

ADAS Subroutine bfdata

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C
SUBROUTINE BFDATA( IUNIT , NDLEV , NDTRN , NDMET ,
&                TITLED , IZ      , IZ0   , IZ1   , BWNO  ,
&                NPL   , BWNOA  , LBSETA, PRTWTA, CPRTA ,
&                IL    ,
&                IA    , CSTRGA , ISA    , ILA    , XJA    , WA    ,
&                CPLA  , NPLA   , IPLA   , ZPLA   ,
&                CIONP ,
&                NV    , SCEF    ,
&                ITRAN , MAXLEV ,
&                TCODE , I1A    , I2A    , AVAL   , SCOM
&                )
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C ***** FORTRAN77 SUBROUTINE: BFDATA *****

C

C PURPOSE: TO FETCH DATA FROM INPUT ADF04 DATA SET, INCLUDING
C MULTIPLE PARENTS ON FREE-ELECTRON AND CHARGE EXCHANGE
C ON RECOMBINATION, INCLUSION OF EXPLICIT CONTRIBUTIONS BY
C IONISATION.

C

C IMPROVEMENT OF AUTOMATIC IONISATION CALC. BY INCLUDING
C ASSIGNMENT OF FINAL STATE PARENT. KEEPS IONISATION
C POTENTIAL INFORMATION FROM LEVEL TERMINATOR LINE.

C

C CALLING PROGRAM: ADAS215

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C DATA:

C THE 'REAL' DATA IN THE FILE IS REPRESENTED IN AN ABBREVIATED
C FORM WHICH OMITTS THE "D" OR "E" EXPONENT SPECIFIER.
C e.g. 1.23D-06 or 1.23E-06 IS REPRESENTED AS 1.23-06
C 6.75D+07 or 6.75E+07 IS REPRESENTED AS 6.75+07

C

C THEREFORE THE FORM OF EACH 'REAL' NUMBER IN THE DATA SET IS:
C N.NN+NN or N.NN-NN

C

C THE UNITS USED IN THE DATA FILE ARE TAKEN AS FOLLOWS:

C

C IONISATION POTENTIAL: WAVE NUMBER (CM-1)
C INDEX LEVEL ENERGIES: WAVE NUMBER (CM-1)
C TEMPERATURES : KELVIN
C A-VALUES : SEC-1
C GAMMA-VALUES :
C RATE COEFFT. : CM3 SEC-1

C

C

C SUBROUTINE:

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C INPUT : (I*4) IUNIT = UNIT TO WHICH INPUT FILE IS ALLOCATED
C INPUT : (I*4) NDLEV = MAXIMUM NUMBER OF LEVELS THAT CAN BE READ
C INPUT : (I*4) NDTRN = MAX. NUMBER OF TRANSITIONS THAT CAN BE READ

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C INPUT : (I*4) NDMET = MAX. NUMBER OF METASTABLES ALLOWED
C
C OUTPUT: (C*3) TITLED = ELEMENT SYMBOL.
C OUTPUT: (I*4) IZ = RECOMBINED ION CHARGE READ
C OUTPUT: (I*4) IZ0 = NUCLEAR CHARGE READ
C OUTPUT: (I*4) IZ1 = RECOMBINING ION CHARGE READ
C (NOTE: IZ1 SHOULD EQUAL IZ+1)
C OUTPUT: (R*8) BWNO = IONISATION POTENTIAL (CM-1) OF LOWEST PARENT
C OUTPUT: (I*4) NPL = NUMBER OF PARENTS ON FIRST LINE AND USED
C IN LEVEL ASSIGNMENTS
C OUTPUT: (R*8) BWNOA() = IONISATION POTENTIAL (CM-1) OF PARENTS
C OUTPUT: (L*4) LBSETA() = .TRUE. - PARENT WEIGHT SET FOR BWNOA()
C .FALSE. - PARENT WEIGHT NOT SET FOR BWNOA()
C OUTPUT: (R*8) PRTWTA() = PARENT WEIGHT FOR BWNOA()
C OUTPUT: (C*9) CPRTA() = PARENT NAME IN BRACKETS
C
C OUTPUT: (I*4) IL = INPUT DATA FILE: NUMBER OF ENERGY LEVELS
C
C OUTPUT: (I*4) IA() = ENERGY LEVEL INDEX NUMBER
C OUTPUT: (C*18) CSTRGA() = NOMENCLATURE/CONFIGURATION FOR LEVEL 'IA()'
C OUTPUT: (I*4) ISA() = MULTIPLICITY FOR LEVEL 'IA()'
C NOTE: (ISA-1)/2 = QUANTUM NUMBER (S)
C OUTPUT: (I*4) ILA() = QUANTUM NUMBER (L) FOR LEVEL 'IA()'
C OUTPUT: (R*8) XJA() = QUANTUM NUMBER (J-VALUE) FOR LEVEL 'IA()'
C NOTE: (2*XJA)+1 = STATISTICAL WEIGHT
C OUTPUT: (R*8) WA() = ENERGY RELATIVE TO LEVEL 1 (CM-1) FOR LEVEL
C 'IA()'
C OUTPUT: (C*1) CPLA() = CHAR. SPECIFYING 1ST PARENT FOR LEVEL 'IA()'
C INTEGER - PARENT IN BWNOA() LIST
C 'BLANK' - PARENT BWNOA(1)
C 'X' - DO NOT ASSIGN A PARENT
C OUTPUT: (I*4) NPLA() = NO. OF PARENT/ZETA CONTRIBUTIONS TO IONIS.
C OF LEVEL
C OUTPUT: (I*4) IPLA(,) = PARENT INDEX FOR CONTRIBUTIONS TO IONIS.
C OF LEVEL
C 1ST DIMENSION: PARENT INDEX
C 2ND DIMENSION: LEVEL INDEX
C OUTPUT: (I*4) ZPLA(,) = EFF. ZETA PARAM. FOR CONTRIBUTIONS TO IONIS.
C OF LEVEL
C 1ST DIMENSION: PARENT INDEX
C 2ND DIMENSION: LEVEL INDEX
C
C OUTPUT: (C*92) CIONP = STRING CONTAINING LEVEL TERMINATOR LINE
C
C OUTPUT: (I*4) NV = INPUT DATA FILE: NUMBER OF GAMMA/TEMPERATURE
C PAIRS FOR A GIVEN TRANSITION.
C OUTPUT: (R*8) SCEF() = INPUT DATA FILE: ELECTRON TEMPERATURES (K)
C (INITIALLY JUST THE MANTISSA. SEE 'ITPOW()')
C (NOTE: TE=TP=TH IS ASSUMED)
C
C OUTPUT: (I*4) ITRAN = INPUT DATA FILE: NUMBER OF TRANSITIONS
C OUTPUT: (I*4) MAXLEV = HIGHEST INDEX LEVEL IN READ TRANSITIONS
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C OUTPUT: (C*1) TCODE() = TRANSITION: DATA TYPE POINTER:
C ' ' => Electron Impact Transition
C 'P' => Proton Impact Transition
C 'H' => Charge Exchange Recombination
C 'R' => Free Electron Recombination
C 'I' => Coll. ionisation from lower stage ion
C OUTPUT: (I*4) I1A() = TRANSITION:
C LOWER ENERGY LEVEL INDEX (CASE ' ' & 'P')
C SIGNED PARENT NDEX (CASE 'H','R' & 'I')
C OUTPUT: (I*4) I2A() = TRANSITION:
C UPPER ENERGY LEVEL INDEX (CASE ' ' & 'P')
C CAPTURING LEVEL INDEX (CASE 'H','R' & 'I')
C OUTPUT: (R*8) AVAL() = TRANSITION:
C A-VALUE (SEC-1) (CASE ' ')
C NEUTRAL BEAM ENERGY (CASE 'H')
C NOT USED (CASE 'P','R' & 'I')
C OUTPUT: (R*8) SCOM(,) = TRANSITION:
C GAMMA VALUES (CASE ' ' & 'P')
C RATE COEFFT.(CM3 SEC-1)(CASE 'H','R' & 'I')
C 1ST DIMENSION - TEMPERATURE 'SCEF()'
C 2ND DIMENSION - TRANSITION NUMBER
C
C (I*4) NVMAX = PARAMETER = MAX. NUMBER OF TEMPERATURES
C THAT CAN BE READ IN.
C (I*4) MTIED = PARAMETER = MUST BE GREATER THAN OR EQUAL TO
C THE MAX. NO. OF LEVELS.
C (R*8) DZERO = PARAMETER = MINIMUM VALUE FOR 'AVAL()' AND
C 'SCOM()' ARRAYS = 1.0D-30
C
C (I*4) I4UNIT = FUNCTION (SEE ROUTINE SELECTION BELOW)
C (I*4) IQS = X-SECT DATA FORMAT SELECTOR
C NOTE: IQS=3 ONLY ALLOWED IN THIS PROGRAM
C (I*4) IFAIL = FAILURE NUMBER FROM B9PARS AND B9PRS1
C (I*4) I = GENERAL USE.
C (I*4) IABT = RETURN CODE FROM 'R(FCTN' (0 => NO ERROR)
C OR FROM INTERROGATION OF 'C7'
C (I*4) J = GENERAL USE.
C (I*4) J1 = INPUT DATA FILE - SELECTED TRANSITION:
C LOWER ENERGY LEVEL INDEX (CASE ' ' & 'P')
C (I*4) J2 = INPUT DATA FILE - SELECTED TRANSITION:
C UPPER ENERGY LEVEL INDEX (CASE ' ' & 'P')
C CAPTURING LEVEL INDEX (CASE 'H' & 'R')
C (I*4) LENCST = BYTE LENGTH OF STRING CSTRGA()
C (I*4) ILINE = ENERGY LEVEL INDEX FOR CURRENT LINE
C (I*4) IRECL = RECORD LENGTH OF INPUT DATASET (<=128)
C (I*4) IAPOW = EXPONENT OF 'AVALM'
C (I*4) IGPOW() = EXPONENT OF 'GAMMA()'
C (I*4) ITPOW() = TEMPERATURES - EXPONENT
C NOTE: MANTISSA INITIALLY KEPT IN 'SCEF()'
C
C (R*4) ZF = SHOULD BE EQUIVALENT TO 'IZ1'
C
C (R*8) AVALM = INPUT DATA FILE - SELECTED TRANSITION:

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C          MANTISSA OF: ('IAPOW' => EXPONENT)
C          A-VALUE (SEC-1)          (CASE ' ')
C          NEUTRAL BEAM ENERGY     (CASE 'H')
C          NOT USED                   (CASE 'P','R' & 'I')
C      (R*8)  GAMMA() = INPUT DATA FILE - SELECTED TRANSITION:
C          MANTISSA OF: ('IGPOW()' => EXPONENT)
C          GAMMA VALUES             (CASE ' ' & 'P')
C          RATE COEFFT.(CM3 SEC-1) (CASE 'H','R' & 'I')
C          DIMENSION => TEMPERATURE 'SCEF()'
C
C      (C*7)  C7      = USED TO PARSE VALUE FOR XJA()
C      (C*7)  CDELIM  = DELIMITERS FOR INPUT OF DATA FROM HEADERS
C      (C*18) C18     = USED TO PARSE VALUE TO CSTRGA()
C      (C*18) C18T   = COPY OF C18 : UNSATISFACTORY METHOD OF
C          AVOIDING COMPILER REFERENCE ERROR :
C          DHB 07.04.95
C      (C*80) CLINE   = CURRENT ENERGY LEVEL INDEX PARAMETER LINE
C      (C*75) STRING  = TAIL STRING OF 1ST DATA LINE FOR PARSING
C      (C*44) STRG1   = TAIL STRING OF LEVEL SPEC LINES FOR PARSING
C      (C*128) BUFFER = GENERAL STRING BUFFER STORAGE
C      (C*3)  CITPOW() = USED TO PARSE VALUES TO ITPOW()
C      (C*5)  CSCEF() = USED TO PARSE VALUES TO SCEF()
C
C      (L*4)  LDATA   = IDENTIFIES WHETHER THE END OF AN INPUT
C          SECTION IN THE DATA SET HAS BEEN LOCATED.
C          (.TRUE. => END OF SECTION REACHED)
C      (L*4)  LTCHR   = .TRUE.  => CURRENT 'TCODE()' = 'H' OR 'R'
C          OR 'I'
C          = .FALSE. => CURRENT 'TCODE()' .NE. 'H' OR 'R'
C          OR 'I'
C      (L*4)  LTCPR   = .TRUE.  => CURRENT 'TCODE()' = 'P' OR 'R'
C          OR 'I'
C          = .FALSE. => CURRENT 'TCODE()' .NE. 'P' OR 'R'
C          OR 'I'
C      (L*4)  LERROR  = .TRUE.  => UNTIED LEVEL FOUND
C          = .FALSE. => ALL LEVELS TIED
C      (L*4)  LTIED() = .TRUE.  => SPECIFIED LEVEL TIED
C          = .FALSE. => SPECIFIED LEVEL IS UNTIED
C          DIMENSION => LEVEL INDEX

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C NOTE:          LTCHR          LTCPR          TCODE()
C -----
C          .TRUE.          .TRUE.    =>    'R','I'
C          .TRUE.          .FALSE.   =>    'H'
C          .FALSE.         .TRUE.    =>    'P'
C          .FALSE.         .FALSE.   =>    ' '

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C          FOR A-VALUES & GAMMA-VALUES ENTRIES LESS THAN 'DZERO' ARE TAKEN
C          AS BEING EQUAL TO DZERO. THIS AFFECTS THE 'AVAL()' AND 'SCOM()'
C          ARRAYS.

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C ROUTINES:

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ROUTINE	SOURCE	BRIEF DESCRIPTION
I4UNIT	ADAS	FETCH UNIT NUMBER FOR OUTPUT OF MESSAGES
R8FCTN	ADAS	CONVERTS FROM CHARACTER TO REAL VARIABLE
I4FCTN	ADAS	CONVERTS FROM CHAR. TO INTEGER VARIABLE
XXWORD	ADAS	PARSES A STRING INTO SEPARATE WORDS FOR ' () <> {} ' DELIMITERS

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C DATE: 04/06/98

C UPDATE: 1.1 DATE: 09/08/98

C MODIFIED: RICHARD MARTIN

C - PUT UNDER SCCS CONTROL.

C UPDATE: 1.2 DATE: 17/05/07

C MODIFIED: Allan Whiteford

C - Updated comments as part of subroutine documentation
 C procedure.

CHARACTER*92	CIONP
CHARACTER	CPLA (NDLEV)
CHARACTER*9	CPRTA (NDMET)
CHARACTER* (*)	CSTRGA (NDLEV)
CHARACTER	TCODE (NDTRN)
CHARACTER*3	TITLED
INTEGER	I1A (NDTRN) , I2A (NDTRN) , IA (NDLEV) , IL
INTEGER	ILA (NDLEV) , IPLA (NDMET, NDLEV)
INTEGER	ISA (NDLEV) , ITRAN, IUNIT, IZ
INTEGER	IZ0, IZ1, MAXLEV, NDLEV
INTEGER	NDMET, NDTRN, NPL
INTEGER	NPLA (NDLEV) , NV
LOGICAL	LBSETA (NDMET)
REAL*8	AVAL (NDTRN) , BWNO, BWNOA (NDMET)
REAL*8	PRTWTA (NDMET) , SCEF (NVMAX)
REAL*8	SCOM (NVMAX, NDTRN) , WA (NDLEV)
REAL*8	XJA (NDLEV) , ZPLA (NDMET, NDLEV)