

ADAS Subroutine b4spln

```
      SUBROUTINE B4SPLN( ITA      , ITVAL      ,  
&                      BWNO      ,  
&                      TETA      , TEVA      ,  
&                      SZD        ,  
&                      SZDA      , ESZDA     ,  
&                      LTRNG  
&                      )
```

```
C-----  
C  
C *****  
C ***** FORTRAN77 SUBROUTINE: B4SPLN *****  
C *****  
C  
C PURPOSE:  
C     PERFORMS CUBIC SPLINE ON LOG(TEMPERATURE <EV> ) VERSUS  
C     LOG(SCALED IONIZATION RATE COEFFICIENTS).  
C     INPUT DATA FOR A GIVEN IONIZING ION COMBINATION DATA-BLOCK.  
C  
C     USING ONE-WAY SPLINES IT CALCULATES THE IONIZATION RATE  
C     COEFFICIENT FOR 'ITVAL' ELECTRON TEMPERATURE VALUES FROM  
C     THE LIST OF ELECTRON TEMPERATURES READ IN FROM THE INPUT FILE  
C  
C     IF A VALUE CANNOT BE INTERPOLATED USING SPLINES IT IS  
C     EXTRAPOLATED VIA 'XXSPLE'. (SEE NOTES BELOW).  
C  
C CALLING PROGRAM: ADAS204/B4SSZD  
C  
C  
C SUBROUTINE:  
C  
C INPUT : (I*4)  ITA      = INPUT DATA FILE: NUMBER OF ELECTRON TEMPERA-  
C                   TURES READ FOR THE DATA-BLOCK BEING ASSESSED  
C INPUT : (I*4)  ITVAL    = NUMBER OF ISPF ENTERED ELECTRON TEMPERATURE  
C                   VALUES FOR WHICH IOINIZATION RATE COEFFFTS  
C                   ARE REQUIRED FOR TABULAR/GRAPHICAL OUTPUT.  
C  
C INPUT : (R*8)  BWNO     = INPUT DATA FILE: IONIZATION POTENTIAL (cm-1)  
C                   FOR THE DATA-BLOCK BEING ASSESSED.  
C  
C INPUT : (R*8)  TETA()   = INPUT DATA FILE: ELECTRON TEMPERATURES (EV)  
C                   FOR THE DATA-BLOCK BEING ASSESSED.  
C                   DIMENSION: ELECTRON TEMPERATURE INDEX  
C INPUT : (R*8)  TEVA()   = USER ENTERED: ELECTRON TEMPERATURES (EV)  
C                   DIMENSION: TEMPERATURE/DENSITY PAIR INDEX  
C  
C INPUT : (R*8)  SZD()    =INPUT DATA FILE: FULL SET OF ZERO DENSITY  
C                   IONIZATION RATE COEFFFTS FOR THE DATA-BLOCK  
C                   BEING ANALYSED.  
C                   1ST DIMENSION: ELECTRON TEMPERATURE INDEX  
C OUTPUT: (R*8)  SZDA()   = SPLINE INTERPOLATED OR EXTRAPOLATED ZERO  
C                   DENSITY IONIZATION RATE COEFFICIENTS FOR  
C                   THE USER ENTERED ELECTRON TEMPERATURES.  
C                   DIMENSION: ELECTRON TEMPERATURE INDEX  
C OUTPUT: (R*8)  ESZDA()  = EXP ( (BWNO/109737.3) * (IH/KTE) ) *SZDA ( )
```

C
C OUTPUT: (L*4) LTRNG() = .TRUE. => OUTPUT 'SZDA()' VALUE WAS INTER-
C POLATED FOR THE USER ENTERED
C ELECTRON TEMPERATURE 'TEVA()'.
C .FALSE. => OUTPUT 'SZDA()' VALUE WAS EXTRA-
C POLATED FOR THE USER ENTERED
C ELECTRON TEMPERATURE 'TEVA()'.
C DIMENSION: ELECTRON TEMPERATURE INDEX
C
C (I*4) NIN = PARAMETER = MAX. NO. OF INPUT TEMPERATURE
C VALUES. MUST BE >= 'ITA'
C (I*4) NOUT = PARAMETER = MAX. NO. OF OUTPUT TEMPERATURE
C PAIRS. MUST BE >= 'ITVAL'
C (I*4) L1 = PARAMETER = 1
C
C (R*8) BCONST = PARAMETER = 1/(SCALED BOLTZMANN'S CONSTANT)
C
C (I*4) IET = ARRAY SUBSCRIPT USED INPUT FILE ELECTRON
C TEMPERATURES.
C (I*4) IT = ARRAY SUBSCRIPT USED FOR USER ENTERED
C TEMPERATURE VALUES.
C (I*4) IOPT = DEFINES THE BOUNDARY DERIVATIVES FOR THE
C SPLINE ROUTINE 'XXSPLE', SEE 'XXSPLE'.
C (VALID VALUES = <0, 0, 1, 2, 3, 4)
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) R8FUN1 = FUNCTION - (SEE ROUTINES SECTION BELOW)
C (R*8) SCONST = SCALING CONSTANT USED TO SCALE THE IONIZA-
C TION RATE COEFFT. WHEN SPLINNING.
C = IONIZATION POTENTIAL / BOLTZMANN CONST.
C
C (R*8) XIN() = LOG(DATA FILE ELECTRON TEMPERATURES)
C (R*8) YIN() = LOG(DATA FILE SCALED ION. RATE COEFFTS.)
C (R*8) XOUT() = LOG(USER ENTERED ELECTRON TEMPS.)
C (R*8) YOUT() = LOG(OUTPUT GENERATED SCALED ION. RATE COEF)
C (R*8) DF() = SPLINE INTERPOLATED DERIVATIVES
C
C
C

C NOTE:

C ONE-DIMENSIONAL SPLINE CARRIED OUT BY THIS SUBROUTINE:

C LOG(EXP(<ion.pt.>/<k>.<Te>) . Szd) vs. LOG(Te)

C ion.pt. = ionization potential (units: cm-1)

C k = Boltzmann's constant (= 1/1.23977E-04)

```

C      Te      = electron temperature (units: eV)
C      Szd     = zero density ionization rate coefficient
C                  (units: cm**3/sec)
C
C      Extrapolation criteria:
C
C      Low Te: zero gradient extrapolation (i.e. DY(1) = 0.0)
C      High Te: zero curvature extrapolation (i.e. DDY(N) = 0.0)
C
C      (These criteria are met by calling XXSPLE with IOPT=4)
C
C
C ROUTINES:
C      ROUTINE      SOURCE      BRIEF DESCRIPTION
C      -----
C      XXSPLE      ADAS        SPLINE SUBROUTINE (EXTENDED DIAGNOSTICS)
C      R8FUN1      ADAS        REAL*8 FUNCTION: ( X -> X )
C
C AUTHOR:  PAUL E. BRIDEN (TESSELLA SUPPORT SERVICES PLC)
C          K1/0/37
C          JET EXT. 2620
C
C DATE:    07/06/91
C
C UPDATE:  17/02/97 HP SUMMERS - ADDED EXP(IP/KTE)*SZD AS AN OUTPUT
C                  PARAMETER.
C
C VERSION: 1.1 DATE: 05-03-98
C MODIFIED: H.SUMMERS, L.HORTON, M.OMULLANE, R.MARTIN
C          - BASED ON E2SPLN.FOR v1.2. PUT UNDER SCCS CONTROL.
C
C-----
C
C-----
C
C      INTEGER      ITA,      ITVAL
C      LOGICAL      LTRNG(ITVAL)
C      REAL*8       BWNO,      ESZDA(ITVAL)
C      REAL*8       SZD(ITA),  SZDA(ITVAL), TETA(ITA)
C      REAL*8       TEVA(ITVAL)

```