

ADAS Subroutine dxspl3

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      SUBROUTINE DXSPL3( ISWIT , LSWIT , IZ1 ,
&                      NDOUT , NTOUT ,
&                      NTIN ,
&                      MAXD , ITE ,
&                      MAXT , TIN , TOUT , EIAVAL ,
&                      TINTRP ,
&                      ATTY
&                      )
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C
C ***** FORTRAN77 SUBROUTINE: DXSPL3 *****
C
C PURPOSE: PERFORMS THE THIRD PART OF A THREE WAY SPLINE ON INPUT DATA.
C          GENERATES A TABLE OF LOG10( COEFFTS./POWERS ) COVERING
C          'MAXT' TEMPERATURES AND 'MAXD' DENSITIES FOR THE ELEMENT
C          RECOMBINING ION CHARGE GIVEN BY 'IZ1'.
C
C CALLING PROGRAM: D1SPLN/D4DATA
C DATA:
C
C          THE SOURCE DATA ORIGINATES 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          IT IS PASSED IN A MODIFIED FORM AFTER PROCESSING BY
C          DXSPL1 AND DXSPL2.
C
C SUBROUTINE:
C
C INPUT : (I*4) ISWIT = DATA TYPE SELECTOR (1 -> 8)
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) NTIN = INPUT - MAXIMUM NUMBER OF TEMPERATURES
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C
C INPUT : (I*4)  MAXD    = INPUT  - NUMBER OF REDUCED DENSITIES
C INPUT : (I*4)  ITE     = INPUT  - NUMBER OF REDUCED TEMPERATURES
C
C INPUT : (I*4)  MAXT    = OUTPUT  - NUMBER OF REDUCED TEMPERATURES
C                                     ( <= NTOUT )
C INPUT : (R*8)  TIN()   = INPUT  - SET OF 'ITE' REDUCED ELECTRON TEMPS
C INPUT : (R*8)  TOUT()  = OUTPUT  - SET OF 'MAXT' ELECTRON TEMPERATURES
C                                     (UNITS: KELVIN) .
C INPUT : (R*8)  EIAVAL  = IONISATION POTENTIAL (RYDBERGS) FOR
C                                     THE ION CHARGE GIVEN BY 'IZ1' .
C
C OUTPUT: (L*4)  TINTRP() = .TRUE.  => 'ATTY(,)' VALUE FOR TEMPERATURE
C                                     INDEX INTERPOLATED.
C                                     = .FALSE. => 'ATTY(,)' VALUE FOR TEMPERATURE
C                                     INDEX EXTRAPOLATED.
C                                     1ST DIMENSION: TEMPERATURE INDEX
C
C IN/OUT: (R*8)  ATTY(,) = WORKING SPACE FOR 3-WAY SPLINE ITERPOLATION
C                                     (STORES LOG10 (INTERPOLATED VALUES))
C                                     INPUT 'ATTY' VALUES ARE ASSIGNED TO 'YIN' &
C                                     THEN 'YOUT' VALUES ARE ASSIGNED TO 'ATTY' .
C                                     1ST DIMENSION: TEMPERATURE
C                                     2ND DIMENSION: DENSITY
C
C          (I*4)  NTDIM1  = PARAMETER = MUST BE EQUAL TO OR GREATER THAN
C                                     THE MAXIMUM NUMBER OF INPUT TEMPERATURES.
C          (I*4)  NTDIM2  = PARAMETER = MUST BE EQUAL TO OR GREATER THAN
C                                     THE MAXIMUM NUMBER OF OUTPUT TEMPERATURES.
C
C          (R*8)  EIACON  = PARAMETER = -68570.7
C
C          (I*4)  IT      = ARRAY SUBSCRIPT USED FOR TEMPERATURE VALUES
C          (I*4)  ID      = ARRAY SUBSCRIPT USED FOR DENSITY VALUES
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)  Z1R2    = 1.0 / (IZ1**2)
C          (R*8)  SCON1   = SCALING CONSTANT
C          (R*8)  SCON2   = SCALING CONSTANT
C          (R*8)  XOUT()  ='TOUT()' CONVERTED TO REDUCED TEMPERATURE
C          (R*8)  YIN()   ='ATTY(,)' AT FIXED DENSITY -
C                                     DIMENSION => TEMPERATURE
C          (R*8)  DF()    = SPLINE INTERPOLATED DERIVATIVES

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C (R*8) YOUT() = SPLINE INTERPOLATED 'ATTY(,,)' VALUES FOR
C REDUCED TEMPERATURE EQUAL TO 'XOUT()' AT A
C FIXED DENSITY.
C

C NOTE:

C SPLINE IS CARRIED OUT ON:
C 'ATTY(,,)' VALUES AT FIXED DENSITY
C VERSUS
C LOG10(REDUCED DENSITY)
C

C ROUTINES:

ROUTINE	SOURCE	BRIEF DESCRIPTION
XXSPLE	ADAS	SPLINE SUBROUTINE (WITH EXTRAP. INFO)
I4UNIT	ADAS	UNIT NUMBER FOR WARNING MESSAGES

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

C DATE : 13/06/91 - PE BRIDEN: ADAS91 VERSION OF 'D4SPL3'

C UNIX-IDL PORT:

C VERSION: 1.1 DATE: 06-09-95
C MODIFIED: TIM HAMMOND (TESSELLA SUPPORT SERVICES PLC)
C - FIRST RELEASE
C VERSION: 1.2 DATE: 25-10-97
C MODIFIED: LORNE HORTON (JET)
C - ADDED FURTHER CHECKS ON LOW T EXTRAPOLATIONS
C

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INTEGER	ISWIT,	ITE,	IZ1,	MAXD
INTEGER	MAXT,	NDOUT,	NTIN,	NTOUT
LOGICAL	LSWIT,	TINTRP (NTOUT)		
REAL*8	ATTY (NTOUT, NDOUT) ,	EIAVAL		
REAL*8	TIN (NTIN) ,	TOUT (NTOUT)		