO JOB (0048,101,P6M1A),LANGNER,MSGLEVEL=(1,1),CLASS=A
/*SETUP DEVICE=TAPE9,ID=DVC645
/*SETUP DEVICE=TAPE9,ID=(NDC429,NORING,,NL)
// EXEC PGM=SORT,PARM='MSG=AP'
//SORTLIB DD DSN=SYS1.SORTLIB,DISP=SHR
//SYSOUT DD SYSOUT=A
//SYSLIN DD UNIT=SYSDA,SPACE=(8C,(10,10))
//SYSLMOD DD UNIT=SYSDA,SPACE=(360C,(20,20,1))
//SYSUT1 DD UNIT=(SYSDA,SEP=(SORTLIB,SYSLMOD,SYSLIN)),
// SPACE=(10C,(60,20))
//SYSPRINT DD DUMMY
//SORTIN DD UNIT=TAPE9,VOL=SER=ND0429,LABEL=(2,NL),
// DCB=(RECFM=FB,LRECL=75,BLKSIZE=750),DISP=(CLD,PASS)
//SORTOUT DD DSN=LNDSOR,UNIT=TAPE9,VOL=SER=DV0645,
// LABEL=(1,SL),DISP=(,PASS),
// DCB=(RECFM=FB,LRECL=75,BLKSIZE=750)
//SORTWK01 DD UNIT=SYSDA,SPACE=(TRK,(100),,CCNTIG)
//SORTWK02 DD UNIT=SYSDA,SPACE=(TRK,(100),,CCNTIG)
//SORTWK03 DD UNIT=SYSDA,SPACE=(TRK,(100),,CONTIG)
//SYSIN DD *
SORT FIELDS=(1,9,CH,A,48,4,FL,A,6C,4,FL,A),SIZE=E16000
```

Note: SIZE=E... depends on the number of data sets in the file that is to be sorted.

#### 4.2.2 Modifications in SCØRE

The source programs for the SCØRE subroutines exist as "members" of a "partitioned data set" /10/. At present this data set consists of 23 members.

New additions are the two "members" NAP and KNDF. Changes were made in the members OVERLAY, ONE, FOUR, FIVE, SIX and EIGHT.

##### 4.2.2.1 Changes in the overlay structure

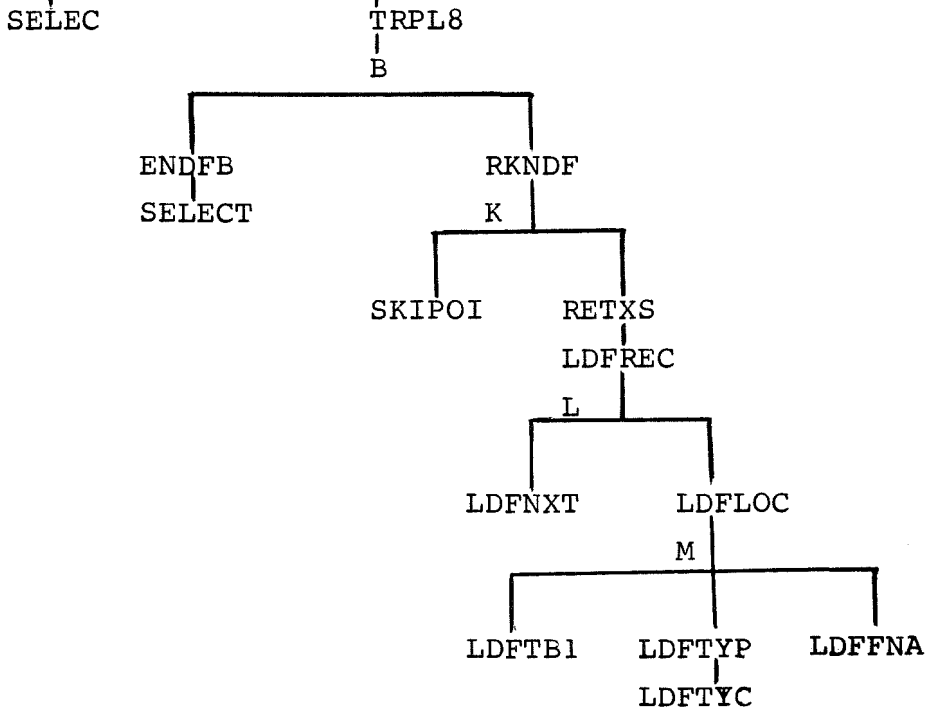
The object program of SCØRE needs 300 kBytes of memory space without, 126 kBytes with overlay. In the overlay structure the memory space is subdivided into four regions. The changes and the addition of subroutines for processing of KEDAK data resulted in the following modifications in Region I and II.

Changes in Overlay Structure

Region I

Root Segment  
LDFOPN (neu)

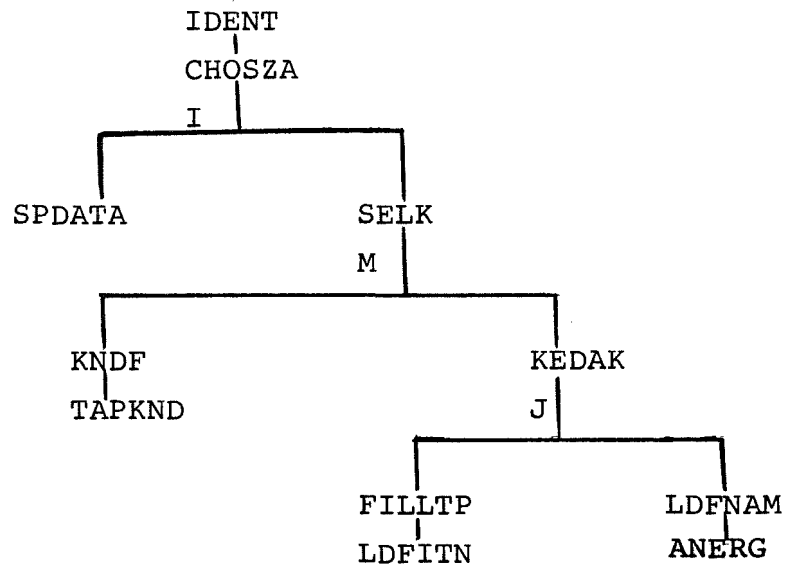
A



Region II

C (REGION)

LDFCTB  
LDFMAT  
LDFMAC



The "root segment" /11/ contains the programs MAIN, GIO, GTVDD, WTOR, DACON, CMBCD, CENTER, INFILQ, ADDBLK, EXINGR, STIMPT, PLOTG, PLOTE, CHGCHR, STOBUF, GREAD, SETKEY, ENLP, INCUR, IFNDBF, RESTSCL, SCALAR and LDFOPN.

List of all SCORE overlay statements:

```
OVERLAY A
INSERT SELEC
OVERLAY A
INSERT FIG2
OVERLAY B
INSERT SORT
OVERLAY B
INSERT DISPAR
OVERLAY A
INSERT DELETE,CORECT
OVERLAY A
INSERT DALIST,PRIND
OVERLAY A
INSERT CAFIT
OVERLAY B
INSERT MODNOD,SOND
OVERLAY B
INSERT CURVFT
OVERLAY B
INSERT LEGCAL
OVERLAY A
INSERT RESFIT
OVERLAY A
INSERT TRPLB
OVERLAY B
INSERT ENDFB,SELECT
OVERLAY B
INSERT RKNDF
OVERLAY K
INSERT SKIPOJ
OVERLAY K
INSERT RETXS,LDFREC
OVERLAY L
INSERT LDFNXT
OVERLAY L
INSERT LDFLOC
OVERLAY M
INSERT LDFTB1
OVERLAY M
INSERT LDFTYP,LDFTYC
OVERLAY M
INSERT LDFFNA
OVERLAY A
INSERT WAITR,GT2250,IHCFEXIT
OVERLAY C(REGION)
INSERT LDFCTB,LDFMAT,LDFMAC
OVERLAY C
INSERT IDENT,CHOSZA
```

OVERLAY I  
INSERT SPDATA, INPUT, LABEL, RECORD, PCINT  
OVERLAY I  
INSERT SELK  
OVERLAY H  
INSERT KNDF, TAPKND  
OVERLAY H  
INSERT KEDAK  
OVERLAY J  
INSERT FILLTP, LDFITN  
OVERLAY J  
INSERT LDFNAM, ANERG  
OVERLAY C  
INSERT RETREV, RECDRE, IDS, TAPER, FL3  
OVERLAY D  
INSERT TAPERS, RITDSK, REACT, TAB1, TABLL, TABLE3, TABLE4, TRAN1  
OVERLAY D  
INSERT FL5  
OVERLAY D  
INSERT FL4, XERIES, MUTERP  
OVERLAY C  
INSERT BLK10  
OVERLAY D  
INSERT DISPLA, BORDER, GRIDT, GRID, WPLLOT, NOPT  
OVERLAY D  
INSERT BSETUP, SPLSUP, INTERP  
OVERLAY E  
INSERT LEGEND, SERIES  
OVERLAY E  
INSERT SPLINE, CHISQR  
OVERLAY E  
INSERT ENGSPL  
OVERLAY D  
INSERT COLPSE, ANERES, EXPAND, RESISC  
OVERLAY E  
INSERT SELETY, SELRES, SELDDP, ROWSET  
OVERLAY E  
INSERT SELZA, ANEWZA  
OVERLAY E  
INSERT RESMOD, RSETUP  
OVERLAY E  
INSERT PKFIT, VALFIT  
OVERLAY D  
INSERT SAP, SIGMAS  
OVERLAY E  
INSERT CCMBCO, LAIN  
OVERLAY E  
INSERT ZEROIN, PD, DDPBR  
OVERLAY D  
INSERT ADLER, SIG, WIDTH, CHIPSI, IHCLXP, IHCLSQRT  
OVERLAY D  
INSERT DELCUR, LIMITS, CHIP, MULP, UPDTIL, ESHIFT  
OVERLAY F(REGION)  
INSERT DISPLAY, PFKEYS, WRTXTM

```
OVERLAY F
INSERT IHCSEXP,IHCFRXPR,IHCFRXPI,IHCSSCN,IHCLSCN
OVERLAY G(REGION)
INSERT ERRCDR,HELP,CENTER,CMBCD
OVERLAY G
INSERT TRACK
OVERLAY G
INSERT FLINE,PLOTP
OVERLAY G
INSERT FIG1
OVERLAY G
INSERT ROWISC,ROWTAB
OVERLAY G
INSERT CADALM,CALIN,LCML
ENTRY MAIN
NAME SCORE(R)
```

#### 4.2.2.2 SCORE routines

##### 4.2.2.2.1 MAIN PROGRAM SCORE

The changes permit access to KEDAK data /3/, /4/ from the KNDF library and display of the cross sections as curves on the screen. Call of the subroutine LDFOPN /6/ makes the KNDF library (Karlsruhe Nuclear Data File) on unit 1 available to SCORE and fills COMMON/CDFMT/ with the material names /5/ of the isotopes stored in KEDAK format. The newly added call of subroutine SELEC permits selection between the KEDAK and NEUDADA option. The option KEDAK was added to the option list MAIN 3/3 /1/. Call of the subroutine KNDF makes the KEDAK data available for display in curve form simultaneously with NEUDADA data. Call of the subroutine KEDAK yields KEDAK data also if experimental data are not available or not transferred to core storage.

##### 4.2.2.2.2 Subroutines ERRCDR, SPDATA, CHOSZA

An error message for reading of KEDAK data was added to subroutine ERRCDR. In subroutine SPDATA the section for the figure "Case Identification" (cf. /1/, Vol. II, p. 16) was replaced by subroutine IDENT (cf. 4.2.3.1.2). Subroutine CHOSZA, which provides for interactive choice of isotope and reaction type on the screen, was changed so that also the 8-byte alphameric KEDAK isotope and reaction type names /5/ can be processed.

#### 4.2.2.3 Program NAP - compilation of SCORE library from experimental (NEUDADA) data

NAP (NEUDADA Adaptation Program) was rewritten and tested.

Experimental data from other centers reach Karlsruhe via CCDN (OECD/NEA Centre de Compilation de Données Neutroniques), Saclay. The data are transmitted on magnetic tape in NEUDADA format /7/.

The user must sort the data received on a NEUDADA tape in ascending order of Z, A and energy with the SORT/MERGE utility program /8/ (see 4.2.1.4), if necessary. The NAP program reformats the data thus prepared and writes them onto a library tape. This library tape contains an identification record and a directory of addresses for the content of the tape. The data themselves are organized in such a way that one data block corresponds to one record of the direct-access library which is also generated by NAP. This library contains a table of available isotopes and the tape names and file numbers of the corresponding library tapes. This table is updated by NAP. Transfer of the data from tape to direct-access library is performed by SCORE /1/ if necessary.

#### 4.2.3 New subroutines for SCORE

##### 4.2.3.1.1 SELEC - choice between NEUDADA and KEDAK

Subroutine SELEC, called from MAIN PROGRAM SCORE, generates the first figure for SCORE. The following text is displayed on the screen:

SCORE

CROSS SECTION EVALUATION SYSTEM  
ADAPTED TO NEUDADA AND KEDAK DATA

SELECT BY LIGHT PEN DATA TYPE  
TO BE RETRIEVED FIRST

KEDAK

NEUDADA

Thus the user can choose between KEDAK or NEUDADA data. If he chooses the NEUDADA option the SCORE version described in /1/ is at his disposal with the additional possibility to display KEDAK data in curve form together with the experimental NEUDADA data - provided that data for the chosen isotope, reaction type and energy range exist in the KNDF library /3/, /4/. If he chooses the KEDAK option he gets the cross sections from the KNDF library displayed on the screen, regardless whether NEUDADA data are available for the chosen isotope, reaction type and energy range, or not.

The option chosen is displayed in large size on the screen. Pressing of the END key or the equivalent key no. 1 on the programmed-function keyboard causes transmittal of the information to the main program. Erroneous pressing of the END key without prior selection of the format has no consequence, SELEC will still wait for the light pen.

SELEC is called by

```
CALL SELEC(IOPT),
```

where IOPT is set by SELEC. (IOPT=1: KEDAK, IOPT=2: NEUDADA.)

SELEC utilizes the following subroutines from the "graphic subroutine package" of SCORE: INPICT, WRTXTM, SETKEY, ENDKEY, STPICT, EXINGR, CHGCHR whose functions are as follows:

INPICT	initializes the figure,
WRTXTM	writes a character string onto a specified place on the screen,
SETKEY	indicates the beginning of a series of graphic commands associated with a key,
ENDKEY	indicates the end of the series,
STPICT	brings a block of graphic commands into the buffer of the IBM 2250 and starts the figure,



EXINGR            controls the input to the IBM 2250,  
CHGCHR            changes the mode of a character string from protected  
                  to unprotected and vice versa

#### 4.2.3.1.2 IDENT - user and case identification

The subroutine IDENT is essentially a program section taken out of subroutine SPDATA so as to be available also to KEDAK. More specifically it is the part that generates the first figure for SCORE (cf. /1/, Vol. I, p. 16) in which the case identification is requested.

IDENT is called by

```
CALL IDENT(IER),
```

where IER is an error flag:

IER = 1 : no error, continue;  
IER = 2 : jump back to the start of the program;  
IER = 3 : error, terminate execution.

IDENT calls the SCORE routines FIG1, DACON, ERRCOR and utilizes COMMON/BLK1/:

FIG1            generates the figure,

DACON            converts numbers from the internal fixed- or floating-point  
                  representation to alphameric or vice versa,

ERRCOR          is a routine for error messages and error response.

```
COMMON/BLK1/ABCD(5,4),ICASE,CNUM
```

The array ABCD contains user identification and date, ICASE is the case number and CNUM the figure number for this case.

#### 4.2.3.2 Subroutines for the organization of KEDAK data

##### 4.2.3.2.1 KNDF - KEDAK data for curve display

Subroutine KNDF organizes cross sections from the KNDF library /3/, /4/ for curve display together with experimental data. Proton number and reaction type number are taken over from the experimental data (option NEUDADA), likewise the energy range is fixed by the experimental data. For differential elastic-scattering cross sections the data for the requested energy are calculated, if necessary, by linear interpolation between the angular distributions existing on KEDAK.

Three quantities are taken from COMMON/BLK4/:

ENER            incident energy for the differential elastic-scattering cross section,

QX             excitation energy for the partial inelastic-scattering cross section,

NREC           KEDAK number of the reaction type chosen.

The energy limits EMIN, EMAX are taken from COMMON/BLK5/. Subroutine KNDF stores cross sections from the KNDF library in COMMON/BLK8/NPØINT,X,Y,NKNDF,ZLIM(4), where

NPØINT        is the number of energies,

X             is the abscissa (energy) array,

Y             is the ordinate (cross section) array,

NKNDF        is taken as equal to NPØINT,

ZLIM         are the minimum and maximum energy and cross section values calculated by subroutine CALIMA.

Subroutine KNDF utilizes subroutines TAPKND,RKNDF (cf. 4.2.3.2.2, 4.2.3.2.3) and the SCORE subroutines CALIMA and ERRCØR. It is called by

```
CALL KNDF(IER)
```

where IER is an error flag as explained in 4.2.3.1.2 above.

4.2.3.2.2 TAPKND - calculation of numeric KEDAK names

Subroutine TAPKND calculates numeric KEDAK names /5/ for the isotope and the reaction type selected for the NEUDADA data for which the KEDAK curve is to be displayed together with the experimental values. The isotope name is calculated from the proton number (IZ) and the nucleon number (IA) as follows

$$\text{NAMES}(1) = \text{IZ} * 10000 + \text{IA}.$$

The reaction type number is calculated as

$$\text{NAMES}(2) = \text{MF} * 10000 + \text{MT} * 10 \quad \text{with MF} = 3$$

according to the ENDF conventions /2/ for all reaction types listed in table MTS (see below), with the exception of the differential elastic-scattering cross section and the inelastic-scattering cross section for a specific excitation of the residual nucleus (NREC = 14, NREC = 15). For these latter reaction types one has

$$\text{NAMES}(2) = \text{MF} * 10000 + \text{MT} * 10 + 2 \quad \text{with MF} = 4.$$

MT is determined by NREC (number of the reaction type in table MTS) from COMMON/BLK4/ and table MTS. The following reaction types are available in SCORE at present:

MTS ENDF/B	KEDAK numeric	KEDAK alphanumeric	Error message from ERRCOR on screen <sup>+) </sup>
107	31070	SGALP	SGAL
104	31040	SGD	SGD
19	30190	SGF	SGF
102	31020	SGG	SGG
106	31060	SGHE3	SHE3
103	31030	SGP	SGP
105	31050	SGH3	SGH3
4	30040	SGI	SGI
16	30160	SG2N	SG2N

<sup>+)</sup> For error messages on screen the type name must be shortened to four characters where necessary.

MTS ENDF/B	KEDAK numeric	KEDAK alphanumeric	Error message from ERRCØR on screen <sup>+</sup> )
17	30170	SG3N	SG3N
2	30020	SGN	SGN
5	30050	SGIZ	SGIZ
1	30010	SGT	SGT
2	40022	SGNC	SGNC
5	40052	SGICZ	SICZ

<sup>+</sup>) For error messages on screen the type name must be shortened to four characters where necessary.

TAPKND is called by

```
CALL TAPKND(NAMES,NZAM,MF,IQ,IEF)
```

where

NAMES is a 5-word array with the KEDAK names inserted in words 1 and 2 by the subroutine,

NZAM is the number of KEDAK names as set by TAPKND,

IQ is the additional KEDAK identifier for angular distributions or inelastic-scattering cross sections,

IEF is an error code (IEF=0 without errors).

#### 4.2.3.2.3 RKNDF - reading of KEDAK data

Subroutine RKNDF prepares the KEDAK data for the energy range EMIN...EMAX in the arrays X and Y. If there are more than 150 energy values all those curve points are deleted by subroutine SKIPØI which are not required for a 1 % accuracy of the cross section curve. The cross sections for the lower and upper energy limits EMIN, EMAX are calculated by linear interpolation in the SCØRE subroutine TRPL8. Differential elastic-scattering cross sections are prepared by subroutine SELK (see 4.2.3.2.5). Reading of KEDAK data from the KNDF file is controlled by subroutine RETXS /6/. Thus the subroutines called by RKNDF are SELK,RETXS,SKIPØI,TRPL8.

RKNDF is called by

```
CALL RKNDF (NAMES, NZNAM, MF, EMIN, EMAX, X, Y, NPØINT, IEF)
```

where

NAMES is the field for the KEDAK names of isotope and reaction type,  
NZNAM is the number of KEDAK names,  
MF = 3 or 4 according to ENDF conventions /2/,  
EMIN is the lower limit of the requested energy range,  
EMAX is the upper limit of the requested energy range,

Note that these arguments must be available, whereas the following arguments are supplied by RKNDF:

X 1-dimensional array of KEDAK energies,  
Y 1-dimensional array of KEDAK cross sections,  
NPØINT number of data points transferred from the KNDF library  
IEF error code for subroutine ERRCØR (IEF=0 without error).

#### 4.2.3.2.4 SKIPØI - deletion of data points

Subroutine SKIPØI deletes energy grid points of the KEDAK data if their number exceeds a limit that depends on the array dimensions in SCØRE (150 at present). Only those points are deleted which are not required for a curve which represents the KEDAK values with  $\epsilon = 1\%$  accuracy.

SKIPØI is called by

```
CALL SKIPØI (XX, YY, NUMX, X, Y, NPØINT, LENX, EPS)
```

where

XX is a 1-dimensional array for the energy values,  
YY is a 1-dimensional array for the cross section values,  
NUMX is the number of XX values (or YY values),

X is the energy array after deletion of superfluous values,  
Y is the cross section array after deletion of superfluous values,  
NPØINT is the number of X (or Y) values,  
LENX is the maximum number of X (or Y) values  
EPS is the accuracy limit  $\epsilon$ .

SKIPØI is repetitively called until the X (and Y) array is filled completely.

#### 4.2.3.2.5 SELK - calculation of the differential elastic-scattering cross section

The differential elastic-scattering cross sections are calculated from the angle-integrated elastic-scattering cross section  $\sigma_n(E)$  and the elastic angular distributions (center-of-mass system)  $p(\mu, E)$  that are stored in the KEDAK library /3/, /4/. For a given incident energy E the angle-integrated cross section is calculated by linear interpolation, if necessary. The angular distribution is also determined by linear interpolation between the distributions stored on KEDAK. Here the interpolation is between scattering-angle cosines  $\mu$  on one hand and incident energies on the other. The angular grid is taken from the lowest incident energy on KEDAK. The differential elastic-scattering cross section is then

$$\frac{d\sigma_n(E, \Omega)}{d\Omega} = \sigma_n(E) \cdot p(\mu, E),$$

in units of mb/sterad. Linear interpolation is performed by the SCØRE subroutine TRPL8. The subroutines RETXS, IDFLOC, IDFNXT /6/ are used for reading of KEDAK data.

SELK is called by

```
CALL SELK(NARG, NAMEN, EMI1, EMI2, XX, YY, MAXNUM, NUM1, IEF)
```

where

NARG is the number of KEDAK names for the requested isotope,  
NAMEN is an array for the KEDAK names in numeric representation,  
EMI1 is the lower limit of the requested energy range,  
EMI2 is the upper limit of the requested energy range,

XX is the array of scattering-angle cosines,  
YY is the array for the differential elastic-scattering cross sections,  
MAXNUM is the maximum number of cosines,  
NUM1 is the actual number of cosines,  
IEF is an error code for the SCORE routine ERRCOR (IEF=0 without errors)

#### 4.2.3.2.6 KEDAK - KEDAK data without experimental (NEUDADA) data

Representation of KEDAK data independent of NEUDADA data (option KEDAK) is organized in subroutine KEDAK.

First KEDAK calls subroutine IDENT which supplies user and case identification. The list of isotope names on KNDF is written into COMMON/LDFMT/ by LDFOPN /6/ and thus transferred to the modified SCORE routine CHOSZA via entry CHOSMA. This permits display on the screen and selection of the desired isotope by light pen.

Subroutine FILLTP /6/ yields the alphameric and numeric /5/ reaction type names for the selected isotope. The alphameric names are displayed on the screen via entry CHOSRE in CHOSZA so that a reaction type can be chosen. In case SGNC (the differential elastic-scattering cross section) or SGIZ (the partial inelastic-scattering cross section) is selected subroutine LDFNAM /6/ is called. This subroutine supplies the incident energies (for SGNC) or excitation energies (for SGIZ) for which the KNDF file contains data. Subroutine ANERG lists these energies on the screen allowing for selection of the appropriate energy by light pen.

Subroutine RKNDF (cf. 4.2.3.2.3) supplies the KEDAK data for the chosen isotope, reaction type, energy range (and - for SGNC - incident energy or - for SGIZ - excitation energy).

Subroutine CALIMA provides minimum and maximum energy and cross section values for the plot. Isotope and reaction type names and case number are written into the appropriate field for figure caption and printout.

Subroutine KEDAK utilizes the subroutines FILLTP,LDFNAM /6/, subroutine IDENT (cf. 4.2.3.1.2), RKNDF (cf. 4.2.3.2.3) and ANERG (cf. 4.2.3.2.7) and the SCORE subroutines CHØSZA, DACØN, CENTER and ERRCØR. It is called by

```
CALL KEDAK(IER)
```

where IER is an error code (see 4.2.3.1.2).

The KEDAK data are stored in CØMMØN/BLK8/ (cf. 4.2.3.2.1).

#### 4.2.3.2.7 ANERG - listing of energies on the screen

Subroutine ANERG displays a list of either

- incident energies for differential elastic-scattering cross sections (reaction type SGNC), or
- excitation energies for partial inelastic-scattering cross sections (reaction type SGIZ),

as available on KEDAK. The desired energy can then be selected by light pen. The display contains isotope and reaction type name, 15 energies and the word PAGE. Touching the word PAGE with the light pen the user can switch to the next 15 energies. If there are no further energies for the requested isotope the first 15 energies are shown again etc. (At present up to 200 different energies can be displayed in this way.)

The selected energy is displayed in large size. The information is transferred to subroutine KEDAK by pressing of the END key.

ANERG utilizes the following subroutines of the "graphic subroutine package" of SCORE: INPICT,WRTCTM,SETKEY,ENDKEY,STPICT,EXINGR,CHGCHR and the SCORE routine DACØN. It is called by

```
CALL ANERG(NAM1,N2,XNAM1,NNAM,IER)
```



where

NAM1 is a field of 2 double words which must contain the alpha-  
meric KEDAK isotope and reaction type names /5/,  
N2 is the number of energies,  
XNAM1 is an array containing the energies,  
NNAM is the running number of the selected energy as  
determined by ANERG,  
IER is an error code (IER=0 without errors).

References

- /1/ C.L. Dunford, SCORE, Vol. I und II, Atomics International, AI-AEC-12994, 1971
  
- /2/ M.K. Drake, Data Formats and Procedures for the ENDF Neutron Cross Section Library National Neutron Cross Section Center, BNL 50274 (T-601), 1970
  
- /3/ J.J. Schmidt, Neutron Cross Sections for fast Reactor Materials, 1966, KFK 120
  
- /4/ B. Goel, report KFK 2234 (1975)
  
- /5/ D. Woll, Card Image Format of the Karlsruhe Evaluated Nuclear Data File 1968, KFK 880 (1968)
  
- /6/ Present Compendium, Chs. III.2, III.4
  
- /7/ NEUDADA System Description, NEA/CCDN Newsletter CCDN/SYS-2, 1969
  
- /8/ IBM SYSTEM/360, Operating System SORT/MERGE, Form C28-6543
  
- /9/ IBM SYSTEM/360 Component Description IBM 2250 Display Unit Model 1, Form A27-2701-2
  
- /10/ IBM SYSTEM/360 Operating System Supervisor and Data Management Services, Form GC28-6646
  
- /11/ IBM SYSTEM/360 Operating System Linkage Editor, Form C28-6538
  
- /12/ H. Alter and C.L. Dunford, Proc. 2nd Intern. Conf. on Nuclear Data for Reactors, Helsinki, 1970; Vol. II, p. 643

Appendix:

A1. List of changes in SCORE routines

MEMBER ONE - MAIN PROGRAM

```
./ DELETE SEQ1=01003600,SEQ2=01003600
  DATA TASK3 /*MAIN 3/3RESTART KEDAK DATA FITENDF/B RES CALC 01003600
./ NUMBER SEQ1=01004750,NEW1=01004800,INCR=100,INSERT=YES
  CALL LDFOPN(1,NDAT,&1000)
./ NUMBER SEQ1=01007000,NEW1=01007100,INCR=100,INSERT=YES
  CALL SELEC(IEF)
  ICATA=IER
  GO TO (21,22),IER
```

```
C
  21 CALL KEDAK(IEF)
  GO TO (30,10,1000),IER
./ DELETE SEQ1=01007300,SEQ2=01007300
  22 CALL SPDATA(IFR)
  01007300
./ DELETE SEQ1=01009500,SEQ2=01009500
  60 GO TO (35,10,36,300,350,400,35,120,35,1000),INDM
  01009500
./ DELETE SEQ1=01012900,SEQ2=01012900
  135 CALL PRIND(ICATA)
  01012900
./ NUMBER SEQ1=01017700,NEW1=01017800,INCR=100,INSERT=YES
C
C KNDF OVERLAY
C
  36 IER=1
  IF(ICATA.EQ.1) GO TO 21
  CALL KNDF(IEF)
  GO TO (425,10,1000),IFR
```

MEMBER ONE - BLOCK DATA

```
./ DELETE SEQ1=01020500,SEQ2=01020500
  DATA VERDAT/'AUGUST 1972'/,ABCD/20*' /*,NRECP,IMT/13,0/,
  01020500
```

MEMBER ONE - SUBROUTINE ERRCOR

```
./ NUMBER SEQ1=01052800,NEW1=01052900,INCR=100,INSERT=YES
  DATA EFR13/'* NO DATA FOUND FOR ON KNDF
  8 /*
  IF(IEF.EQ.13) GO TO 11
  GO TO 12
  11 ERR13(6)=RTYPE
  CALL IDACDN (ERR13(7),DUM1(14),2,1,1,1,IFR)
  12 IER=1
```

MEMBER FOUR - SUBROUTINE GRID

```
./ NUMBER SEQ1=04025800,NEW1=04025900,INCR=100,INSERT=YES
  IF(JEXP.LT.10) GO TO 21
./ NUMBER SEQ1=04025800,NEW1=04025900,INCR=100,INSERT=YES
  GO TO 22
  21 WRITE (99,23) BL,TEN,S,JEXP,BL
```

23 FORMAT (A1,A4,A1,I1,A1)  
22 CALL INFILQ (A,I2)

MEMBER FIVE - SUBROUTINE SPDATA

```
./ NUMBER SEQ1=05003400,NEW1=05003500,INCR=100,INSERT=YES
  IDATA=IER
./ DELETE SEQ1=05004400,SEQ2=05006000
  2 'ATOMICS INTERFNATIONAL AND IBM - PALO ALTO'////29X,'GFK - KARLSR05004400
  3UFE VERSION III-K',5X,3A4) 05004500
C 05004600
  CALL IDENT (IER) 05004700
  IF(IER.NE.1) GO TO 1000 05004800
./ DELETE SEQ1=05008700,SEQ2=05008700
  12 CALL CHOSZA (IZT,IAT,NISO,I,IDATA) 05008700
./ DELETE SEQ1=05023300,SEQ2=05023300
  75 CALL CHOSRT (IRT,NRT,BCDRT,RTYPE,NREC,EL,IDATA) 05023300
```

MEMBER FIVE - SUBROUTINE CHOSZA

```
./ DELETE SEQ1=05047100,SEQ2=05047100
  SUBROUTINE CHOSZA (IZ,IA,N,I,IDAT) 05047100
./ DELETE SEQ1=05048300,05048300
  DIMENSION TITLE(9),TITLE1(12),TITLE2(8,3),IZ(1),IA(1),XKED(3), 05048300
./ NUMBER SEQ1=05048400,NEW1=05048500,INCR=100,INSERT=YES
  REAL*8 NAME,SGIZ,XZ(1),R1,XB(1)
  EQUIVALENCE (TITLE1(9),NAME)
./ DELETE SEQ1=05048500,SEQ2=05048500
  DATA TITLE/'ISOTUPES AVAILABLE FROM NEUDADA ARE '/,TITLE1/'REACTIO5048500
./ DELETE SEQ1=05048700,SEQ2=05048700
  2X1,X4,Y1,Y4/2.,2.,10.75,2./,Y0,DY1,X5/9.5,.5,4./,X00,Y00/4.5,10./ 05048700
./ NUMBER SEQ1=05049000,NEW1=05049000,INCR=100,INSERT=YES
  DATA XKED/'KEDAK ARE '/,BLANK/' '/,SGIZ/'SGIZ '/
C
  DIMENSION XEUD(3),TIT1(5)
  DATA XEUD/'NEUDADA ARE '/,TIT1/' Z = A = ARE'/'
  DO 40 L=1,3
  40 TITLE(L+6)=XEUD(L)
  DO 50 L=1,5
  50 TITLE1(L+7)=TIT1(L)
./ NUMBER SEQ1=05049400,NEW1=05049500,INCR=100,INSERT=YES
  ENTRY CHOSMA (XZ,IA,N,I,IDAT)
  IST=(I-1)/15
  25 CALL INPICT
  IF(ICAT.EQ.2) GO TO 1
  TITLE(7)=XKFD(1)
  TITLE(8)=XKED(2)
  TITLE(9)=XKED(3)
  1 CALL WRTXTM(2,36,TITLE,X1,Y1)
  IF(ICAT.EQ.1) GO TO 2
./ NUMBER SEQ1=05052050,NEW1=05052100,INCR=100,INSERT=YES
  GO TO 3
  DISPLAY MATERIAL NAMES FROM KEDAK
  2 Y2=YC
  IL=IST*15+1
  IF (IL.LE.N) GO TO 4
  IL=1
```

```
IST=0
4 IU=IL+14
  IU=MINO(IU,N)
  NN=IU-IL+1
  CALL SETKEY (7,1,8)
  DO 10 K=IL,IU
  ISZ=1
  IF(I.EQ.K) ISZ=2
  NAME=XZ(K)
  TITLE1(8)=BLANK
  CALL WRTXTM(ISZ,8,TITLE1(9),X5,Y2)
10 Y2=Y2-DY1
  3 CALL ENDKEY
./ NUMBER SEQ1=05055200,NEW1=05055300,INCR=100,INSERT=YES
  IF(ICAT.EQ.1) GO TO 11
./ NUMBER SEQ1=05055500,NEW1=05055600,INCR=100,INSERT=YES
11 NAME=XZ(I)
  GO TO 1000
./ NUMBER SEQ1=05055700,NEW1=05055800,INCR=100,INSERT=YES
  ENTRY CHDSRE (IR,NR,XB,R1,NREC,XX,IDAT)
./ NUMBER SEQ1=05058200,NEW1=05058300,INCR=100,INSERT=YES
  L=4
  IF(ICAT.EQ.1) L=8
  CALL SETKEY (7,1,L)
./ NUMBER SEQ1=05058500,NEW1=05058600,INCR=100,INSERT=YES
  NRE=16
  IF(ICAT.EQ.1) NRE=NR
./ DELETE SEQ1=05058800,SEQ2=05058800
  DO 250 I=1,NRE
./ NUMBER SEQ1=05059300,NEW1=05059400,INCR=100,INSERT=YES
  IF(ICAT.EQ.1) GO TO 5
./ NUMBER SEQ1=05059500,NEW1=05059600,INCR=100,INSERT=YES
  5 CALL WRTXTM (1,8,XB(I),X3,Y3)
  GO TO 6
./ NUMBER SEQ1=05061400,NEW1=05061500,INCR=100,INSERT=YES
  IF(ICAT.EQ.1) GO TO 7
./ NUMBER SEQ1=05061600,NEW1=05061700,INCR=100,INSERT=YES
  7 R1 = XB(I)
  GO TO 380
./ NUMBER SEQ1=05063700,NEW1=05063800,INCR=100,INSERT=YES
  IF(R1.EQ.SGIZ) I2=3
```

MEMBER SIX - SUBROUTINE PRIND

```
./ DELETE SEQ1=06060700,SEQ2=06060700
  SUBROUTINE PRIND (IDAT)
./ DELETE SEQ1=06065400,SEQ2=06065400
  30 IF(IDAT.EQ.1) GO TO 240
  WRITE (6,35) ABCRE
./ DELETE SEQ1=06069000,SEQ2=06069000
  240 IF(NENDFB.EQ.0) GO TO 1000
  IF(IDAT.EQ.1) GO TO 270
./ NUMBER SEQ1=06069200,NEW1=06069200,INCR=100,INSERT=YES
  GO TO 260
  270 WRITE (6,280)
  280 FORMAT (////////30X,'KEDAK DATA'///)
  260 IT=1
```

06060700  
06065400  
06065500  
06069000  
06069100

MEMBER EIGHT - SUBROUTINE SOND

```
./ DELETE SEQ1=08030600,SEQ2=08030700
  CALL ROUND (ZX,1,5)
  CALL ROUND (ZX,1,5)
```

08030600  
08030700

MEMBER NAP - PROGRAM NAP

```
./ DELETE SEQ1=51701300,SEQ2=51701300
  COMMON /BLK176(16,55),C(16),IV,ITSC,IR,IC,NPRINT,RECT(175),
./ NUMBER SEQ1=51702300,NEW1=51702400,INCR=100,INSERT=YES
  DATA BCDN/'N'/
C
  DEFINE FILE 17(520,880,U,IV)
./ DELETE SEQ1=51702700,SEQ2=51702700
  READ (IN,500) DATX,NF,IFILE1,IDACS,IUPD,NPRINT
./ DELETE SEQ1=51704900,SEQ2=51705900
./ DELETE SEQ1=51707800,SEQ2=51708200
  WRITE (17'1) NIS,NTS,(ISOTOP(I),ISNUM(I),DATEIS(1,I),DATEIS(2,I),
  1 NRECS(I),I=1,NIS),(TAPE(1,J),TAPE(2,J),J=1,NTS)
./ DELETE SEQ1=51708800,SEQ2=51709100
  5 READ (17'1) NIS,NTS,(ISOTOP(I),ISNUM(I),DATEIS(1,I),DATEIS(2,I),
  1 NRECS(I),I=1,NIS),(TAPE(1,J),TAPE(2,J),J=1,NTS)
./ NUMBER SEQ1=51712400,NEW1=51712500,INCR=100,INSERT=YES
  BLK=PAC(BT(7),BL,BL,BL)
./ DELETE SEQ1=51734800,SEQ2=51735100
  WRITE (17'1) NIS,NTS,(ISOTOP(I),ISNUM(I),DATEIS(1,I),DATEIS(2,I),
  1 NRECS(I),I=1,NIS),(TAPE(1,J),TAPE(2,J),J=1,NTS)
```

51701300  
51702700  
51707800  
51707900  
51708800  
51708900  
51734800  
51734900

MEMBER NAP - SUBROUTINE PRINT

```
./ DELETE SEQ1=51741700,SEQ2=51741700
./ DELETE SEQ1=51745100,SEQ2=51745100
  38 IU=55
```

51745100

MEMBER NAP - SUBROUTINE CREATE

```
./ DELETE SEQ1=51754000,SEQ2=51754000
```

MEMBER NAP - SUBROUTINE RDISK

```
./ DELETE SEQ1=51756200,SEQ2=51756200
```

MEMBER NAP - SUBROUTINE DPOINT

```
./ DELETE SEQ1=51757900,SEQ2=51760900
./ ENDUP
```

Appendix:

A2. Source program lists of the new subroutines

SELEC  
IDENT  
KNDF  
RKNDF  
TAPKND  
SELK  
RETXS  
SKIPOI  
KEDAK  
ANERG  
FILLTP



```
      SUBROUTINE SELEC (IOPT)
C
C   PERMITS THE SELECTION OF NEUDADA OR KEDAK DATA
C
      DIMENSION TEXT(14),TEXT1(2,2),TITLE(19)
      DATA Y1/5./,DY/.75/,X1/2./,TEXT1/'KEDA','K ','NEUD','ADA '/
      DATA TITLE/'SCORE  CROSS SECTION EVALUATION SYSTEM ADAPTED TO NEU
IDADA AND KEDAK DATA  '/
      DATA TEXT/'SELECT BY LIGHT-PEN DATA TYPE  TO BE RETRIEVED FIRST
1  '/
C
      ICPT=0
      IC = 0
C
C   INITIALIZE DISPLAY AND WRITE TEXT
C
1  CALL INPICT
   CALL WRTXTM (2,8,TITLE(1),5.,10.75)
   CALL WRTXTM (2,32,TITLE(3),2.,10.)
   CALL WRTXTM (2,36,TITLE(11),2.,9.25)
   CALL WRTXTM (2,32,TEXT(1),2.,7.75)
   CALL WRTXTM (2,24,TEXT(9),2.,7.)
   CALL SETKEY (7,1,8)
   Y=Y1
   DO 10 I=1,2

      ISZ=1
      IF(ID.EQ.I) ISZ=2
      CALL WRTXTM(ISZ,8,TEXT1(1,I),X1,Y)
10  Y=Y-DY
      CALL ENDKEY
C
C   GENERATE DISPLAY AND AWAIT USER ACTION
C
      CALL STPICT
20  CALL EXINGR (IKEY,IVAL,IOPT)
      GO TO (3,2,1,1000,20,20),IKEY
C
C   PROCESS USER ACTION AND RETURN
C
3  IF(IVAL.GT.2) GO TO 20
   IF(IOPT.EQ.0) GO TO 1
   GO TO (1000,1),IVAL
2  IF(IOPT.GT.2) GO TO 20
   IF(IO.LT.1.OR.ID.GT.2) GO TO 5
   CALL CHGCHR (IVAL,IO,1)
5  CALL CHGCHR (IVAL,IOPT,2)
   IC=IOPT
   GO TO 20
1000 RETURN
      END
```

```
C      SUBROUTINE IDENT (IER)
C
      DIMENSION T1(4,4)
      COMMON/BLK1/ABCD(5,4),ICASE,CNUM,DUM1(43)
      DATA T1/'LABORATORY:      EVALUATOR:      DATE:      PROBLEM I
1.D.:      '/
C
C      USER SUPPLIES IDENTIFICATION FOR PROBLEM
      NOUT=6
1  CALL FIG1 (4,16,20,T1,ABCD,IKEY)
      CALL DACON (ABCD(1,4),CASE,5,1,1,0,2,IE)
      IF(IE.EQ.0) GO TO 2
      CALL ERRCOR (8,IER)
      GO TO (1,1000,1000),IER
2  ICASE=100*IFIX(CASE+0.1)
      CALL IDACON (CNUM,ICASE,1,1,1,1,IE)
      WRITE (NOUT,101) ICASE
101 FORMAT(1H1,/' CASE NUMBER',2X,I8//)
      DC 10 I=1,3
      WRITE (NOUT,102) (T1(J,I),J=1,4),(ABCD(K,I),K=1,5)
102 FORMAT (1H0,5X,4A4,5X,5A4)
      10 CONTINUE
1000 RETURN
      END
```

```
MEMBER NAME KNDF
SUBROUTINE KNDF (IER) 25400100
C 25400200
C THIS IS THE CONTROL ROUTINE FOR THE HANDLING OF THE KNDF-DATA 25400300
C 25400400
COMMON /BLK4/ DUM1(3),ENER,QX,NREC,NRECP,IZ,IAND,IZA,DUM2(5) 25400500
COMMON /BLK5/ DUM5(4),XPLOTP(3),DUM6(7) 25400600
COMMON /BLK8/ NPCINT,X(150),Y(150),NKNDF,ZLIM(4) 25400700
DIMENSION NAMES(5) 25400900
EQUIVALENCE (Q,IQ) 25401000
C 25401100
IEF=0 25401200
IER=1 25401300
NPCINT=0 25401400
C 25401500
IF(NREC.GT.NRECP) GO TO 1 25401600
EMIN=XPLOTP(1)*1.0E+06 25401700
EMAX=XPLOTP(2)*1.0E+06 25401800
GO TO 2 25401900
C 25402000
1 EMIN=ENER*1.0E+06 25402100
EMAX=0. 25402200
2 Q=QX*1.0E+06 25402300
C 25402400
CALL TAPKND (NAMES,NZNAM,MM,IQ,IEF) 25402500
IF(IEF.NE.0) GO TO 55 25402600
C 25402700
3 CALL RKNDP(NAMES,NZNAM,MM,EMIN,EMAX,X,Y,NPCINT,IEF) 25402800
IF(IEF.NE.0) GO TO 55 25402900
C 25403000
CHANGE ENERGY UNIT FROM EV TO MEV 25403100
C 25403200
IF(MM.EQ.4) GO TO 5 25403300
DO 20 I=1,NPCINT 25403400
20 X(I)=X(I)*1.0E-06 25403500
C 25403600
6 IF(X(2).LT.1.000001*X(1).AND.X(2).GT.0.999999*X(1)) GO TO 5 25403700
IF(Y(1).NE.0.0) GO TO 4 25403800
5 DO 30 I=2,NPCINT 25403900
I1=I-1 25404000
X(I1)=X(I) 25404100
30 Y(I1)=Y(I) 25404200
NPCINT=NPCINT-1 25404300
C 25404400
4 CONTINUE 25404500
C 25404600
IF(X(1).LT.XPLOTP(1).AND.NREC.LE.NRECP) X(1)=XPLOTP(1) 25404700
1000 NKNDF=NPCINT 25404800
C 25404900
CALL CALIMA (NPCINT,X,Y,ZLIM) 25405000
C 25405100
100 RETURN 25405200
55 CALL ERRCOR (IEF,IER) 25405300
GO TO 100 25405400
END 25405500
C 25405600
SUBROUTINE TAPKND (NAMES,NZNAM,MM,IQ,IEF) 25405700
C 25405800
THIS ROUTINE CALCULATES THE KNDF-NAMES 25405900
```

```

MEMBER NAME  KNDF
C   OF THE MATERIAL AND THE REACTION TYPE                25406000
C                                                         25406100
C   INTEGER DUM3                                          25406200
COMMON /BLK4/ DUM1(4),IQX,NREC,NRECP,IZ,IA,IZA,DUM3(4),RTYPE 25406300
C   DIMENSION NAMES(5),NNAM(16),MTS(16),TYPNA(16)        25406400
C                                                         25406500
C   DATA NNAM/11*2,1*3,1*2,1*3,1*4,1*0/                25406600
C                                                         25406700
C   DATA TYPNA/'SGAL','SGD ','SGF ','SGG ','SFE3','SGP ','SGH3','SGI ' 25406800
C 2,'SG2N','SG3N','SGN ','SGIZ','SGT ','SGNC','SICZ','  '/ 25406900
C                                                         25407000
C   DATA MTS /107,104,19,102,106,103,105,4,16,17,2,5,1,2,5,29/ 25407100
C                                                         25407200
C   DC 30 I=1,5                                           25407300
30 NAMES(I)=0                                             25407400
C   NAMES(1)=IZ*10000+IA                                  25407500
C   NAMES(3)=IQ                                           25407600
C                                                         25407700
C                                                         25407800
C   DC 20 I=1,16                                          25407900
C   MT=MTS(I)                                             25408000
C   ITY =I                                                25408100
C   NZNAM=NNAM(I)                                         25408200
C   IF(NREC.EQ.1) GO TO 2                                  25408300
C                                                         25408400
C   20 CONTINUE                                           25408500
C   IEF=13                                                25408600
C   GO TO 1000                                            25408700
C                                                         25408800
C   4 MF=4                                                 25408900
C   GO TO 3                                               25409000
C   2 IF(NREC.GT.NRECP) GO TO 4                          25409100
C   MF=3                                                  25409200
C   3 NAMES(2)=MF*10000+MT*10                             25409300
C   IF(NREC.EQ.14.OR.NREC.EQ.15) NAMES(2)=NAMES(2)+2    25409400
C   RTYPE=TYPNA(ITY)                                     25409500
C   DUM3(4)=NAMES(1)                                     25409600
C   IEF=0                                                 25409700
1000 RETURN                                              25409800
C                                                         25409900
C   END                                                    25410000
C   SUBROUTINE RKNDP(NAMES,NZNAM,MF,EMIN,EMAX,X,Y,NPOINT,IEF) 25410100
C                                                         25410200
C   THIS ROUTINE FILLS THE ARRAYS X AND Y WITH KNDF-DATA FROM THE 25410300
C   REQUIRED ENERGY RANGE                                25410400
C                                                         25410500
C   DIMENSION X(150),Y(150),NARG(3),NAMES(5),XX(150),YY(150) 25410600
C   DATA MAXNUM,EPS/150,0.01/                            25410700
C                                                         25410800
C   IEF=0                                                 25410900
C   NOUT=6                                                25411000
C   IF(MF.EQ.4) GO TO 400                                 25411100
C   IF(MF.EQ.3) GO TO 300                                 25411200
C   GO TO 1                                               25411300
400 CALL SELK (NARG,NAMES,EMIN,EMAX,X,Y,MAXNUM,NPOINT,IEF) 25411400
C   GO TO 8                                               25411500
300 NPOI1=0                                              25411600
C   NARG(1)=NZNAM                                         25411700

```

```
MEMBER NAME KNDF
C
CALL RETXS(NARG,NAMES,EMIN,EMAX,XX,YY,NUMX,MAXNUM,NR) 25411800
C
I1=2 25411900
21 IF(NR.EQ.0) GO TO 11 25412000
IF(NR.EQ.1) GO TO 12 25412100
IF(NR.EQ.2) GO TO 13 25412200
IF(NR.EQ.3) GO TO 1 25412300
IF(NR.EQ.4) GO TO 1 25412400
IF(NR.EQ.5) GO TO 1 25412500
GO TO 16 25412600
C
11 I1=2 25412700
IF(NPOI1.NE.0) GO TO 13 25412800
18 DO 10 I=1,NUMX 25412900
X(I)=XX(I) 25413000
10 Y(I)=YY(I) 25413100
NPCINT=NUMX 25413200
GO TO 17 25413300
C
19 CALL REPXS(NARG,NAMES,EMIN,EMAX,XX,YY,NUMX,MAXNUM,NR) 25413400
GO TO 21 25413500
C
16 WRITE (NOUT,102) NR 25413600
102 FORMAT(1X,'THE RETURN-CODE FROM RETXS = ',I10) 25413700
GO TO 1 25413800
C
12 IF(NPOI1.EQ.0) GO TO 18 25413900
13 CALL SKIP01(XX,YY,NUMX,X,Y,NPOI1,MAXNUM,EFS) 25414000
IF(NPOI1.LT.MAXNUM.AND.NR.EQ.2) GO TO 19 25414100
NPCINT=NPOI1 25414200
IF(NPOINT.LT.2) GO TO 1 25414300
C
INTERPOLATE CROSS SECTION TO LOWER ENERGY BOUNDARY 25414400
C
17 IF(EMIN.GT.0.999999*X(1).AND.EMIN.LT.1.000001*X(1)) X(1)=EMIN 25414500
IF(EMIN.LE.X(1)) GO TO 7 25414600
CALL TRPL8(X(1),X(2),EMIN,I1,Y(1),Y(2),Z) 25414700
X(1)=EMIN 25414800
Y(1)=Z 25414900
7 IF(EMAX.EQ.0.0) GO TO 8 25415000
IF(EMAX.GE.X(NPOINT))GO TO 8 25415100
C
INTERPOLATE CROSS SECTION TO UPPER ENERGY BOUNDARY 25415200
C
CALL TRPL8 (X(NPCINT-1),X(NPCINT),EMAX,I1,Y(NPOINT-1),Y(NPOINT),Z 25415300
1) 25415400
X(NPOINT)=EMAX 25415500
Y(NPOINT)=Z 25415600
IEF=0 25415700
GO TO 8 25415800
1 IEF=13 25415900
8 RETURN 25416000
END 25416100
SUBROUTINE SELK (NARG,NAMEN,EM11,EM11,XX,YY,MAXNUM,NUM1,IEF) 25416200
C
NEUTRON ELASTIC SCATTERING ANGULAR DISTRIBUTIONS FROM KNDF 25416300
C
END 25416400
C
END 25416500
C
END 25416600
C
END 25416700
C
END 25416800
C
END 25416900
C
END 25417000
C
END 25417100
C
END 25417200
C
END 25417300
C
END 25417400
C
END 25417500
```

```
MEMBER NAME KNDF
DIMENSION XX(150),YY(150),NARG(3),NAMES(5),X1(101),Y1(101),X2(101) 25417600
1,Y2(101),XN(2),YN(2),F(10),NAMEN(1) 25417700
INTEGER*4 SGN/30020/,SGNC/40022/ 25417800
EQUIVALENCE(NAMES(3),XNAM) 25417900
C 25418000
IEF=0 25418100
NOUT=6 25418200
MXN=101 25418300
NAMES(1)=NAMEN(1) 25418400
NAMES(2)=SGN 25418500
NARG(1)=2 25418600
EMIN=EM11 25418700
C 25418800
CALL RETXS(NARG,NAMES,EMIN,EMIN,XN,YN,NUMX,2,NR) 25418900
IF(NR.GT.2) GO TO 13 25419000
C 25419100
IF(XN(1).EQ.EMIN) GO TO 11 25419200
IF(XN(2).EQ.EMIN) GO TO 12 25419300
CALL TRPL3(XN(1),XN(2),EMIN,2,YN(1),YN(2),2) 25419400
WSGN=Z 25419500
GO TO 14 25419600
C 25419700
11 WSGN=YN(1) 25419800
GO TO 14 25419900
C 25420000
12 WSGN=YN(2) 25420100
14 ESGN=EMIN 25420200
WRITE (NOUT,105) ESGN,WSGN 25420300
C 25420400
IGO=0 25420500
NAMES(1)=NAMEN(1) 25420600
NAMES(2)=SGNC 25420700
XNAM=0. 25420800
NARG(1)=3 25420900
7 CALL IDFLOC (NERR,NARG,NAMES,F) 25421000
IF(XNAM.EQ.0.) GO TO 19 25421100
C 25421300
NAMES(1)=NAMEN(1) 25421400
NAMES(2)=SGNC 25421500
NARG(1)=3 25421600
EMAX=1. 25421700
EMIN=-1. 25421800
E1=XNAM 25421900
XNAM1=XNAM 25422000
CALL RETXS(NARG,NAMES,EMIN,EMAX,XX,YY,NUMX,MAXNUM,NR) 25422100
IF(NR.GT.2) GO TO 23 25422200
105 FORMAT(1X,10E13.5) 25422400
LEVEL=1 25422500
NUM1=NUMX 25422600
XNAM2=XNAM 25422700
E2=XNAM 25422800
C 25422900
15 IF(E1.LT.ESGN) GO TO 16 25423000
IF(E1.GT.ESGN) GO TO 19 25423100
LEVEL=2 25423200
IF(IGO.NE.0) GO TO 18 25423300
IEF=0 25423400
GO TO 1000 25423500
```

```
MEMBER NAME KNDF
C
13 WRITE (NDUT,101) NAMES(1),EMIN 25423600
101 FORMAT(1X,' FOR ',110,' SGN, EMIN= ',E16.8,' NO DATA FOUND ON KNDF 25423700
1') 25423800
GC TO 19 25423900
23 WRITE (NDUT,102) NAMES(1),E1 25424000
102 FORMAT(1X,' FOR ',110,' SGNC LEVEL ENERGY= ',E16.3,' NO DATA FOUND 25424100
IGN KNDF') 25424200
19 IEF=13 25424300
GC TO 1000 25424400
C
16 IF(E2.LT.ESGN) GO TO 17 25424500
IF(E2.GT.ESGN) GO TO 18 25424600
17 CALL IDFNXT(NERR,NARG,NAMES,F) 25424700
IF(NERR.NE.0) GO TO 17 25424800
IF(XNAM2.EQ.XNAM) GO TO 18 25424900
XNAM1=XNAM2 25425000
E1=E2 25425100
XNAM2=XNAM 25425200
E2=XNAM 25425300
IGC=1 25425400
GO TO 15 25425500
C
18 XNAM=XNAM1 25425600
NAMES(1)=NAMEN(1) 25425700
NAMES(2)=SGNC 25425800
CALL RETXS(NARG,NAMES,EMIN,EMAX,X1,Y1,NUMX,MXN,NR) 25425900
IF(NR.GT.2) GO TO 23 25426000
NUM2=NUMX 25426100
NAMES(1)=NAMEN(1) 25426200
NAMES(2)=SGNC 25426300
XNAM=XNAM2 25426400
IF(LEVEL.EQ.2) GO TO 24 25426500
CALL RETXS(NARG,NAMES,EMIN,EMAX,X2,Y2,NUMX,MXN,NR) 25426600
IF(NR.GT.2) GO TO 21 25426700
LEVEL=3 25426800
NUM3=NUMX 25426900
C
24 NUM1=NUM2-1 25427000
DO 10 I=1,NUM1 25427100
DO 20 J=1,NUM1 25427200
IF(X1(J).LT.XX(I).AND.X1(J+1).GT.XX(I)) GO TO 25 25427300
IF(X1(J).EQ.XX(I)) GO TO 32 25427400
20 CCNTINUE 25427500
IF(X1(NUM2).LE.XX(I)) GO TO 21 25427600
GC TO 29 25427700
32 YY(I)=Y1(J) 25427800
GO TO 10 25427900
C
21 J=NUM1 25428000
25 CALL TRPL8 (X1(J),X1(J+1),XX(I),2,Y1(J),Y1(J+1),YY(I)) 25428100
10 CCNTINUE 25428200
IF(LEVEL.EQ.2) GO TO 31 25428300
C
26 NUM1=NUM3-1 25428400
DO 30 I=1,NUM1 25428500
DO 40 J=1,NUM1 25428600
IF (X2(J).LT.XX(I).AND.X2(J+1).GT.XX(I)) GO TO 27 25428700
25428800
25428900
25429000
25429100
25429200
25429300
25429400
25429500
```

MEMBER NAME	KNDF	
	IF(Z(1).LE.EMIN) GOTO 5	25435400
	IF(Z(1).GE.EMAX) GOTO 36	25435500
11	I=I+1	25435600
	X(I)=W(1)	25435700
	Y(I)=W(2)	25435800
12	I=I+1	25435900
	X(I)=Z(1)	25436000
	Y(I)=Z(2)	25436100
	GOTO NST,(20,200)	25436200
20	CALL IDFNXT(NERR,NARG,NAMES,Z)	25436300
	IF(NERR.EQ.0) GOTO 22	25436400
	IF(Z(1).GE.EMAX) GOTO 24	25436500
	IF(I.EQ.MAXNUM) GOTO 34	25436600
21	I=I+1	25436700
	X(I)=Z(1)	25436800
	Y(I)=Z(2)	25436900
	GOTO 20	25437000
C		25437100
	ENTRY REPRS(NARG,NAMES,EMIN,EMAX,X,Y,NUMX,MAXNUM,NR)	25437200
	I=0	25437300
	NAMZ=NARG(1)	25437400
	IF(NAMZ.LE.2) GOTO 21	25437500
	DO 19 J=3,NAMZ	25437600
19	NAMSV(J-2)=NAMES(J)	25437700
	GOTO 21	25437800
C		25437900
22	IF(I.LT.1) GOTO 23	25438000
	NR=1	25438100
	GOTO 198	25438200
23	NR=5	25438300
	I=1	25438400
	X(I)=Z(1)	25438500
	Y(I)=Z(2)	25438600
	GOTO 198	25438700
C		25438800
24	IF(I.EQ.MAXNUM) GOTO 34	25438900
	I=I+1	25439000
	X(I)=Z(1)	25439100
	Y(I)=Z(2)	25439200
	NR=0	25439300
	GOTO 200	25439400
26	I=1	25439500
	X(I)=W(1)	25439600
	Y(I)=W(2)	25439700
	NR=5	25439800
	GOTO 198	25439900
C		25440000
30	NR=3	25440100
	GOTO 200	25440200
C		25440300
32	NR=4	25440400
	I=1	25440500
	X(I)=Z(1)	25440600
	Y(I)=Z(2)	25440700
	GOTO 200	25440800
C		25440900
34	NR=2	25441000
	GOTO 200	25441100



```
MEMBER NAME KNDF
36 ASSIGN 200 TO NST 25441200
NR=0 25441300
GOTO 7 25441400
38 ASSIGN 200 TO NST 25441500
NR=0 25441600
GOTO 11 25441700
198 IF(NAMZ.LE.2) GOTO 200 25441800
DC 199 J=3,NAMZ 25441900
199 NAMES(J)=NAMSV(J-2) 25442000
200 NUMX=I 25442100
RETURN 25442200
END 25442300
SUBROUTINE SKIPDI (XX,YY,NUMX,X,Y,NPCINT,LENX,EPS) 25442400
DIMENSION XX(1),YY(1),X(1),Y(1) 25442500
C 25442600
C SKIPI SKIPS POINTS NOT NECESSARY TO REPRESENT CURVE TO AN 25442700
C ACCURACY OF EPS. 25442800
C 25442900
J=1 25443000
I=0 25443100
IF(NPOINT.EQ.0) GOTO 10 25443200
I=NPOINT 25443300
IF(NUMX.GT.1) GOTO 11 25443400
10 I=I+1 25443500
IF(I.GT.LENX) GOTO 99 25443600
X(I)=XX(J) 25443700
Y(I)=YY(J) 25443800
IF(J.GE.NUMX) GOTO 100 25443900
11 IF(J+1.GE.NUMX) GOTO 98 25444000
JA=J+1 25444100
JB=J+2 25444200
DO 30 JK=JB,NUMX 25444300
XK=(YY(JK)-Y(I))/(XX(JK)-X(I)) 25444400
JC=JK-1 25444500
IS=0 25444600
DC 20 JJ=JA,JC 25444700
YI=XK*(XX(JJ)-X(I))+Y(I) 25444800
IF(ABS(YI-YY(JJ)).LT.EPS*YY(JJ)) GOTO 18 25444900
GOTO 22 25445000
18 IF(ABS(YI-YY(JJ)).GT.1.E-5*YY(JJ)) GOTO 20 25445100
IS=IS+1 25445200
20 CONTINUE 25445300
JA=JA+IS 25445400
GOTO 30 25445500
22 J=JC 25445600
GOTO 10 25445700
30 CONTINUE 25445800
J=NUMX 25445900
GOTO 10 25446000
98 J=J+1 25446100
GOTO 10 25446200
99 I=I-1 25446300
100 NPCINT=I 25446400
RETURN 25446500
END 25446600
SUBROUTINE KEDAK (IER) 25446700
C 25446800
C RETRIEVES KEDAK DATA FOR SCORE 25446900
```

MEMBER NAME	KNDF	
C		25447000
	CCOMDN /LDFMT/ IDUM(4), MAT(70), IMA(70), DUM(105), NMAT	25447100
	COMMON /BLK1/ ABCD(5,4), ICASE, ISHIFT, E(5), REAC(10), BGTIL(18), U(10)	25447200
	COMMON /BLK2/ NULL, P(1500), I1, DU(1004)	25447300
	CCOMDN /BLK3/ NR, BCDR(5,30), TCHAR(30), ISEL(30)	25447400
	COMMON /BLK4/ DUM1(4), QX, ID(5), IUM3(4), RTYPE	25447500
	CCOMDN /BLK5/ LI(4), XY3(6), XY4(4)	25447600
	COMMON /BLK6/ NUL1, XYS(90), NUL2	25447700
	CCOMDN /BLK7/ NUL3, XY(300), NUL4, NUL5, YX(300), NUL6	25447800
	CCOMDN /BLK8/ NPCINT, X(150), Y(150), NKNDP, ZLIM(4)	25447900
	COMMON /BLK13/ NEAD, EAD(15), GAMD(15), NPE(16)	25448000
C		25448100
	REAL*8 MAT, MATER, TYPES(70), RTYP, NAM1(2)	25448200
	DIMENSION ITYP(70), IR(70), XX(3), NAMES(5), IAT(70), NDAT(70), REA(10)	25448300
	DIMENSION XMAT(5), DUMC(2), BGTIL2(18), XNAM1(200)	25448400
	DATA IR/70*2/, BLANK/' '/	25448500
	EQUIVALENCE (REA(1), RTYP), (XMAT(1), MATER), (NAMES(3), XNAM)	25448600
C		25448700
	NULL = 0	25448800
	NUL1=0	25448900
	NUL2=0	25449000
	NUL3=0	25449100
	NUL4=0	25449200
	NUL5=0	25449300
	NUL6=0	25449400
	NZ=2	25449500
	NPCINT=0	25449600
	NCUT=6	25449700
	IDAT=IER	25449800
	IER=1	25449900
	ID(1)=1	25450000
	NR=0	25450100
	DC 120 I=1,10	25450200
	REA(I)=BLANK	25450300
120	U(1)=BLANK	25450400
	DC 130 I=1,18	25450500
	BGTIL2(I)=BLANK	25450600
130	BGTIL(I)=BLANK	25450700
	DC 10 J=1,30	25450800
	DC 10 I=1,5	25450900
	XMAT(I)=BLANK	25451000
10	BCDR(I,J)=BLANK	25451100
C		25451200
C	USER IDENTIFICATION	25451300
C		25451400
	CALL IDENT(IER)	25451500
	IF(IER.NE.1) GO TO 1000	25451600
1	IMAT=0	25451700
C		25451800
C	SELECT ISOTPE	25451900
C		25452000
	CALL CHOSMA(MAT, IAT, NMAT, IMAT, IDAT)	25452100
	IF(IMAT.LT.1.OR.IMAT.GT.70) GO TO 1	25452200
	NAMES(1)=IMA(IMAT)	25452300
	IUM3(4)=NAMES(1)	25452400
	MATER=MAT(IMAT)	25452500
	IATER=IMA(IMAT)	25452600
	ID(3) =IATER/10000	25452700

```
MEMBER NAME  KNDF
WRITE (NOUT,101) MATER
101 FFORMAT (1H0,5X,A3)
C
C   FILL LIST WITH REACTION TYPES FOR SELECTED ISOTOPE
C
CALL FILLTP (NER,MATER,TYPES,ITYP,NTYP)
IF(NER.NE.1) GO TO 33
IF(NTYP.EQ.0) GO TO 1
XX(3)=0.
C
C   CHOOSE REACTION TYPE
C
CALL CHOSRE (IR,NTYP,TYPES,RTYP,NREC,XX,ICAT)
IF(NREC.LT.1.OR.NREC.GT.32) GO TO 1
RTYPE=RTYP
ITYPE=ITYP
NAMES(2)=ITYPE
MM=ITYPE/10000
IF(MM.EQ.3) GO TO 11
IF(ITYPE.EQ.40022) GO TO 12
GO TO 1000
C
12 ID(1)=14
11 EMIN=XX(1)*1.0E+6
EMAX=XX(2)*1.0E+6
XNAM=XX(3)*1.0E+6
DUM1(1)=EMIN
DUM1(2)=EMAX
DUM1(3)=XNAM
C
IF(ITYPE.EQ.40022.OR.ITYPE.EQ.30050) GO TO 13
GO TO 200
C
13 NAM1(1)=MATER
NAM1(2)=RTYP
MAX=200
NZ=3
C
CALL LDFNAM (NER,NAM1,XNAM1,N1,N2,MAX,2)
C
IF(NER.NE.1) GO TO 33
C
CALL ANERG(NAM1,N2,XNAM1,NNAM,IER)
C
IF (IER.NE.0) GO TO 55
IER=1
IF (ITYPE.EQ.40022) GO TO 110
QX=XNAM1(NNAM)*1.0E-6
XNAM=XNAM1(NNAM)
DUM1(3)=XNAM
GO TO 200
110 EMIN=XNAM1(NNAM)
DUM1(1)=EMIN
DUM1(3)=EMIN
NEAD=1
EAD(1)=EMIN
C
C   RETRIEVE CROSS SECTION FROM KEDAK
```

```

MEMBER NAME   KNDF
C
200 CALL RKNDP (NAMES,NZ,MM,EMIN,EMAX,X,Y,NPCINT,IEF)
      IF (IEF.NE.0) GO TO 55
      IF (MM.EQ.4) GO TO 2
C
      DC 20 I=1,NPCINT
20 X(I)=X(I)*1.E-6
      IF (X(2).LT.1.000001*X(1).AND.X(2).GT.C.999999*X(1)) GO TO 5
      IF (Y(1).NE.0.0) GO TO 4
C
      DC 30 I=2,NPCINT
      I1=I-1
      X(I1)=X(I)
30 Y(I1)=Y(I)
      NPCINT=NPCINT-1
      CONTINUE
      NKNDP=NPCINT
C
      CALL CALIMA(NPCINT,X,Y,ZLIM)
C
      DC 70 I=1,4
      XY4(I)=ZLIM(I)
70 DL(I)=ZLIM(I)
C
      DC 40 I=1,10
      BGTIL(I+5)=REA(I)
40 REAC(I)=REA(I)
      DC 50 I=1,5
      BGTIL(I)=XMAT(I)
50 E(I)=XMAT(I)
      ICASE=ICASE+1
      CALL IDACON (DUMC,ICASE,2,1,1,1,IE)
      CNUM=DUMC(2)
      ICNUM=17
      BGTIL(17)=CNUM
      DC 90 I=1,18
50 BGTIL2(I)=BGTIL(I)
      CALL CENTER(BGTIL2,18,BGTIL,ISHIFT)
      ISHIFT=ISHIFT+4*(ICNUM-1)+1
C
1000 RETURN
      DC 50 CALL ERRORD(IEF,IER)
      GO TO 1000
33 IEF=13
      GO TO 55
      END
C
      SUBROUTINE FILLTP(NR,MAT,TYP,ITYP,NT)
C
      REAL*8 TYP(32)/'SGT','SGN','SGX','SGI','SGIZ','SGIZ','S52N','SG32
IN','SGIA','SGI3A','SG2NA','SG3NA','SGIP','SGNI','SGA','SGF','SGG',
2'SGP','SGD','SGH3','SGHE3','SGALP','SG2FE','SGTR','MUEL','ETA','AL
3PHA','NUC','NUEP','CHIF','CHIFD','SGNC',TYP(70),MAT,NAMES(3)
      INTEGER NP(70),MTYP/70/,MTYP5/32/,NARG(3)
      DIMENSION ITYP(70)
      COMMON/LDFTT/MC(3)
C
      FILL TABLE OF SINGLE VALUED,ENERGYDEPENDENT TYPES.

```

25458600  
25458700  
25458800  
25458900  
25459000  
25459100  
25459200  
25459300  
25459400  
25459500  
25459600  
25459700  
25459800  
25459900  
25460000  
25460100  
25460200  
25460300  
25460400  
25460500  
25460600  
25460700  
25460800  
25460900  
25461000  
25461100  
25461200  
25461300  
25461400  
25461500  
25461600  
25461700  
25461800  
25461900  
25462000  
25462100  
25462200  
25462300  
25462400  
25462500  
25462600  
25462700  
25462800  
25462900  
25463000  
25463100  
25463200  
25463300  
25463400  
25463500  
25463600  
25463700  
25463800  
25463900  
25464000  
25464100  
25464200  
25464300

MEMBER NAME	KNDF	
	KCUT=6	25464400
	CALL LDFITN(NR,MAT,TYP,NP,NT,MTYP,2)	25464500
	IF(NR.EQ.0.AND.NT.EQ.0) GOTO 100	25464600
	IMAT=MO(3)	25464700
	CALL LDFITN(NR,IMAT,ITYP,NP,NT,MTYP,1)	25464800
	IF(NR.EQ.0.AND.NT.EQ.0) GOTO 99	25464900
	J=0	25465000
	DO 12 N=1,NT	25465100
	DC 10 K=1,MTYPS	25465200
	IF(TYP(N).NE.TYPS(K)) GOTO 10	25465300
	J=J+1	25465400
	TYP(J)=TYP(N)	25465500
	ITYP(J)=ITYP(N)	25465600
	NP(J)=NP(N)	25465700
	GOTO 12	25465800
10	CONTINUE	25465900
12	CONTINUE	25466000
	NT=J	25466100
	GOTO 100	25466200
99	WRITE(KOUT,600)	25466300
600	FORMAT(///' PROGRAMMING-ERROR IN LDFITN.CALL PROGRAMMER.')	25466400
	STOP	25466500
100	CONTINUE	25466600
	RETURN	25466700
	END	25466800
C		25466900
	SUBROUTINE ANERG (NAM1,N2,XNAM1,NNAM,IER)	25467000
C		25467100
C	PERMITS THE SELECTION OF THE LEVEL ENERGY FOR INELASTIC SCATTERING	25467200
C	OR THE INCIDENT ENERGY FOR ELASTIC SCATTERING	25467300
C		25467400
	REAL*8NAM1(2)	25467500
	DIMENSION XNAM1(1),BCD(3)	25467600
	DATA PAGE/'PAGE',Y1,X00,Y00,DY1/10.75,4.5,10.,.5/	25467700
C		25467800
	I=0	25467900
	NNAM=0	25468000
	IST=0	25468100
C		25468200
100	CALL INPICT	25468300
	CALL WRTXTM (2,16,NAM1, 3.,Y1)	25468400
	X3=X00	25468500
	Y3=Y00	25468600
	IL=15*IST+1	25468700
	IF(IL.LT.N2) GO TO 111	25468800
	IL=1	25468900
	IST=0	25469000
111	IU=IL+14	25469100
	IL=MINO(IU,N2)	25469200
	NN=IU-IL+1	25469300
C		25469400
	CALL SETKEY (7,1,12)	25469500
	DC 10 K=IL,IU	25469600
	ISZ=1	25469700
	IF (I.EQ.K) ISZ=2	25469800
	DNUM=XNAM1(K)	25469900
	CALL DACON (BCD,DNUM,3,1,1,6,1,IER)	25470000
	IF(IER.NE.0) GO TO 1000	25470100

MEMBER NAME	KNDF	
	CALL WRTXTM (1SZ,12,BCD,X3,Y3)	25470200
10	Y3=Y3-DY1	25470300
	CALL ENDKEY	25470400
C		25470500
	CALL SETKEY (1,1,4)	25470600
	CALL WRTXTM (2,4,PAGE,5.5,1.)	25470700
	CALL ENDKEY	25470800
C		25470900
	CALL STPICT	25471000
150	CALL EXINGR(IKEY,IVAL,ISEQ)	25471100
C		25471200
	GO TO (120,200,120,120,150,150),IKEY	25471300
200	IF (IVAL.EQ.2) GO TO 110	25471400
	IF (ISEQ.GT.NN)GO TO 150	25471500
	IF (I.LT.IL.OR.I.GT.IU) GO TO 105	25471600
	ICH=I-IL+1	25471700
	CALL CHGCHR (1,ICH,1)	25471800
105	CALL CHGCHR (1,ISEQ,2)	25471900
	I=IL+ISEQ-1	25472000
	GO TO 150	25472100
110	IST=IST+1	25472200
	GO TO 100	25472300
C		25472400
120	NNAM=I	25472500
1000	RETURN	25472600
	END	25472700

A3. Input Summary for Karlsruhe Version of SCØRE

Input from the IBM 2250 terminal.

Note: END key means key no. 1,

CANCEL key means key no. 2 of the programmed-function keyboard.

Figure 1:

- (1) Chose between options KEDAK and NEUDADA by the light pen.
- (2) Press END key.

Figure 2: User and case identification

- (1) Type in user identification, date and running number of problem (case identification) from keyboard.
- (2) Press END key.

Figure 3: List of isotope names

- (1) Selection of PAGE with light pen turns page
- (2) Select isotope with light pen
- (2) Press END key

Figure 4: List of reaction types

- (1) Select reaction type with light pen
- (2) Type in energy limits from keyboard
- (3) Press END key

Figure 5: List of energies for SGNC or SGIZ (only after selection of option KEDAK in Figure 1)

- (1) Selection of PAGE by light pen turns page
- (2) Select energy with light pen
- (3) press END key

Figure 6: Option lists MAIN 1/3, 2/3, 3/3

Options can be chosen by light pen or from programmed-function keyboard.

Notation used:

- k nn : programmed-function key no. nn,
- N : option valid for NEUDADA only,
- K : option valid for KEDAK only,
- N+K : option valid for both NEUDADA and KEDAK,

where NEUDADA and KEDAK refer to the choice made in Figure 1.

Option list MAIN 1/3

N+K	MAIN 1/3	k 11	Display option list MAIN 2/3.
N+K	RESTART	k 12	Jump back to start of program, Figure 1.
N+K	PRINT	k 13	Print data displayed on screen.
N	LETTERS(PØINTS)	k 14	Replace curve points by letters distinguishing references (or not).
N	HATS(NØ HATS)	k 15	Add error bars (or not).
N+K	EXPAND	k 16	Expand the graph portion of the display to full screen size, omitting caption, references and option lists. Pressing of the END key restores the unexpanded picture.
N	NØ CURVE	k 17	Delete curves of evaluated data.
N	LISTDATA	k 18	Tabulate energy and cross section values (plus references), 20 data points at one time on the screen. Choosing one of the numbers appearing in the right-



most column under the + (or -) sign by light pen causes forward (or backward) shifting of the list by so many energies. Light pen selection of (+) produces (-). Pressing of the END key terminates the LISTDATA option.

N+K            QUIT                            k 20        End terminal session .

Option list MAIN 2/3

N+K            MAIN 2/3                            k 11        Display option list MAIN 3/3 .

N+K            RESTART                            k 12        Jump back to start, Figure 1 .

N + K            REPLØT                            k 13        Change co-ordinate system .

N                AUTØ LIM                            k 14        Recalculate plot scale with modified data (after CØRRECT or DELETE).

N                SØRT                                k 15        Delete data points belonging to the references selected from the following Figure 7 .

N                DELETE                            k 17        Delete points indicated with light pen on expanded graph (curve points being represented by letters associated with references). Deletion of the last point can be canceled with the CANCEL key. Pressing of the END key terminates the DELETE option.

N                RESTØRE                            k 18        Restore all points previously deleted .

N                CØRRECT                            k 19        Earmark points indicated by light pen for correction. Erroneous assignment can be canceled with the CANCEL key. After selecting all points to be corrected one gets them listed upon pressing the END key in Figure 8.

N+K	QUIT	k 20	End terminal session .
<u>Option list MAIN 3/3</u>			
N+K	MAIN 3/3	k 11	Display option list MAIN 1/3 .
N+K	RESTART	k 12	Jump back to start, Figure 1 .
N	KEDAK	k 13	Display KEDAK curve together with experimental data (for option NEUDADA from Figure 1).
K	KEDAK	k 13	Jump back to Figure 2: Treat next case (for option KEDAK from Figure 1).
N	DATA FIT	k 14	Call module for data fit .
N	ENDF/B	k 15	Display ENDF/B curve together with experimental data .
N	RES CALC	k 16	Call module for resonance fits .
N	NØ CURVE	k 18	Delete curves or points. Requested option must be indicated by light pen and is then displayed in large size. Cancellation possible with CANCEL key. Option is terminated with END key.
N+K	QUIT	k 20	End terminal session .

Figure 7: List of references

This figure appears after the SORT option has been chosen from option list MAIN 2/3.

If more than 13 references exist for the selected data type and energy range one can turn the page by pressing the END key. Points from the references indicated by light pen are discarded, the selected references are displayed in large size. They can be restored by pointing to them once more. One can also type in

a reduced energy interval and get a corresponding new co-ordinate system by pointing to 'YES' with the light pen.

Figure 8: List of uncorrected and corrected data

This figure tabulating the first 15 points to be corrected appears after the CØRRECT option has been chosen from option list MAIN 2/3. One can indicate after LIGHT PEN DATA whether a value is to be multiplied by a factor or increased (decreased) by a given amount which then must be typed in. The values (energies or cross sections) to be corrected in this way, can then be indicated by light pen. Erroneous selection can be canceled with the CANCEL key. Alternatively one can set a cursor on the energy value of the incorrect point (right-hand pair of columns) and then type in the correct values. After completion of all corrections in the right-hand columns one must press the END key to get a display of the next 15 points earmarked for correction, etc., until eventually the corrected plot appears on the screen.

DATA FIT: Module for fitting non-resonance data

For this module there are two option lists,  
DATA 1/2 and 2/2.

Option list DATA 1/2

N	DATA 1/2	k 11	Display option list DATA 2/2.
N	LETTERS(POINTS)	k 12	Replace curve points by letters distinguishing references (or not).
N	HATS(NO HATS)	k 13	Add error bars (or not).
N	EXPAND	k 14	Expand the graph portion of the display to full screen size, cf. option list MAIN 1/3.
N	CØMPARE	k 15	Show curves fitted with different methods for comparison.
N	AUTO LIM	k 16	Recalculate plot scale with modified data (after CØRRECT or DELETE).

N	PRINT	k 17	Print fitted data.
N	LEG CØEF	k 18	Analyse differential angular distributions in terms of Legendre coefficients. END key is to be pressed after completion.
N	NØ CURVE	k 19	Delete curve or points. Requested option must be indicated by light pen and is then displayed in large size. Cancellation possible with CANCEL key. Option is terminated with END key.
N	MAIN ØPT	k 20	Jump back to main program SCØRE.

Option list DATA 2/2

N	DATA 2/2	k 11	Display option list DATA 1/2.
N	CHG NØDE	k 12	Nodes can be added, deleted or moved. On the expanded graph the curve points are shown with error bars together with a tracking pattern (3 concentric circles) which facilitates manipulation of nodes. Three programmed function keys can be activated:
	(ADD NØDE)	k 11	Add node with light pen. Each time the END key is pressed the center of the tracking pattern becomes a node of the fit curve, represented by an asterisk. Up to 30 node points can be generated. Pressing of the CANCEL key terminates this option.
	(REMOØVE)	k 12	Delete nodes with light pen. Pressing of the CANCEL key terminates this option.

	(MØVE)	k 13	Move a point: The point is indicated by light pen whereupon it is replaced by the tracking pattern which can be moved around with the light pen. Upon pressing of the END key the center is replaced by an asterisk which fixes a node. Pressing of the CANCEL key terminates this option.
N	COMPARE	k 15	Show curves fitted with different methods for comparison.
N	LINEAR	k 16	Interpolate linearly between nodes.
N	SPLINE	k 17	Interpolate with cubic spline between nodes.
N	AUTØ SPL	k 18	Interpolate with cubic spline determined by least-squares method.
N	NØ CURVE	k 19	Delete curve or points. Requested option must be indicated by light pen and is then displayed in large size. Cancellation is possible with CANCEL key. Option is terminated with END key.
N	MAIN OPT	k 20	Jump back to main program SCØRE.

RES CALC: Module for resonance fitting

For evaluation in the resonance region three resonance formalisms are available:

Breit-Wigner single-level formalism,  
Reich-Moore multi-level formalism,  
Adler-Adler multi-level formalism.

The fit module for the resonance region is called by selection of the option RES CALC in option list MAIN 3/3. A series of special figures is displayed:

Figure 1: Selection of resonance formalism

- (1) Select one of the three available formalisms by light pen.

Figure 2: User information

- (1) Any time a new resonance formalism is selected Doppler kernel, resolution function and sample composition must be respecified.
- (2) The selected options are displayed in large size.
- (3) After each selection the END key must be pressed.

Figure 3: Input of Doppler-broadening parameters

- (1) Select Doppler kernel with the light pen.
- (2) The chosen kernel is displayed in large size together with a list of parameters needed.
- (3) Type in the appropriate parameters.
- (4) Press END key.

Figure 4: Input of resolution-broadening parameters

- (1) Select resolution function with the light pen.
- (2) The chosen resolution function is displayed in large size together with a list of parameters needed.
- (3) Type in the appropriate parameters.
- (4) Press END key.

Figure 5: Specification of isotopic mixture

- (1) Select isotopes by light pen from a list of available isotopes. The chosen isotope is displayed in large size.
- (2) A list of needed parameters is shown. Type in the appropriate values.
- (3) By pointing with the light pen to DELETE the user can delete the selected isotope.
- (4) Press END key to terminate input for the isotope.
- (5) Renewed pressing of the END key calls Figure 6.

Figure 6:

- (1) Selection of YES by light pen calls Figure 7.
- (2) Selection of NØ by light pen terminates input of new isotopes.

Figure 7: Isotope identification

- (1) Type in identification parameters for the isotope.
- (2) Press END key. If resonance parameters for the isotope are already in the library Figure 5 reappears, otherwise Figure 8.

Figure 8: Resonance parameter input

- (1) Type in resonance parameters
- (2) After completing input for one resonance press END key. Figure 8 reappears for the next resonance. Up to 10 new resonances can be entered.
- (3) Press CANCEL key to terminate resonance parameter input.

Two options exist for modification of resonance parameters during the calculation without changing them in the disc library:

- ADD RES permits addition of new parameters
- MOD RES permits modification of parameters in core storage.

Both options start with Figure 9.

Figure 9: Selection of isotope for resonance modification

- (1) A list of isotopes in core storage is displayed allowing light-pen selection of those isotopes for which resonance parameters are to be added or modified.
- (2) Selection of ALL ZA means that changes are to be made for all isotopes.
- (3) Press END key after selection is complete.

If ADD RES was chosen, Figure 8 is displayed for each isotope selected for modification.

If RES MØD was chosen, Fig. 10 appears with a list of resonance energies.

Figure 10: Input of modified resonance parameters (RES MØD option)

- (1) Select resonance to be modified with the light pen. A picture for the chosen resonance similar to Figure 8 is displayed. Parameters must be typed in. After completion press END key and modify next resonance etc. After last modification press END key twice to terminate RES MØD option.

Figure 11:

This figure appears after selection of option E SHIFT from option list RES 3/3. Type in the letter identifying the reference for which the energy scale is to be modified, and also the coefficients  $c_1$  and  $c_2$  of the modification formula (see below). After pressing of the END key corrected energies  $E_c$  are calculated from the uncorrected energies E according to

$$E_c = \frac{c_1}{(1+c_2\sqrt{E})^2} E.$$

The E SHIFT option is terminated with the CANCEL key.

For module RES CALC there are three option lists, RES 1/3, 2/3 and 3/3.

Option list RES 1/3

MODEL	Display Figure 1.
CHANGES	Display Figure 2.
ADD RES	Display Figure 9 for isotope selection, then Figure 8 for addition of resonances.
MØD RES	Display Figure 9 for isotope selection, then Figure 10 for resonance modification.
LØAD RES	Load selected parameters into core memory.
CØMPARE	Display results of preceding and present calculation for comparison.
FIT	Apply iterative algorithm for resonance parameter adjustment thrice.
CALC	Calculate with selected resonance formalism and display resulting curve (no fit).
MAIN OPT	Terminate RES CALC option, jump back to main program SCØRE.



Option list RES 2/3

LETTERS (POINTS) Represent experimental points by the letters distinguishing references (by dots).

HATS(NO HATS) Add error bars (or not).

EXPAND Expand the graph portion of the display to full screen size, cf. option list MAIN 1/3.

STØRE RES Update direct-access library on disc (after completion of fit procedure).

AUTØ LIM Recalculate plot scale with modified data.

CØMPARE Display results of preceding and present calculation simultaneously for comparison.

FIT Apply iterative algorithm for resonance parameter adjustment thrice.

CALC Calculate with selected resonance formalism and display resulting curve (no fit).

MAIN ØPT Terminate option RES CALC, jump back to main program SCØRE.

Option list RES 3/3

P PØINTS Set "peak points" on expanded graph. Three points per resonance must be specified for the fit by means of the programmed-function keys:

- k 1 END key.
- k 2 CANCEL key.
- k 3 Delete all points.
- k 4 Delete previously selected points.
- k 11 Go 1 point further for selection of next point.
- k 12 Go 2 points further for selection of next point.
- k 13 Go 5 points further for selection of next point.
- k 14 Go 10 points further for selection of next point.
- k 15 Go to first point
- k 16 Add or modify "peak points" with light pen and tracking pattern. Select one point at the peak, press END key, select a point on the left-hand slope, press END key, select a point on the right-hand slope, press END key. Three such points are required per resonance.

V PØINTS	Additional "valley points" can be set on the expanded graph to improve the fit in the valleys between resonances. Use programmed-function keys in the same way as for P PØINTS.
E SHIFT	Display Figure 11 to enable energy scale for points from a given reference to be shifted.
NØ CURVE	Delete curves
CØMPARE	Display results of preceding and present calculation for comparison.
FIT	Apply iterative algorithm for parameter adjustment thrice.
CALC	Calculate with selected resonance formalism and display curve.
MAIN ØPT	Terminate RES CALC option, jump back to main program SCØRE.

## 5. PROGRAMS FOR ESTIMATION OF NUCLEAR-MODEL AND CROSS-SECTION PARAMETERS

A good evaluation of cross section data is not possible without utilization of nuclear models and nuclear reaction theory. Codes which serve this purpose are briefly described in the following sections. Since these codes are essentially independent of any evaluated-data format and most of them can be used as stand-alone programs they do not belong, strictly speaking, to the KEDAK software proper. To a certain degree, however, they reflect the methods employed in KEDAK evaluation work. Therefore, and because the programs as such may interest potential users elsewhere, program abstracts were included in the present compendium.

### 5.1 Multi-level shape analysis program for transmission data (FANAL)

Name of program:

FANAL2

Name and establishment of author:

F.H. Fröhner

INR

Kernforschungszentrum

D-7500 Karlsruhe, West Germany

Problem solved:

Resonance parameter extraction from several sets of high-resolution time-of-flight transmission data simultaneously.

Method of solution:

Multi-level shape analysis with two-channel Reich-Moore formalism. Picket-fence approximation for distant levels. Resolution broadening. Doppler broadening of narrow ("p-, d-.wave") levels, but not of s-wave levels (the code was written for structural materials). Iterative least-squares fit. Resulting fit is plotted after each step.

Restrictions:

Doppler broadening of s-wave levels is neglected (but not of "p-, d-...wave" levels, cf. above). Up to 5 transmission data sets with a total of not more than 5120 data points can be fitted simultaneously. The number of cross section parameters (channel radii, distant-level strength functions, resonance energies and partial widths) must not exceed 200 (corresponding to about 45 resonances) of which not more than 50 can be adjusted. The sample material must not contain more than 5 different nuclides.

Unusual features:

Card input is organized for maximum user convenience (no flags, no bookkeeping information), modular structure, no tricky programming.

Related and auxiliary programs:

Non-KFK users must replace the Karlsruhe plotter subroutine PLØTA following the instructions given in the program.

Programming language:

FØRTRAN IV

Computer:

IBM/360-65, /360-91, /370-168 etc.

Machine requirements:

260 kbytes of core storage, card reader, printer, plotter.

Typical running time:

1-20 minutes on IBM/370-168 under OS, depending on number of data points, number of parameters adjusted, quality of first guesses etc.

References

F.H. Fröhner, report KFK 2129 (1976)

5.2 Multi-level shape analysis program for capture data (FANAC)

Name of program:

FANAC

Name and establishment of author:

F.H. Fröhner

INR

Kernforschungszentrum

D-7500 Karlsruhe, West Germany

Problem solved:

Resonance parameter extraction from several sets of high-resolution time-of-flight capture data simultaneously.

Method solved:

Multi-level shape analysis with two-channel Reich-Moore formalism. Picket-fence approximation for distant levels. Resolution broadening. Doppler broadening of narrow ("p-, d-...wave") levels, but not of s-wave levels (the code was written for structural materials). Multiple-collision events are treated by Monte-Carlo simulation. Iterative least-squares fit. Resulting fit is plotted after each step.

Restrictions:

Doppler broadening of s-wave levels is neglected (but not of "p-, d-...wave" levels, cf. above). Up to 5 capture data sets with a total of not more than 512 data points can be fitted simultaneously. The number of cross section parameters (channel radii, distant-level strength functions, resonance energies and partial widths) must not exceed 200 (corresponding to about 45 resonances) of which not more than 20 can be adjusted. The sample material must not contain more than 5 different nuclides.

Unusual features:

Importance sampling and Russian roulette are used in the Monte Carlo calculations to increase efficiency and to avoid bias. Card input is organized for maximum user convenience (no flags, no bookkeeping information). Modular structure, no tricky programming.

Related and auxiliary programs:

Non-KFK users must replace the Karlsruhe plotter subroutine PLØTA following the instructions given in the program. Non-IBM users must replace the random-number generator RANDU by an equivalent subroutine.

Programming language:

FØRTRAN IV

Computer:

IBM/360-65,/360-91,/370-168 etc.

Machine requirements:

458 kbytes of core storage, card reader, printer, plotter.

Typical running time:

2-20 minutes on IBM/370-168 under OS, depending on number of data points, number of parameters adjusted, quality of first guesses etc.

References:

F.H. Fröhner, Proc. 4th Conf. on Nuclear Cross Sections and Technology, Washington D.C., 1975, Vol. 2, p. 929

5.3 Estimation of level-statistical parameters from individual resonance parameters (STARA)

Name of program:

STARA

Name and establishment of author:

F.H. Fröhner

INR

Kernforschungszentrum

D-7500 Karlsruhe, West Germany

Problem solved:

Estimation of level-statistical cross section parameters (average level spacing, average reduced neutron width, strength function, average radiation width) from individual resonance parameters (level energies, neutron widths, radiation widths), with due account of missing levels.

Method of solution:

Maximum-likelihood estimation of mean spacing and average reduced neutron width simultaneously, based on the "crystalline" regularity of level spacings (Dyson and Mehta) and the Porter-Thomas distribution of reduced neutron widths (per channel). The coupled system of 2 maximum-likelihood equations is solved by iteration. Uncertainties are also estimated. Results are plotted.

Restrictions:

The maximum number of resonances is 1000.

Unusual features:

No estimate of the observability threshold or its energy dependence is required, in contrast to the widely used Fuketa-Harvey method.

Related and auxiliary programs:

Non-KFK users must replace the Karlsruhe plotter subroutine PLØTA following the instructions given in the program.

Programming language:

FØRTRAN IV

Computer:

IBM/360-65,/360-91,/370-168 etc.

Machine requirements:

234 kbytes of core storage, card reader, printer, plotter.

Typical running time:

5-10 s per resonance series (on IBM/370-168 under OS)

References:

-

5.4 Estimation of level-statistical parameters from average cross sections  
(FITACS)

Name of program:

FITACS

Name and establishment of author:

F.H. Fröhner

INR

Kernforschungszentrum

D-7500 Karlsruhe, West Germany

Problem solved:

Estimation of level-statistical cross section parameters (average spacings, strength functions, average radiation widths) from average total and capture cross sections.

Method of solution:

Least-squares fit to average total, or average total and capture cross section data simultaneously, with due account of inelastic scattering. Data are weighted



according to uncertainty and/or point density. Average cross sections are calculated with Hauser-Feshbach theory including width fluctuations. Results are plotted.

Restrictions:

Energy dependence of strength function is neglected and only s-, p-, d-wave interactions are considered. Fission cannot be handled by present code version. This restricts applicability to energies up to 100 - 200 keV and excludes fitting of partial cross sections of fissile nuclei. Up to 200 total and up to 200 capture cross section values can be fitted. Only pure nuclides can be handled.

Unusual features:

As the code was written mainly for the interpretation of evaluated average cross section data weighting of points can be chosen according to point density (i.e. inversely proportional to energy interval "covered" by a given point) as an alternative to the more common weighting by inverse variance.

Related and auxiliary programs:

Non-KFK users must replace the Karlsruhe plotter subroutine PLØTA following the instructions given in the program.

Programming language:

FØRTRAN IV

Computer:

IBM/360-65, /360-91, /370-168 etc.

Machine requirements:

102 kbytes of core storage, card reader, printer, plotter.

Typical running times:

About 1 s per nuclide on IBM/370-168 under OS.

References:

-

5.5 INCH3 - a coupled-channels program for inelastic and elastic exchange reactions with automatic parameter optimization

INCH3 is a modified version of the program INCH written by A.D. Hill and D.J. Edens.

Name of program:

INCH3

Name and establishment of author:

B. Goel

INR

Kernforschungszentrum

D-7500 Karlsruhe, West Germany

Problem solved:

Estimation of interaction parameters of coupled-channels problem by fitting angle-integrated and differential elastic and inelastic cross section and polarization data at a given incident energy.

Method of solution:

Weighted least-squares fit. Numerical integration of the coupled-channels equations is performed with the full interaction out to a cut-off radius  $R_1$  for nuclear forces. Beyond  $R_1$  only long-range Coulomb terms (proportional to  $r^{-L-1}$ ) are used up to a radius  $R_2$  where the asymptotic series expansions of the Coulomb wave functions are valid and can be matched.

Restrictions:

Antisymmetrization between incoming particle and target nucleons is neglected and also excitation of the scattered particle. The interaction potential must be local and velocity-independent. No restrictions (other than practicability) limit the number of coupled channels, the spin of the scattered particle or the number of excited target-nuclear states.

Unusual features:

Standard potentials (Coulomb, Saxon-Woods, derivative Saxon-Woods, centrifugal potential) are available as options, Modifications are possible in subroutine NUCLIN following author's instructions /1/.

Related and auxiliary programs:

-

Programming language:

FØRTRAN IV

Computer:

ATLAS, CDC 6600, IBM/370-168

Machine requirements:

240 kbytes of core storage

Typical running time:

40 sec on IBM/370-168 for a parameter optimisation problem of 10 search cycles

References:

/1/ Unpublished description available from author.

5.6 DWBA: Program calculating inelastic-scattering and charge-exchange cross sections

Name of program:

DWBA

Name and establishment of author:

B. Goel

INR

Kernforschungszentrum

D-7500 Karlsruhe, West Germany

Problem solved:

Calculation of differential cross sections for inelastic scattering and charge exchange reactions such as those for (n,p) or ( $^3\text{He}$ ,t) reactions.

Method of solution:

Distorted-wave Born approximation. Numerical integration of Schrödinger equation. A combination of Saxon-Woods form and its derivative for real and imaginary nuclear potential is used. Calculated and experimental data can be plotted together (line printer plot).

Restrictions:

Only angular distribution is calculated.

Unusual features:

A formalism for the calculation of knock-on processes is also included.

Related and auxiliary programs:

-

Programming language:

FØRTRAN IV

Computer:

ATLAS, IBM/370-168

Machine requirements:

128 kbytes of core storage

Typical running time:

about 1 min per problem or angular distribution on IBM/370-168

Reference:

B. Goel, Thesis, University of Freiburg i.Br., 1970

5.7 SECDIST: Program for calculation of energy distributions of secondary neutrons produced by inelastic scattering

Name of program:

SECDIST

Names and establishment of authors:

I. Broeders, C. Broeders

INR

Kernforschungszentrum

D-7500 Karlsruhe, West Germany

Problem solved:

Calculation of angle-integrated energy spectra of secondary neutrons produced by inelastic scattering with account of equilibrium and pre-equilibrium processes /1/, /2/. Comparison with experimental data.

Method of solution:

Equilibrium processes are calculated using a special version of the Hauser-Feshbach expression for continuous channels, which is obtained by angle integration of the corresponding expression used in the computer code HELENE /3/.

Preequilibrium reactions are calculated using Blann's theory /4/ and modified version of Blann's code /5/. Absorption cross sections and transmission coefficients needed for the equilibrium and preequilibrium calculations are obtained by the nonlocal optical model developed by Perey and Buck /6/ and a modified version of their code /7/.

Experimental data for energy distributions of secondary neutrons are often given as functions of the scattering angle. In order to obtain quantities that are comparable with the computer results the angular distributions are fitted and extrapolated by squares of spherical Bessel functions or by Legendre polynomials and integrated.

#### Restrictions:

Only elastic and inelastic neutron scattering channels are considered in the Hauser-Feshbach formula.

#### Unusual features:

Absorption cross sections and transmission coefficients which are needed for the Hauser Feshbach and for the preequilibrium calculations are obtained by the nonlocal-optical-model code /6,7/ and transferred within the system by densely stored tables. The transfer of two-dimensional tables (cross sections versus energy or energy distributions of secondary neutrons versus energy) is performed according to PLOTEASY /8/ formats. In this way automatic plotting of these quantities is made possible. In order to minimize user input the NAMELIST input option is used by all modules.

#### Related and auxiliary programs:

SECDIST consists of the subprograms OPTMØD, PREEQ, SIGMIX and FITMØD and utilizes the plot package PLOTEASY /8/.

#### Programming language:

IBM FØRTRAN IV

#### Computer:

IMB/360 and /370 series (e.g. IBM 370 model 168). Plotting devices: IBM-1627, CALCOMP-Plotter, STATOS-Plotter, SYNETICS-Plotter, CRT-display at Karlsruhe. The program is executed under ØS 370.

Machine requirements:

One disk is needed for transferring data. Without overlay all modules can be run with 122 k core storage with one exception: OPTMOD needs 200 k.

Typical running time:

As example the CPU times on a IBM 370/168 are given for the calculations for 14 MeV neutrons on Fe<sup>56</sup> /1,2/

1. Preparation of the optical model data with the code OPTMOD ≈ 5 min 30 sec  
CPU-Time
2. Preequilibrium calculation with PREEQ 2 sec.
3. Equilibrium and composite data with SIGMIX 5 sec.
4. Fitting of experimental data 5 - 13 sec.

References:

- /1/ H. Jahn, Contributed paper to the IAEA consultants meeting on the use of nuclear theory in neutron nuclear data evaluation, Trieste 8-12 December 75.
- /2/ H. Borgwaldt, C.H.M. Broeders, I. Broeders, H. Jahn, M. Lalović, D. Rusch, Reaktortagung Nürnberg, 1975.
- /3/ S.K. Penny, "HELENE", ORNL-TM-2590 (1969)
- /4/ M. Blann, Phys, Rev. Lett. 28 (1972), 757 and Nucl. Phys. A213, (1973), 570
- /5/ M. Blann, COO\_3494-14
- /6/ F.G. Perey, B. Buck, Nucl. Phys. 32, (1962), 353
- /7/ F.G. Perey, ORNL-3429
- /8/ C.H.M. Broeders, PLOTEASY, (1976) unpublished