

## ADAS Subroutine xxdata\_37

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subroutine xxdata_37( iunit ,
&                      nemax , ntmax ,
&                      title , icateg , nenerg , nblock ,
&                      nform1 , param1 , nform2 , param2 ,
&                      ea     , fa     , teff   , mode   ,
&                      median , filnam , filout , calgeb ,
&                      ealgeb
&                      )
C-----
C
C **** fortran77 subroutine: xxdata_37 ****
C
C purpose: To fetch data from an adf37 data set and detect its main
C           characteristics.
C
C calling program: various
C
C input : (i*4) iunit      = unit to which input file is allocated
C input : (i*4) nemax      = max no of energy points that can be read in
C input : (i*4) ntmax      = max no of effective temps that can be read in
C
C output: (c*80) title      = header for file
C output: (i*4) icateg     = category of file
C                           1 => superposition
C                           2 => numerical
C output: (i*4) nenerg     = type 1 => number of distribution families
C                           type 2 => number of energy points
C output: (i*4) nblock      = type 1 => number of members in output family
C                           type 2 => number of effective temperatures
C output: (i*4) nform1     = type of threshold behaviour
C                           1 => cutoff
C                           2 => energy^param1
C output: (r*8) param1     = parameter of threshold form
C output: (i*4) nform2     = type of high-energy behaviour
C                           1 => cutoff
C                           2 => energy^-param2(1)
C                           3 => exp(-param2(1)*energy)
C                           4 => exp(-param2(1)*energy^param2(2))
C output: (r*8) param2()   = parameter of high-energy form
C output: (r*8) ea(,)      = energy points of tabulation
C output: (r*8) fa(,)      = distribution function tabulation
C output: (r*8) teff()     = effective temperature (eV)
C output: (r*8) mode()     = most probable energy (eV)
C output: (r*8) median()   = median energy (eV)
C output: (c*120)filnam() = file names of input families
C output: (c*120)filout    = file name of output family
C output: (c*25) calgeb(,) = distribution function algebra
C output: (c*25) ealgeb()   = energy parameter algebra
C
C local : (i*4) ieunit     = energy units of distribution function
C                           1 => kelvin
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C                                2 => eV
C local : (i*4) i      = general use
C local : (i*4) j      = general use
C local : (i*4) med_index= energy index of median
C local : (i*4) mode_index() = energy index of mode
C local : (i*4) dummy    = general use
C local : (i*4) ie      = general use
C local : (i*4) iblock   = general use
C local : (r*8) sum     = average energy contribution from i -> i+1
C local : (r*8) contrib()= average energy contribution from i -> i+1
C local : (r*8) de      = energy step from i -> i+1
C local : (i*4) ifirst   = position of first non-blank character in string
C local : (i*4) ilast    = position of last non-blank character in string
C local : (i*4) indx()   = index of algebra
C local : (c*80) blank   = dummy string
C
C routines:
C      routine      source      brief description
C -----
C      i4unit       ADAS        fetch unit number for output of messages
C      xxslen       ADAS        finds string length excluding leading and
C                               trailing blanks
C
C author: Paul Bryans, University of Strathclyde
C
C date: 04/02/04
C
C update:
C
C-----
CHARACTER*25          CALGEB (NTMAX, 4),          EALGEB (NTMAX)
CHARACTER*120         FILNAM (NEMAX),           FILOUT
CHARACTER*80          TITLE
INTEGER                ICATEG,          IUNIT,          NBLOCK,          NEMAX
INTEGER                NENERG,          NFORM1,          NFORM2,          NTMAX
REAL*8                 EA (NTMAX, NEMAX),        FA (NTMAX, NEMAX)
REAL*8                 MEDIAN (NTMAX),        MODE (NTMAX),        PARAM1
REAL*8                 PARAM2 (2),          TEFF (NTMAX)

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