

## ADAS Subroutine escape

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subroutine escape( ndprof, nddens, ndcyl , iz      ,
&                  tg      , j0      , mm      , x0      , x1      ,
&                  y0      , sum1   , sum2   , sum3   ,
&                  sum4   , alpha   , ics    , iden   ,
&                  sum    , iprofile , zlen, y, x, wid,
&                  dmult, modprof)

C
C-----
C
C ****&***** fortran77 subroutine: escape ****&*****
C
C original name: escape.bas
C             (developed by K. H. Behringer)
C
C version: 1.0
C
C purpose: computes escape factors.
C
C calling program: adas214
C
C input:
C
C (i*4)  nddens = parameter = max. number of points over ion profile
C (i*4)  ndcyl  = parameter = max. number of intervals for cylinder
C (i*4)  ndprof  = parameter = max. number of points over line profile
C
C (i*4)  ics     = 1 for slab geometry of plasma
C                 = 2 for cyl. geometry of plasma
C (i*4)  iden    = 1 for homogenous density distribution
C                 = 2 for linear density
C                 = 3 for parabolic density
C (i*4)  iprofile= 1 for Doppler line profile
C                 = 2 for Lorentzian line profile
C                 = 3 for Holtzman line profile
C                 = 10 for Double Doppler line profile
C                 = -2 for Voigt line profile
C                 = -3 for Doppler Holtzman line profile
C (i*4)  iz      = number of points over line profile
C (i*4)  j0      = number of density values for integration over
C                   cylinder/slab
C (i*4)  k0      = number of density values - must be enough for
C                   interpolation
C (i*4)  z0      = number of intervals for cylinder/slab
C
C (r*8)  allam   = Spectral absorption coefficient
C (r*8)  alpha    = neutral density*oscillator strength*length
C (r*8)  domeg   = delta omega in cylinder/slab
C (r*8)  l        = Length of plasma (0 to inf. on one side)
C (r*8)  mm       = atomic mass number
C (r*8)  omega    = solid angle in cylinder/slab
C (r*8)  r        = radius of cylinder
C (r*8)  sum1()   = sum assuming homogeneous density
C (r*8)  sum2()   = sum assuming linear density
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c      (r*8)  sum3() = sum assuming parabolic density
c      (r*8)  sum4() = used in sum for escape factor
c      (r*8)  sum()  = sum1/2/3, depending upon options chosen for density pr
c      (r*8)  tg     = neutral temperature (eV)
c      (r*8)  w      = Full Doppler width calculated for lambda=100nm
c      (r*8)  x0()   = Absorption coefficient at line centre/10
c      (r*8)  x1()   = Absorption coefficient at line centre
c      (r*8)  y      = Line profile intensity
c      (r*8)  y0     = Line profile intensity (?) at line centre
c      (r*8)  ff     = Used in calculation of escape factor
c      (r*8)  xl     = log of x1, used for interpolation
c      (r*8)  yl     = log of sum, used in interpolation
c      (r*8)  zlen   = variable used to select length along plasma integration
c
c
c routines:
c      routine      source      brief description
c      -----
c      faltung      ADAS       computes line profiles
c
c author: K. H. Behringer (IPF, University of Stuttgart)
c
c date: 30/01/94
c
c update: ???? brought up to date with latest Behringer code
c
c VERSION: 1.1 DATE: 18-06-98
c MODIFIED: STUART LOCH
c - CONVERTED TO FORTRAN
c
c VERSION: 1.2 DATE: 26-11-98
c MODIFIED: STUART LOCH
c - CHECKS IF LINE PROFILE IS INTEGRATED FAR ENOUGH,
c   IF NOT THE PROFILE IS RE-EVALUATED TO TWICE THE
c   PREVIOUS WIDTH, USING THE SAME STEP SIZE AS BEFORE.
c VERSION: 1.3                               DATE: 19-02-99
c MODIFIED: STUART LOCH
c           -DIMENSIONS OF X,Y, YT, YL AND MODPROF INCREASED FROM 1000
c             TO 3000, TO ACCOMODATE POSSIBLE INCREASES IN INTEGRATION
c             LIMIT OF LINE PROFILE
c VERSION: 1.4                               DATE: 24-09-99
c MODIFIED: STUART LOCH
c           -REMOVED EMPTY DO LOOP
c
c-----
c-----
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INTEGER	ICS,	IDEN,	IPROFILE,	IZ
INTEGER	J0,	NDCYL,	NDdens,	NDPROF
REAL*8	ALPHA,	DMULT,	MM	
REAL*8	MODPROF(K0,3000),		SUM(NDdens)	
REAL*8	SUM1(NDdens),		SUM2(NDdens)	
REAL*8	SUM3(NDdens),		SUM4(NDdens)	
REAL*8	TG,	WID,	X(3000)	
REAL*8	X0(NDdens),	X1(NDdens),	Y(3000),	Y0

REAL\*8

ZLEN