

## ADAS Subroutine h9ispf

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      SUBROUTINE H9ISPF( LPEND      , NDTIN      ,  
&                      IL         , NDTEM     ,  
&                      NV         , TSCEF    , TSCEF2   ,  
&                      ITRAN     , I1A      , I2A      ,  
&                      FTYPE     , TITLE    ,  
&                      ISTRN     , IFOUT    ,  
&                      MAXT      , TINE     ,  
&                      LFSEL     , LOSEL    ,  
&                      TOLVAL    ,  
&                      DTYPE     , KAP_VAL  , DRU_VAL  ,  
&                      ADF37     ,  
&                      )
```

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C-----  
C  
C ***** FORTRAN77 SUBROUTINE: H9ISPF *****  
C  
C PURPOSE: PIPE COMMUNICATIONS WITH IDL AND TO RETURN USER SELECTED  
C           OPTIONS AND VALUES.  
C  
C CALLING PROGRAM: ADAS809  
C  
C SUBROUTINE:  
C  
C I/O   : (L*4)   LPEND    = .TRUE.  => END ANALYSIS OF CURRENT DATA  
C                                     SETS  
C                                     = .FALSE. => CONTINUE PANALYSIS WITH CURRENT  
C                                     DATA SETS  
C  
C INPUT : (I*4)   NDTIN    = MAX. NUMBER OF TEMPERATURES ALLOWED  
C  
C INPUT : (I*4)   IL       = NUMBER OF ENERGY LEVELS  
C  
C INPUT : (I*4)   NDTEM    = INPUT DATA FILE: MAX. NO. OF TEMPERATURES  
C  
C INPUT : (I*4)   NV        = INPUT DATA FILE: NUMBER OF TEMPERATURES  
C  
C INPUT : (R*8)   TSCEF(,) = INPUT DATA FILE: ELECTRON TEMPERATURES, OR  
C                                     2/3 AVERAGE ENERGY OF A NUMERICAL  
C                                     DISTRIBUTION  
C                                     2ND DIMENSION: 1 => KELVIN (IFOUT=1)  
C                                               2 => EV (IFOUT=2)  
C                                               3 => REDUCED (IFOUT=3)  
C  
C INPUT : (R*8)   TSCEF2(,) = INPUT DATA FILE: MOST COMMON ELECTRON  
C                                     'TEMPERATURE' FOR A NUMERICAL DISTRIBUTION,  
C                                     OR A ZERO ARRAY IN ANALYTIC CASE  
C                                     2ND DIMENSION: 1 => KELVIN (IFOUT=1)  
C                                               2 => EV (IFOUT=2)  
C                                               3 => REDUCED (IFOUT=3)  
C  
C INPUT : (I*4)   ITRAN    = NUMBER OF ELECTRON IMPACT TRANSITIONS  
C
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C INPUT : (I*4) I1A() = ELECTRON IMPACT TRANSITION:
C LOWER ENERGY LEVEL INDEX
C
C INPUT : (I*4) I2A() = ELECTRON IMPACT TRANSITION:
C UPPER ENERGY LEVEL INDEX
C
C INPUT : (C*6) FTYPE = TYPE OF FILE SELECTED AS COMPARISON
C
C OUTPUT: (C*40) TITLE = USER ENTERED GENERAL TITLE FOR PROGRAM RUN
C
C OUTPUT: (I*4) ISTRN = SELECTED ELECTRON IMPACT TRANSITION INDEX
C
C OUTPUT: (I*4) IFOUT = 1 => INPUT TEMPERATURES IN KELVIN
C = 2 => INPUT TEMPERATURES IN EV
C = 3 => INPUT TEMPERATURES IN REDUCED FORM
C
C OUTPUT: (I*4) MAXT = NUMBER OF INPUT TEMPERATURES ( 1 -> 20)
C
C OUTPUT: (R*8) TINE() = ELECTRON TEMPERATURES (UNITS: SEE 'IFOUT')
C
C OUTPUT: (L*4) LFSEL = .TRUE. => CARRY OUT MINIMAX POLYNOMIAL
C FITTING
C = .FALSE. => - DO NOT DO THE ABOVE -
C
C OUTPUT: (R*8) TOLVAL = FRACTIONAL TOLERANCE FOR MINIMAX FIT
C (=0 IF MINIMAX FIT NOT SELECTED)
C
C OUTPUT: (L*4) LOSEL = .TRUE. => CALCULATE INTERPOLATED VALUES
C FOR OUTPUT.
C = .FALSE. => - DO NOT DO THE ABOVE -
C
C OUTPUT: (I*8) DTYPE = DISTRIBUTION TYPE SUCH THAT:
C 0 => MAXWELLIAN
C 1 => KAPPA
C 2 => NUMERICAL
C OUTPUT: (R*8) KAP_VAL = VALUE OF KAPPA
C
C
C
C (I*4) ILOGIC = RETURN VALUE FROM IDL WHICH IS USED TO
C REPRESENT A LOGICAL VARIABLE SINCE IDL
C DOES HAVE SUCH DATA TYPES.
C
C (I*4) I = GENERAL PURPOSE COUNTER
C
C (I*4) J = GENERAL PURPOSE COUNTER
C
C ROUTINES:
C ROUTINE SOURCE BRIEF DESCRIPTION
C -----
C XXFLSH IDL-ADAS CALLS 'FLUSH' TO CLEAR PIPE
C
C AUTHOR: HUGH SUMMERS (UNIV. OF STRATHCLYDE)
C

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C DATE: 30/11/01

C

C DATE: 07/04/05

C MODIFIED: ALLAN WHITEFORD

C - ADDED FORMAT STATEMENT TO READING FROM PIPE

C-----

CHARACTER*80	ADF37			
CHARACTER*6	FTYPE			
CHARACTER*40	TITLE			
INTEGER	DTYPE,	I1A (ITRAN),	I2A (ITRAN),	IFOUT
INTEGER	IL,	ISTRN,	ITRAN,	MAXT
INTEGER	NDTEM,	NDTIN,	NV	
LOGICAL	LFSEL,	LOSEL,	LPEND	
REAL*8	DRU_VAL,	KAP_VAL,	TINE (NDTIN),	TOLVAL
REAL*8	TSCEF (NDTEM, 3),		TSCEF2 (NDTEM, 3)	