

ADAS Subroutine b4sszd

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      SUBROUTINE B4SSZD( dsname , IBSEL , IZ0IN ,
&                      ITVAL , TVAL ,
&                      BWNO , IZ , IZ1 ,
&                      METI , METF ,
&                      SZDA , ESZDA , LTRNG ,
&                      TITLX , IRCODE , OPEN17
&                      )
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C
C ***** FORTRAN77 SUBROUTINE: B4SSZD *****
C
C PURPOSE: TO EXTRACT AND INTERPOLATE ZERO-DENSITY IONIZATION RATE-
C           COEFFICIENTS FOR GIVEN ELEMENT NUCLEAR CHARGE AND DATA-BLOCK
C           FOR AN INPUT SET OF ELECTRON TEMPERATURES (eV).
C
C
C CALLING PROGRAM: GENERAL USE
C
C SUBROUTINE:
C
C INPUT : (C*80) DSNAME = ADF07 DATAFILE NAME UNDER UNIX INCLUDING PATH
C INPUT : (I*4)  IBSEL  = INDEX OF DATA-BLOCK SELECTED FOR ANALYSIS
C INPUT : (I*4)  IZ0IN  = NUCLEAR CHARGE OF REQUIRED ELEMENT
C
C INPUT : (I*4)  ITVAL  = NUMBER OF ELECTRON TEMPERATURE VALUES
C INPUT : (R*8)  TVAL() = ELECTRON TEMPERATURES (UNITS: EV)
C                       DIMENSION: ELECTRON TEMPERATURE INDEX
C
C OUTPUT: (R*8)  BWNO   = INPUT FILE - SELECTED DATA-BLOCK:
C                       EFFECTIVE IONIZATION POTENTIAL (cm-1).
C OUTPUT: (I*4)  IZ     = INPUT FILE - SELECTED DATA BLOCK:
C                       IONIZING ION - INITIAL CHARGE
C OUTPUT: (I*4)  IZ1    = INPUT FILE - SELECTED DATA BLOCK:
C                       IONIZING ION - FINAL CHARGE
C
C OUTPUT: (I*4)  METI   = INPUT FILE - SELECTED DATA-BLOCK:
C                       INITIAL STATE METSTABLE INDEX
C OUTPUT: (I*4)  METF   = INPUT FILE - SELECTED DATA-BLOCK:
C                       FINAL STATE METSTABLE INDEX
C
C OUTPUT: (R*8)  SZDA() = ZERO-DENSITY IONIZATION RATE-COEFFICIENTS
C                       DIMENSION: ELECTRON TEMPERATURE INDEX
C OUTPUT: (R*8)  ESZDA() = EXP((BWNO/109737.3)*(IH/KTE))*SZDA()
C                       DIMENSION: ELECTRON TEMPERATURE INDEX
C OUTPUT: (L*4)  LTRNG() = .TRUE. => OUTPUT 'SZDA()' VALUE WAS INTER-
C                           POLATED FOR THE USER ENTERED
C                           ELECTRON TEMPERATURE 'TVAL()'.
C                           .FALSE. => OUTPUT 'SZDA()' VALUE WAS EXTRA-
C                           POLATED FOR THE USER ENTERED
C                           ELECTRON TEMPERATURE 'TVAL()'.
C                           DIMENSION: ELECTRON TEMPERATURE INDEX
C
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C OUTPUT: (C*120)TITLX = INFORMATION STRING (DSN ETC.)
C OUTPUT: (I*4) IRCODE = RETURN CODE FROM SUBROUTINE:
C 0 => NORMAL COMPLETION - NO ERROR DETECTED
C 2 => DISCREPANCY BETWEEN REQUESTED CHARGES
C AND THOSE IN INPUT FILE.
C 3 => THE SELECTED DATA-BLOCK 'IBSEL' IS OUT
C OF RANGE OR DOES NOT EXIST.
C 4 => INVALID VALUE FOR 'IZ0IN' ENTERED.
C ('IZ0MIN' <= 'IZ0IN' <= 'IZ0MAX')
C 9 => ERROR ENCOUNTERED WHEN TRYING TO OPEN
C INPUT DATA-SET.
C
C (I*4) NSTORE = PARAMETER= MAXIMUM NUMBER OF DATA-BLOCKS
C WHICH CAN BE READ FROM THE INPUT
C DATA-SET.
C (I*4) NTDIM = PARAMETER= MAXIMUM NUMBER OF ELECTRON TEMP-
C ERATURES THAT CAN BE READ FROM
C AN INPUT DATA-SET DATA-BLOCK.
C (I*4) IZ0MIN = PARAMETER: MIN. ALLOWED VALUE FOR 'IZ0IN'
C (I*4) IZ0MAX = PARAMETER: MAX. ALLOWED VALUE FOR 'IZ0IN'
C
C (I*4) IZ0LST = LAST VALUE OF 'IZ0IN' FOR WHICH INPUT
C DATA WAS READ.
C (I*4) IUNIT = UNIT TO WHICH INPUT DATA SET IS ALLOCATED
C (I*4) NBSEL = TOTAL NUMBER OF DATA-BLOCKS READ FROM INPUT
C DATA SET.
C (I*4) IZ0 = INPUT FILE - EMITTING ION - NUCLEAR CHARGE
C
C (L*4) LOPEN = .TRUE. => INPUT DATA SET OPEN.
C .FALSE. => INPUT DATA SET CLOSED.
C
C (C*2) ESYM = INPUT FILE - IONIZING ION - ELEMENT SYMBOL
C (C*3) EXTIN = CURRENT ADAS SOURCE DATA FILE EXTENSION
C (C*3) EXTLST = ADAS SOURCE DATA FILE EXT. USED LAST TIME
C DATA WAS READ.
C
C (I*4) ISELA() = INPUT DATA FILE: DATA-BLOCK ENTRY INDICES.
C DIMENSION: DATA-BLOCK INDEX
C (I*4) ITA() = INPUT DATA SET-NUMBER OF ELECTRON TEMPERA-
C TURES.
C DIMENSION: DATA-BLOCK INDEX
C (I*4) IZOUT() = INPUT DATA FILE: IONIZING ION INITIAL CHARGE
C DIMENSION: DATA-BLOCK INDEX
C (I*4) IZ1OUT() = INPUT DATA FILE: IONIZING ION FINAL CHARGE
C DIMENSION: DATA-BLOCK INDEX
C
C (R*8) BWNOUT() = INPUT DATA FILE: EFFECTIVE IONIZATION POT.
C (UNITS: cm-1).
C DIMENSION: DATA-BLOCK INDEX
C (R*8) TETA(,) = INPUT DATA SET -
C ELECTRON TEMPERATURES (UNITS: eV)
C 1st DIMENSION: ELECTRON TEMPERATURE INDEX
C 2nd DIMENSION: DATA-BLOCK INDEX

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C      (R*8)  SZD(,)    =INPUT DATA SET -
C                      FULL SET OF IONIZATIONS RATE-COEFFICIENTS
C                      1st DIMENSION: ELECTRON TEMPERATURE INDEX
C                      3rd DIMENSION: DATA-BLOCK INDEX
C
C      (C*2)  CICODE()= INPUT DATA FILE - INITIAL STATE META. INDEX
C                      DIMENSION: DATA-BLOCK INDEX
C      (C*2)  CFCODE()= INPUT DATA FILE - FINAL STATE META. INDEX
C                      DIMENSION: DATA-BLOCK INDEX
C      (C*6)  CIION() = INPUT DATA FILE - INITIAL ION
C                      DIMENSION: DATA-BLOCK INDEX
C      (C*6)  CFION() = INPUT DATA FILE - FINAL ION
C                      DIMENSION: DATA-BLOCK INDEX
C
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C

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C ROUTINES:

ROUTINE	SOURCE	BRIEF DESCRIPTION
XXDATA_07	ADAS	FETCH INPUT DATA FROM SELECTED DATA SET
B4SPLN	ADAS	INTERPOLATE DATA WITH ONE-WAY SPLINES
E2TITL	ADAS	CREATE DESCRIPTIVE TITLE FOR OUTPUT

C Original version

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C JET EXT. 6023

C DATE: 07/06/91

C UPDATE: 17/02/97 - H P SUMMERS: RENAME SSZD AS B4SSZD. EXTRACT
C Exp(I/KTE) * S AS WELL AS S

C UPDATE: 04-03-97 - R. MARTIN: ADDED OPEN17 FOR SWITCHING OUTPUT TO
C 'adas204.pass1' ON AND OFF.

C VERSION: 1.2 DATE: 03-12-98

C MODIFIED: Martin O'Mullane:
C Rewritten to account for adf07 filename
C being included in the adf25 namelist. It is
C now much simplified.

C VERSION: 1.3 DATE: 26-03-08

C MODIFIED: Allan Whiteford:
C Changed call from E2DATA to XXDATA_07.

CHARACTER*120	DSNAME,	TITLX		
INTEGER	IBSEL,	IRCODE,	ITVAL,	IZ

INTEGER	IZOIN,	IZ1,	METF,	METI
LOGICAL	LTRNG(ITVAL),		OPEN17	
REAL*8	BWNO,	ESZDA(ITVAL)		
REAL*8	SZDA(ITVAL),	TVAL(ITVAL)		