

ADAS Subroutine b9data

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SUBROUTINE B9DATA( 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   ,
&                NV    , SCEF    ,
&                ITRAN , MAXLEV ,
&                TCODE , I1A    , I2A    , AVAL   , SCOM   , ITYP
&                )
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C
C ***** FORTRAN77 SUBROUTINE: B9DATA *****
C
C PURPOSE:  TO FETCH DATA FROM INPUT COPASE 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.
C
C CALLING PROGRAM: ADAS209
C
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 SUBROUTINE:
C
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
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
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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: (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
C
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

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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:
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)

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C GAMMA VALUES (CASE ' ' & 'P')

C RATE COEFFT. (CM3 SEC-1) (CASE 'H','R' & 'I')

C DIMENSION => TEMPERATURE 'SCEF()'

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*56) 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 (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

C OUTPUT: (I*4) ITYP = RESOLUTION OF PARENT METASTABLES

C 1 - LS RESOLVED

C 2 - LSJ RESOLVED

C 3 - UNIDENTIFIED

C NOTE:

LTCHR	LTCPR	TCODE()
.TRUE.	.TRUE.	'R','I'
.TRUE.	.FALSE.	'H'
.FALSE.	.TRUE.	'P'
.FALSE.	.FALSE.	' '

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.

C ROUTINES:

ROUTINE	SOURCE	BRIEF DESCRIPTION
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C I4UNIT ADAS FETCH UNIT NUMBER FOR OUTPUT OF MESSAGES
C R8FCTN ADAS CONVERTS FROM CHARACTER TO REAL VARIABLE
C I4FCTN ADAS CONVERTS FROM CHAR. TO INTEGER VARIABLE
C XXWORD ADAS PARSES A STRING INTO SEPARATE WORDS
C FOR ' (<>{}' DELIMITERS
C
C AUTHOR: HP SUMMERS (REVISION OF BXDATA BY PE BRIDEN)
C K1/1/57
C JET EXT. 4941
C
C DATE: 11/06/92
C
C UPDATE: 9/07/93 HPS - USE NEW VERSIONS OF PARSING ROUTINES
C B8PARS AND B8PRS1
C UPDATE: 12/07/93 HPS - REVISE TO CONSISTENCY WITH BXDATA
C AT 25/07/93.
C UPDATE: 11/05/95 HPS - ADDED CPRTA TO PARAMETER LIST.ALTERED
C 'READ()BUFFER' TO BE CONSISTENT WITH IDL-ADAS
C
C UNIX-IDL PORT:
C
C VERSION: 1.1 DATE: 27-06-95
C MODIFIED: TIM HAMMOND (TESSELLA SUPPORT SERVICES PLC)
C - PUT UNDER SCCS CONTROL
C
C VERSION: 1.2 DATE: 19-01-96
C MODIFIED: DAVID BROOKS (UNIVERSITY OF STRATHCLYDE)/TIM HAMMOND
C - INCREASED LENGTH OF CPRTA FROM 4 TO 9 &
C STRING FROM 55 TO 75 IN LINE WITH
C MODIFICATIONS TO ACCOMODATE J-RESOLVED
C PARENT METASTABLES IN THE DATASETS.
C - INCREASED LENGTH OF CLINE TO 92 & STRG1 TO
C 56. ALTERED FORMAT NO. 1003 & READING OF
C CLINE FORMAT TO ACCOMMODATE CHANGES.
C
C VERSION: 1.3 DATE: 26-01-96
C MODIFIED: DAVID BROOKS
C - PASSED ITYP THROUGH TO MAIN PROGRAM
C
C
C VERSION: 1.4 DATE: 18/04/96
C UPDATE: WILLIAM OSBORN
C - INCREASED MTIED TO SAME AS NDLEV
C
C VERSION: 1.5 DATE: 18/11/98
C UPDATE: DAVID BROOKS
C - ALLOWED LEVELS TO 250.
C
C VERSION: 1.6 DATE: 01/11/2002
C MODIFIED: Martin O'Mullane
C - Can handle S lines correctly.
C
C UPDATE: 1.5 DATE: 17/05/07

C MODIFIED: Allan Whiteford
C - Updated comments as part of subroutine documentation
C procedure.
C

C-----
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, ITYP, IUNIT
INTEGER IZ, IZ0, IZ1, MAXLEV
INTEGER NDLEV, 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)