

ADAS Subroutine dxspl1

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      SUBROUTINE DXSPL1( ISWIT , LSWIT , IZ1 ,
&                      NDOUT , NTOUT ,
&                      NDIN  , NTIN  , NZIN  ,
&                      IDE   , ITE   , IZE   ,
&                      TIN   , ZIPT  , EIA   ,
&                      AIPT  ,
&                      ZINTRP ,
&                      ATTY  ,
&                      )
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C
C ***** FORTRAN77 SUBROUTINE: DXSPL1 *****
C
C PURPOSE: PERFORMS THE FIRST PART OF A THREE WAY SPLINE ON INPUT DATA.
C           GENERATES A TABLE OF LOG10(SCALED COEF/PWRS) COVERING 'ITE'
C           TEMPERATURES AND 'IDE' DENSITIES FOR THE ELEMENT RECOMBINING
C           ION CHARGE GIVEN BY 'IZ1'.
C
C CALLING PROGRAM: D1SPLN/D4DATA
C
C DATA:
C
C           THE SOURCE DATA IS CONTAINED AS MEMBERS OF PARTITIONED
C           DATA SETS AS FOLLOWS:
C
C           1. JETUID.ACD<YR>.DATA
C           2. JETUID.SCD<YR>.DATA
C           3. JETUID.CCD<YR>.DATA
C           4. JETUID.PRB<YR>.DATA
C           5. JETUID.PRC<YR>.DATA
C           6. JETUID.PRB<YR>.DATA
C           7. JETUID.PRC<YR>.DATA
C           8. JETUID.PLT<YR>.DATA
C           9. JETUID.PLS<YR>.DATA
C           10. JETUID.MET<YR>.DATA
C
C           WHERE <YR> DENOTES TWO INTEGERS FOR THE YEAR SELECTED.
C
C           THE PARTICULAR TYPE OPENED (1-10) IS SELECTED BY 'ISWIT'
C
C SUBROUTINE:
C
C INPUT : (I*4) ISWIT = DATA TYPE SELECTOR (SEE ABOVE) (1 -> 10)
C INPUT : (L*4) LSWIT = .TRUE. => IONISATION POTENTIALS PRESENT
C           .FALSE. => IONS. POTENTIALS NOT PRESENT
C INPUT : (I*4) IZ1 = OUTPUT - ELEMENT RECOMBINING ION CHARGE
C
C INPUT : (I*4) NDOUT = OUTPUT - MAXIMUM NUMBER OF DENSITIES
C INPUT : (I*4) NTOUT = OUTPUT - MAXIMUM NUMBER OF TEMPERATURES
C
C INPUT : (I*4) NDIN = INPUT - MAXIMUM NUMBER OF DENSITIES
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C INPUT : (I*4)  NTIN    = INPUT  - MAXIMUM NUMBER OF TEMPERATURES
C INPUT : (I*4)  NZIN    = INPUT  - MAXIMUM NUMBER OF CHARGE STATES
C
C INPUT : (I*4)  ITE     = INPUT  - NUMBER OF REDUCED TEMPERATURES
C INPUT : (I*4)  IDE     = INPUT  - NUMBER OF REDUCED DENSITIES
C INPUT : (I*4)  IZE     = INPUT  - NUMBER OF CHARGE STATES/RECOMBINING
C                               ION CHARGE
C
C INPUT : (R*8)  TIN()   = INPUT  - SET OF 'ITE' REDUCED ELECTRON TEMPS.
C INPUT : (R*8)  ZIPT()  = INPUT  - SET OF 'IZE' RECOMBINING ION CHARGES
C INPUT : (R*8)  EIA()   = IONISATION POTENTIALS: ()=ION CHARGE
C                               (UNITS: RYDBERGS)
C
C INPUT : (R*8)  AIPT(,,) = INPUT  - COEFFICIENT/POWER ARRAY.
C                               1ST DIMENSION: REDUCED DENSITY ('DENSUR()')
C                               2ND DIMENSION: REDUCED TEMPERATURE ('TR()')
C                               3RD DIMENSION: CHARGE STATE ('ZIPT()')
C
C OUTPUT: (L*4)  ZINTRP(1) = .TRUE.  => 'ATTY(,)' VALUES INTERPOLATED
C                               = .FALSE. => 'ATTY(,)' VALUES EXTRAPOLATED
C
C OUTPUT: (R*8)  ATTY(,) = WORKING SPACE FOR 3-WAY SPLINE ITERPOLATION
C                               (STORES VALUES OF 'ANS(1)')
C                               1ST DIMENSION: TEMPERATURE
C                               2ND DIMENSION: DENSITY
C
C           (I*4)  NZDIM1  = PARAMETER = MUST BE EQUAL TO OR GREATER THAN
C                               THE MAXIMUM NUMBER OF INPUT CHARGE STATES
C           (I*4)  L1      = PARAMETER = 1
C
C           (R*8)  TK2ATE  = PARAMETER = EQUATION CONSTANT = 157890
C
C           (I*4)  I4UNIT  = FUNCTION (SEE ROUTINE SECTION BELOW)
C           (I*4)  IT      = ARRAY SUBSCRIPT USED FOR TEMPERATURE VALUES
C           (I*4)  ID      = ARRAY SUBSCRIPT USED FOR DENSITY VALUES
C           (I*4)  IZ      = ARRAY SUBSCRIPT USED FOR CHARGE STATE VALUES
C           (I*4)  JZ      = RECOMBINING ION CHARGE FOR CHARGE STATE 'IZ'
C           (I*4)  IOPT    = DEFINES THE BOUNDARY DERIVATIVES FOR THE
C                               SPLINE ROUTINE 'XXSPLE', SEE 'XXSPLE'.
C                               (VALID VALUES = 0, 1, 2, 3)
C
C           (L*4)  LSETX   = .TRUE.  => SET UP SPLINE PARAMETERS RELATING
C                               TO 'XIN' AXIS.
C                               .FALSE. => DO NOT SET UP SPLINE PARAMETERS
C                               RELATING TO 'XIN' AXIS.
C                               (I.E. THEY WERE SET IN A PREVIOUS
C                               CALL )
C                               (VALUE SET TO .FALSE. BY 'XXSPLE')
C
C           (R*8)  ATE     = 'TK2ATE' / INPUT REDUCED TEMPERATURE
C           (R*8)  Z2     = 'JZ' SQUARED
C           (R*8)  Z3     = 'JZ' CUBED
C           (R*8)  ZL     = LOG10 OF CURRENT CHARGE STATE

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C      (R*8)  Y      = SCALED 'AIPT(,,)' VALUE
C      (R*8)  Z1(1) = IZ1
C      (R*8)  ANS(1) = SPLINE INTERPOLATED LOG10(SCALED 'AIPT(,,)')
C                      VALUE FOR A RECOMBINING ION CHARGE EQUAL TO
C                      'IZ1' AT FIXED TEMPERATURE AND DENSITY.
C      (R*8)  F()   = LOG10 ( 'Y' ) - DIMENSION => CHARGE STATE
C      (R*8)  DF()  = SPLINE INTERPOLATED DERIVATIVES

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C NOTE:

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C                      SPLINE IS CARRIED OUT ON:
C      LOG10(SCALED 'AIPT(,,)' VALUES AT FIXED TEMP AND DENSITY)
C                      VERSUS
C                      RECOMBINING ION CHARGE

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C ROUTINES:

ROUTINE	SOURCE	BRIEF DESCRIPTION
XXSPLE	ADAS	SPLINE SUBROUTINE (WITH EXTRAP. INFO)
I4UNIT	ADAS	INTEGER*4 FUNCTION - FETCH UNIT NUMBER FOR OUTPUT OF MESSAGES
R8FUN2	ADAS	REAL*8 FUNCTION: (X -> 1/(1+X))

C AUTHOR: PAUL E. BRIDEN (TESSELLA SUPPORT SERVICES PLC)
 K1/0/37
 JET EXT. 2520

C DATE : 13/06/91 - PE BRIDEN: ADAS91 VERSION OF 'D4SPL1'
 C DATE : 15/12/92 - PE BRIDEN: IF IZE<=1 THEN ANS(1)=F(1) INSTEAD
 C OF ANS(1)=F(IZ) (WHICH IN ERROR GIVES
 C ANS(1)=F(2) IF IZE=1.)

C UPDATE: 12/08/93 - HP SUMMERS: INCLUDE ISWIT IN PARAMETER AND ALLOW
 C SEPARATE TREATMENT ACCORDING TO ISWIT
 C UPDATE: 18/08/93 - HP SUMMERS: ALTER ORDER TO TAKE LOG10 FIRST BEFORE
 C ADJUSTING FOR DATA CLASS

C UNIX-IDL PORT:

C VERSION: 1.1 DATE: 06-09-95
 C MODIFIED: TIM HAMMOND (TESSELLA SUPPORT SERVICES PLC)
 C - FIRST RELEASE

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      INTEGER      IDE,      ISWIT,      ITE,      IZ1
      INTEGER      IZE,      NDIN,      NDOUT,      NTIN
      INTEGER      NTOUT,      NZIN
      LOGICAL      LSWIT,      ZINTRP(1)
      REAL*8       AIPT(NDIN,NTIN,NZIN),      ATTY(NTOUT,NDOUT)
      REAL*8       EIA(50),      TIN(NTIN),      ZIPT(NZIN)

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