

## ADAS Subroutine cxbms

SUBROUTINE CXBMS (DSNIN, NSITYP, IOUNIT, SIFRAC, UBMENG, UTDENS,  
& UTTEMP, NREQ, MXREQS, BSTOT)

```
C-----  
C  
C          ***** FORTRAN 77 SUBROUTINE: CXBMS *****  
C  
C PURPOSE: TO ASSEMBLE COMPOSITE BEAM STOPPING OF EMISSION  
C COEFFICIENTS USING THE LINEAR INTERPOLATION  
C AND COMBINATION METHOD.  
C  
C  
C INPUT  
C  
C (R*8)  UBMENG : USER REQUESTED NEUTRAL BEAM ENERGIES  
C UNITS: EV/AMU  
C DIMENSION: NREQ  
C (R*8)  UTDENS : USER REQUESTED TARGET DENSITIES  
C UNITS: CM-3  
C DIMENSION: NREQ  
C (R*8)  UTTEMP : USER REQUESTED TARGET TEMPERATURES  
C UNITS: EV  
C DIMENSION: NREQ  
C (R*8)  SIFRAC : FRACTIONAL IMPURITY CONTENT.  
C 1ST DIMENSION: MXREQS  
C  
C 2ND DIMENSION: NSITYP  
C (I*4)  NSITYP : NUMBER OF PLASMA IMPURITY IONS.  
C (I*4)  NREQ : NUMBER OF REQUESTED ENERGIES,  
C DENSITIES AND TEMPERATURES.  
C (I*4)  MXREQS : SIZE OF FIRST DIMENSION OF SIFRAC  
C  
C (I*4)  IOUNIT : UNIT NUMBER EMPLOYED TO READ ADF21  
C AND ADF22 TYPE FILES.  
C  
C (CHR)  DSNIN() : ARRAY CONTAINING NAME OF EACH FILE TO BE READ.  
C DIMENSION: NSITYP  
C  
C  
C OUTPUT  
C  
C (R*8)  BSTOT() : TOTAL BEAM STOPPING COEFFICIENTS.  
C DIMENSION: NREQ  
C  
C GENERAL  
C  
C (R*8)  BREF() : REFERENCE BEAM ENERGIES.  
C UNITS: EV/AMU  
C DIMENSION: MXIT  
C (R*8)  TDREF() : REFERENCE TARGET DENSITIES.  
C UNITS: CM-3  
C DIMENSION: MXIT  
C (R*8)  TTREF() : REFERENCE TARGET TEMPERATURES.
```

```

C          UNITS: EV
C          DIMENSION: MXIT
C (R*8)  SVREF ( ) : STOPPING COEFFT. AT REFERENCE BEAM ENERGY,
C          TARGET DENSITY AND TEMPERATURE.
C          UNITS: CM3 S-1
C          DIMENSION: MXIT
C (R*8)  BE ( , ) : BEAM ENERGIES.
C          UNITS: EV/AMU
C          1ST DIMENSION: MXBE
C          2ND DIMENSION: MXIT
C (R*8)  TDENS ( , ) : TARGET DENSITIES.
C          UNITS: CM-3
C          1ST DIMENSION: MXTD
C          2ND DIMENSION: MXIT
C (R*8)  TTEMP ( , ) : TARGET TEMPERATURES.
C          UNITS: EV
C          1ST DIMENSION: MXTT
C          2ND DIMENSION: MXIT
C (R*8)  SVT ( , ) : STOPPING COEFFT. AT REFERENCE BEAM ENERGY
C          AND TARGET DENSITY.
C          UNITS: CM3 S-1
C          1ST DIMENSION: MXTT
C          2ND DIMENSION: MXIT
C (R*8)  SVED ( , , ) : STOPPING COEFFT. AT REFERENCE TARGET
C          TEMPERATURE.
C          UNITS: CM3 S-1
C          1ST DIMENSION: MXBE
C          2ND DIMENSION: MXTD
C          3RD DIMENSION: MXIT
C (R*8)  BSION ( , ) : BEAM STOPPING COEFFICIENTS FOR INDIVIDUAL
C          IONS.
C          1ST DIMENSION: MXREQ
C          2ND DIMENSION: MXIT
C
C (R*8)  FACT1      : FACTOR USED IN CALCULATING ZEFF
C (R*8)  FACT2      : SIMILAR TO FACT1.
C (R*8)  ZEFF ( ) : USED SO THAT THE BEAM STOPPING
C          COEFFICIENTS FOR INDIVIDUAL
C          IMPURITY IONS ARE EVALUATED AT THE
C          CORRECT EFFECTIVE ELECTRON DENSITY.
C (R*8)  WT : WEIGHTING FACTOR.
C (R*8)  TFRAC : GENERAL VARIABLE.
C
C (I*4)  MXREQ      : MAXIMUM NUMBER OF REQUESTED ENERGIES
C          DENSITIES AND TEMPERATURES
C (I*4)  MXIT       : MAXIMUM POSSIBLE NUMBER OF DIFFERENT
C          PLASMA IONS.
C (I*4)  MXBE      : MAXIMUM NUMBER OF BEAM ENERGIES WHICH CAN
C          BE READ FROM THE ADF21/22 TYPE FILES.
C (I*4)  MXTD      : MAXIMUM NUMBER OF TARGET DENSITIES WHICH CAN
C          BE READ FROM THE ADF21/22 TYPE FILES.
C (I*4)  MXTT      : MAXIMUM NUMBER OF TARGET TEMPERATURES WHICH
C          CAN BE READ FROM THE ADF21/22 TYPE FILES.

```

C (I\*4) ITZ() : TARGET ION CHARGE.  
 C DIMENSION: MXIT  
 C (I\*4) NBE() : NUMBER OF BEAM ENERGIES.  
 C DIMENSION: MXIT  
 C (I\*4) NTDENS() : NUMBER OF TARGET DENSITIES.  
 C DIMENSION: MXIT  
 C (I\*4) NTTEMP() : NUMBER OF TARGET TEMPERATURES.  
 C DIMENSION: MXIT  
 C  
 C (CHR) TSYM() : TARGET ION ELEMENT SYMBOL.  
 C DIMENSION: MXIT  
 C  
 C (L\*4) LIBMA(,) : FLAGS IF INTERPOLATION OR EXTRAPOLATION  
 C USED FOR REQUESTED BEAM ENERGIES.  
 C .TRUE. => INTERPOLATION USED.  
 C .FALSE. => EXTRAPOLATION USED.  
 C 1ST DIMENSION: MXREQ  
 C 2ND DIMENSION: MXIT

C NOTE: USE OF FLAGS NOT IMPLEMENTED

C (L\*4) LIDNA(,) : FLAGS IF INTERPOLATION OR EXTRAPOLATION  
 C USED FOR REQUESTED ION DENSITIES.  
 C .TRUE. => INTERPOLATION USED.  
 C .FALSE. => EXTRAPOLATION USED.  
 C 1ST DIMENSION: MXREQ  
 C 2ND DIMENSION: MXIT

C NOTE: USE OF FLAGS NOT IMPLEMENTED

C (L\*4) LITIA(,) : FLAGS IF INTERPOLATION OR EXTRAPOLATION  
 C USED FOR REQUESTED ION TEMPERATURES.  
 C .TRUE. => INTERPOLATION USED.  
 C .FALSE. => EXTRAPOLATION USED.  
 C 1ST DIMENSION: MXREQ  
 C 2ND DIMENSION: MXIT

C NOTE: USE OF FLAGS NOT IMPLEMENTED

C ROUTINES:

ROUTINE	SOURCE	BRIEF DESCRIPTION
C4DATA	ADAS	READS INPUT DATA SET IN ADF21/22 FORMAT.
C4SPLN	ADAS	PERFORMS SPLINE ON INPUT DATA.
I4UNIT	ADAS	RETURNS UNIT NO. FOR OUTPUT OF MESSAGES.

C NOTE : THE IMPURITY FRACTIONS ARE RENORMNALISED TO  
 C ENSURE THAT THE TOTAL SUM OF EACH IMPURITY  
 C FRACTION DOES NOT EXCEED THE VALUE OF ONE.

C AUTHOR: HARVEY ANDERSON  
 C UNIVERSITY OF STRATHCLYDE  
 C ANDERSON@PHYS.STRATH.AC.UK

C DATE : 30/09/99

```

C
C VERSION : 1.1
C     DATE      : 3-6-2000
C MODIFIED: Martin O'Mullane
C   Repositioned declaration of variables into standard
C             ADAS convention and to satisfy g77.
C
C VERSION : 1.2
C     DATE      : 19-3-2003
C MODIFIED: Lorne Horton
C   Increased MXREQ to 1024.  Added check on internal
C             matrix sizes.  Added LSET flag to allow faster
C             splines on repeated calls to C4SPLN.
C
C VERSION : 1.3
C     DATE      : 03-12-2004
C MODIFIED: Martin O'Mullane
C   - Replace c4data with xxdata_21.
C   - Increase dsnin length to 132 characters.
C   - Merge L Horton's changes into central version.
C
C VERSION : 1.4
C     DATE      : 08-12-2004
C MODIFIED: Martin O'Mullane/Allan Whiteford
C   - Increase size of dsnsav to 132 and set initial
C     values to ' ' rather than ''.
C
C-----

```

CHARACTER*132	DSNIN(NSITYP)			
INTEGER	IOUNIT,	MXREQS,	NREQ,	NSITYP
REAL*8	BSTOT(NREQ),	SIFRAC(MXREQS,NSITYP)		
REAL*8	UBMENG(NREQ),		UTDENS(NREQ)	
REAL*8	UTTEMP(NREQ)			