

ADAS Subroutine d9rdnm

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C Copyright (c) 1997, Strathclyde University.
  SUBROUTINE D9RDNM( DSNINC , LPART , IFAIL ,
&                   IZ0    , NPART , IPRTD , IGRDD , ICLASS ,
&                   IZ1    , ITMAX  , IDMAX  ,
&                   ISDIMD , IZDIMD , ITDIMD ,
&                   ISMAXD , IZMAXD , ITMAXD , IDMAXD , NPARTR,
&                   DTEV   , DDENS  ,
&                   DTEVD  , DDENSD , DRCOFD , ZDATA  ,
&                   DRCOFI
&                   )
C
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C
C ***** FORTRAN77 SUBROUTINE: D9RDNM *****
C
C PURPOSE : TO EXTRACT COLLISIONAL DIELECTRONIC DATA FOR A
C           (TEMPERATURE, DENSITY) GRID FROM
C           EITHER PARTIAL (METASTABLE/PARENT RESOLVED) OR STANDARD
C           (UNRESOLVED) ISONUCLEAR MASTER FILES
C
C NOTE    : THE SOURCE DATA IS CONTAINED AS SEQUENTIAL DATASETS
C           WITH THE FOLLOWING NAMING CONVENTIONS:
C
C           (1) JETSHP.ACD<YR>#<EL>.<CODE>DATA
C           (2) JETSHP.SCD<YR>#<EL>.<CODE>DATA
C           (3) JETSHP.CCD<YR>#<EL>.<CODE>DATA
C           (4) JETSHP.PRB<YR>#<EL>.<FILT>.<CODE>DATA
C           (5) JETSHP.PRC<YR>#<EL>.<FILT>.<CODE>DATA
C           (6) JETSHP.QCD<YR>#<EL>.<CODE>DATA
C           (7) JETSHP.XCD<YR>#<EL>.<CODE>DATA
C           (8) JETSHP.PLT<YR>#<EL>.<CODE>DATA
C           (9) JETSHP.PLS<YR>#<EL>.<CODE>DATA
C
C           WHERE, <YR>   = TWO DIGIT YEAR NUMBER
C                   <EL>   = ONE OR TWO CHARACTER ELEMENT SYMBOL
C                   <CODE> = R       => PARTIAL DATA
C                           U       => PARTIAL DATA
C                           OMITTED => STANDARD DATA
C                   <FILT> = SIX CHARACTER POWER FILTER CODE
C
C           AND DATA OF CLASSES 6 AND 7 DO NOT EXIST FOR THE PARTIAL CASE.
C
C INPUT   : (C*120) DSNINC   = ISONUCLEAR MASTER FILE NAME - VERIFIED
C                   AND READY FOR DYNAMIC ALLOCATION.
C INPUT   : (L*4)  LPART    = .TRUE.  => PARTIAL (RESOLVED) MASTER DATA
C                   . FALSE. => UNSRESOLVED MASTER DATA
C INPUT   : (I*4)  IZ0      = NUCLEAR CHARGE
C INPUT   : (I*4)  NPART()  = METASTABLE PARTITION. I.E. NUMBER OF
C                   METASTABLES FROM CHARGE STATE IZ1MIN-1 TO
C                   IZ1MAX ON INPUT
C INPUT   : (I*4)  IPRTD    = REQUIRED PARENT INDEX
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C INPUT  : (I*4)  IGRDD      = REQUIRED GROUND INDEX
C INPUT  : (I*4)  ICLASS     = CLASS OF DATA (1 - 9 )
C INPUT  : (I*4)  IZ1        = REQUIRED ION CHARGE + 1
C INPUT  : (I*4)  ITMAX      = NUMBER OF DTEV() VALUES
C INPUT  : (I*4)  IDMAX      = NUMBER OF DDENS() VALUES
C INPUT  : (I*4)  ISDIMD     = MAXIMUM NUMBER OF (CHARGE, PARENT, GROUND)
C                               BLOCKS IN ISONUCLEAR MASTER FILES
C INPUT  : (I*4)  IZDIMD     = MAXIMUM NUMBER OF CHARGE STATES
C                               IN ISONUCLEAR MASTER FILES
C INPUT  : (I*4)  ITDIMD     = MAXIMUM NUMBER OF TEMP OR DENS VALUES IN
C                               ISONUCLEAR MASTER FILES
C INPUT  : (R*8)  DTEV()     = DLOG10(ELECTRON TEMPERATURES (EV))
C INPUT  : (R*8)  DDENS()    = DLOG10(ELECTRON DENSITIES (CM-3))
C
C OUTPUT  : (I*4)  IFAIL     = 0    IF ROUTINE SUCCESSFUL - DATA FOR THE
C                               REQUESTED YEAR USED.
C                               = 1    IF ROUTINE OPEN STATEMENT FAILED
C                               = 2    IF FILE EXISTS BUT REQUIRED DATA
C                               BLOCK DOES NOT
C OUTPUT  : (I*4)  ISMAXD     = NUMBER OF (CHARGE, PARENT, METASTABLE)
C                               BLOCKS IN SELECTED MASTER FILE
C OUTPUT  : (I*4)  IZMAXD     = NUMBER OF ZDATA() VALUES IN SELECTED
C                               MASTER FILE
C OUTPUT  : (I*4)  ITMAXD     = NUMBER OF DTEVD() VALUES IN SELECTED
C                               MASTER FILE
C OUTPUT  : (I*4)  IDMAXD     = NUMBER OF DDENSD() VALUES IN SELECTED
C                               MASTER FILE
C OUTPUT  : (I*4)  NPARTR()   = METASTABLE PARTITION. I.E. NUMBER OF
C                               METASTABLES FROM CHARGE STATE IZ1MIN-1 TO
C                               IZ1MAX FOUND IN MASTER FILE
C OUTPUT  : (R*8)  DTEVD()    = DLOG10(DATA ELECTRON TEMPERATURES (EV))
C                               IN SELECTED MASTER FILE
C OUTPUT  : (R*8)  DDENSD()   = DLOG10(DATA ELECTRON DENSITIES (CM-3))
C                               IN SELECTED MASTER FILE
C OUTPUT  : (R*8)  DRCOFD(,,) = DLOG10(DATA RATE COEFFICIENTS (CM-3/S))
C                               IN SELECTED MASTER FILE
C                               1ST DIM: (CHARGE,META,GRD) BLOCK INDEX
C                               2ND DIM: TEMPERATURE INDEX
C                               3RD DIM: DENSITY INDEX
C OUTPUT  : (R*8)  ZDATA()    = CHARGE + 1 FOR IONS IN SELECTED MASTER
C                               FILE
C                               1ST DIM: (CHARGE,META,GRD) BLOCK INDEX
C OUTPUT  : (R*8)  DRCOFI(,)  = INTERPOLATION OF DRCOFD(,,) FOR
C                               DTEV() & DDENS()
C
C PROGRAM: (C*80) DSNOLD     = FILE NAME USED IN PREVIOUS CALL
C           (C*80) CLINE     = GENERAL CHARACTER VARIABLE
C           (C*80) CTERM     = TERMINATOR LINE - '-' FILLED VARIABLE
C           (C*4)  CPATRN()  = PATTERN USED TO DETECT DATA CLASS
C           (I*4)  IZ0D      = NUCLEAR CHARGE READ FROM MASTER FILE
C           (I*4)  IZ1MIN    = MINIMUM CHARGE+1 READ FROM MASTER FILE
C           (I*4)  IZ1MAX    = MAXIMUM CHARGE+1 READ FROM MASTER FILE
C           (I*4)  IABT      = ABORT CODE

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C      (I*4)  INDSEL    = LOCATION OF (CHARGE, PRNT, GRND)
C
C      (I*4)  IZDAT     = CURRENT DATA BLOCK ION CHARGE +1
C      (I*4)  ISEL      = GENERAL INDEX
C      (I*4)  I          = GENERAL INDEX
C      (I*4)  IT         = GENERAL INDEX
C      (I*4)  ID         = GENERAL INDEX
C      (I*4)  IZCHK     = INDEX TO VERIFY DATA Z1 SET COMPLETE
C      (I*4)  IPRTR()   = PARENT INDICES IN DATA SET
C      (I*4)  IGRDR()   = GROUND INDICES IN DATA SET
C      (I*4)  LCK        = MUST BE GREATER THAN 'ITMAXD' & 'IDMAXD'
C
C      (R*8)  A()        = GENERAL ARRAY
C      (R*8)  DRCOF0(, ) = INTERPOLATION OF DRCOFD(, ,) W.R.T DTEV()
C      (L*8)  LEXIST     = TRUE --- FILE TO OPEN EXISTS ELSE NOT
C      (I*4)  L1         = PARAMETER = 1
C      (I*4)  IOPT       = DEFINES THE BOUNDARY DERIVATIVES FOR THE
C
C      (L*4)  LSETX      = .TRUE. => SET UP SPLINE PARAMETERS RELATING
C
C      .FALSE. => DO NOT SET UP SPLINE PARAMETERS
C
C      RELATING TO X-AXIS.
C
C      (I.E. THEY WERE SET IN A PREVIOUS
C
C      CALL )
C
C      (VALUE SET TO .FALSE. BY 'XXSPLN')
C      (R*8)  DY()       = SPLINE INTERPOLATED DERIVATIVES

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C ROUTINES:

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C      ROUTINE      SOURCE      BRIEF DESCRIPTION
C      -----
C      I4UNIT       ADAS        FETCH UNIT NUMBER FOR OUTPUT OF MESSAGES
C      I4FCTN       ADAS        CONVERT STRING TO INTEGER FORM
C
C      (R*8 ADAS FUNCTION - 'R8FUN1' ( X -> X ) )

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C AUTHOR : Alessandro Lanzafame

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- PUT UNDER SCCS CONTROL

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C-----
C      CHARACTER*120      DSNINC
C      INTEGER           ICLASS,      IDMAX,      IDMAXD,      IFAIL
C      INTEGER           IGRDD,      IPRTD,      ISDIMD,      ISMAXD
C      INTEGER           ITDIMD,      ITMAX,      ITMAXD,      IZ0
C      INTEGER           IZ1,      IZDIMD,      IZMAXD
C      INTEGER           NPART(IZDIMD),      NPARTR(IZDIMD)

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LOGICAL	LPART	
REAL*8	DDENS (IDMAX) ,	DDENSD (ITDIMD)
REAL*8	DRCOFD (ISDIMD, ITDIMD, ITDIMD)	
REAL*8	DRCOFI (ITMAX, IDMAX) ,	DTEV (ITMAX)
REAL*8	DTEVD (ITDIMD) ,	ZDATA (ISDIMD)