

APPENDIX L: TIPS AND OPERATIONAL HINTS THOUGHT USEFUL TO NEW USERS

This appendix comprises a collection of useful bits of information, extracted primarily from the source code, and thought by the code sponsor to be of potential utility to new users.

Usefull text extracted from source code

```
*!  
*      Real xareas = areasp (flowrate, xtclout, hrepo, flwcnst, poroi)  
*****  
***** Equivalent A_R_E_A of S_tuck P_ipe, repository removed *****  
***** - - - - - *****  
*****  
*  
* PURPOSE:      Calculates equivalent area of repository removed on stuck  
* pipe:  
*  
* AUTHORS:      Robert A. Cole  
*  
* UPDATED:      02-FEB-1996 --RAC: Deleted write for DBSPALL  
*                -JUL-1994 --RAC: Created  
*  
* CALLED BY:    CUTTINGS {main program}  
*  
* CALLS:        NONE  
*!  
*      The following hardwired constants are not presently used  
* DATA ACCNST/1.128E+13/  
* DATA AVOGAD/6.022E+23/  
*!  
*  
*      ...Source: Newton circa (1642-1727)  
* DATA GRAV/9.81/  
*  
*      Seconds per based on a Tropical Year of 31556926 sec  
* DATA YRPSEC/3.168876E-08/  
*  
*!  
*      Subroutine blowout (dia, hf, perm, porof, pg, partdia, xareas)  
*****  
***** Equivalent B_L_O_W O_U_T, calculates the equivalent area *****  
***** ejected by a blowout *****  
***** - - - - - *****  
*  
* PURPOSE: Calculates final blowout radius, and back pressure,  
* and returns equivalent waste area removed  
*  
*      Calculates entrance mach number for different  
* Pressure ratios for isothermal flow in pipes with  
* Friction....see eq 3.9 page 63 binder, advanced fluid  
* Mechanics vol1.  
*      Assumes two diameters equivalent to the annular areas
```

```

* adjacent to the collars and drillpipe. when mach10
* =mach2 then the value for p2 is correct and the
* entrance mach number= mach1=mach2.

Includes additional flow equations to calculate
Final blowout radius and back pressure
*corrected steady state flow out of waste
*annular area and waste surface relationship
*terminal velocity of spherical waste particle
    includes specific collar diameters
    modified to read from input file

* AUTHORS: Jerry W. Berglund & Robert A. Cole
* CALLED BY: CUTTINGS {main program}
* CALLS: NONE
*****!
* Subroutine brgcdb( maxc, maxq, iq, lq, q, c, ierrdb )
*****
* B_R_G_C_D_B set up read of a BRAGFLO CDB file
* ****
* PURPOSE: Process dynamic memory allocation for input
* BRAGFLOW CDB file etc...
* AUTHOR: Robert A. Cole
* CREATED: 02-FEB-1996 --RAC: Created from a copy of PROCDB
* CALLED BY: CUTTINGS {main program}
* CALLS: DBISIZES
*         DBITITLE
*         ADDMEM
*         DYNMEM
*         ZERODA
*         DYNMEMC
*         READBRAG
*
* ARGUMENTS:
* ENTRY/
*   --through subroutine call
*     MAXC   = Total no. of CHAR*1 words allocated to dynamic storage
*     MAXQ   = Total no. of words allocated to dynamic storage
*     IQ     = INTEGER Dynamic Q-array (equivalenced to Q)
*     LQ     = <LOGICAL> Dynamic Q-array (equivalenced to Q)
*     Q      = REAL Dynamic memory base array
*     C      = CHARACTER Dynamic memory base array
*     IERRDB = INTEGER error code
*
* LOCAL/
*     MORMEMC= No. additional words to be allocated to C array
*     MORMEMR= No. additoinal words to be allocated to Q array

```

```
*  
* EXIT/  
  
*****  
*****!  
* Subroutine cal_bra  
*****  
***** CALIBRATION checks out model2 with experiments *****  
***** --- --- *****  
  
*  
* PURPOSE: To calibrate model2 with experimental data or WIPP data  
* also validate's model2 over possible WIPP variable  
* ranges  
* This subroutine gets activated by defining the LOGICAL  
* variable CUSP$CHECK_OUT to TRUE. This subroutine  
* expects to find the first keyword on the TXT1 input file  
* either CALIBRATE, EXPERIMENT, or VALIDATE  
  
*  
* AUTHOR: Robert A. Cole  
* CREATED: 02-FEB-1996 --RAC:  
* CALLED BY: CUTTINGS {main program}  
* CALLS: CAL_INP  
* SOLID_OUT  
  
* LOCAL/  
*****  
*!  
* Subroutine cal_inp (test_funct)  
*****  
***** CALibration INPut Reads input for calibration *****  
***** --- --- *****  
  
*  
* PURPOSE: To read input required for CALIBRATION, VERIFICATION,  
* & EXPERIMENT  
*  
* AUTHOR: Robert A. Cole  
* CREATED: 02-FEB-1996 --RAC:  
* CALLED BY: CAL_BRA  
* CALLS: NEXT0  
* NEXTW  
* NEXT_R  
* NEXT_I  
  
*****  
*!  
* Subroutine drill( iwasflg, ihit, xtint, dbdiam, tunits, xnewarea)  
*****  
***** D_R_I_L_L_ing through radwaste repository module *****
```

```
***** - - - - - *****  
*  
* PURPOSE: Calculates new borehole area based on drilling and  
* penetrating either:  
*  
* (1) WIPP repository containing CH or RH type  
*      transuranic waste, or  
* (2) Hypothetical INEL/WINCO repository containing  
*      HLW/SNF waste.  
*  
* AUTHORS: Jerry W. Berglund  
*          Jonathan S. Rath  
*  
* CALLED BY: CUTTINGS {main program}  
*  
* CALLS: QAABORT  
*        INTERFACE  
*        FRICTION {REAL function}  
*        F11 {REAL function}  
*        F12 {REAL function}  
*        F13 {REAL function}  
*        F21 {REAL function}  
*        F22 {REAL function}  
*        F23 {REAL function}  
*        F31 {REAL function}  
*        F32 {REAL function}  
*        F33 {REAL function}  
*        EF1 {REAL function}  
*        EF2 {REAL function}  
*        EF3 {REAL function}  
*  
* ARGUMENTS:  
* ENTRY/  
* --through subroutine call  
* IWASFLG= WIPP TRU waste type flag [INTEGER]  
*       = 1; Contact Handled (CH) TRU waste  
*       = 2; Remote Handled (RH) TRU waste  
* IHIT   = Intrusion (or hit) index [INTEGER]  
* XTINT  = Time of this intrusion (or hit) [REAL]  
* DBDIAM = Drill Bit Diameter for this intrusion (or hit);  
*           (Outside diameter of drillstring annulus) [REAL]  
* TUNITS = Units of time (i.e. seconds, years, etc...) [CHAR*8]  
*  
*****  
*! Integer function fgrad (input)  
*****  
*****  
*  
* PURPOSE: To determine the number of unique radionuclides  
*  
* AUTHOR: Robert A. Cole  
*  
* CREATED: 02-FEB-1996 --RAC:  
*  
* CALLED BY: PROCIN  
*  
*****  
*****  
*!  
* Real function gasers (xdbdia, flowrate, drate, depth, hrepo,
```

```
*                               flwcnst, poroi)
*****
*****   Equivalent G_A_S_ E_R_o_S_i_o_n,      repository removed *****
*****   - - - - - *****
*
* PURPOSE:      Calculates equivalent area of repositoy removed because
*               gas erosion:
*
* AUTHORS:      Robert A. Cole
*
* UPDATED:      02-FEB-1996 --RAC: Deletet write for DBSPALL
*                 xx-JUL-1994 --RAC: Created
*
* CALLED BY:    CUTTINGS {main program}
*
* CALLS:        NONE
*
* ARGUMENTS:
* ENTRY/
*   --through function arguments
*     DHOLE = Radius of borehole
*     FLOWRATE=Recommended (from handbook formula) drilling fluid
*               (mud) flowrate based on drillstring diameter [REAL].
*     TCLOUT = Cleanout time for stuck pipe (s)
*     HREPO = Height of repository (m)
*     POROI = Initial porosity
*
* EXIT/
*   --through fucntion value
*   GASERS = Equivalent area of repository removed (REAL)
*
*****
*!
* Subroutine gencut( iwasflg, diammod, aintc, tint, nnic,
* $                      radchn, nucnam, MAPSAV, idmtr1, halfy,
* $                      awt, actcnv, xmoles, eparel, cipmsq,
* $                      cizero, sumci, radhis, xnorm, xinvt0,
* $                      cizersv, xporo, xprgs, xprbr,
* $                      xsatb, mapinv, c, q, kulp )
*****
*****
*****   G_E_N_erate amount of C_U_T_tings module *****
*****   - - - - - *****
*
* PURPOSE:      Determines the drill cuttings (in Curies) brought to
*               the surface. Using the intrusion borehole diameter
*               and intrusion time, the routine determines the amount
*               of cuttings, and ages the cuttings to the intrusion
*               time.
*
* AUTHORS:      Harold J. Iuzzolino
*                 Jonathan S. Rath
*
* MODIFIED:     Robetr A. Cole
*
* CALLED BY:    CUTTINGS {main program}
*
* CALLS:        WRTANA
```

```
*          QAMAXERR
*
*          WRPLFIL
*
* FUNCTIONS: QUANKG
*             GINDX
*             GJNDX
*
* ARGUMENTS:
* ENTRY/
* --through subroutine call
* IWASFLG= radWASTE type FLaG (INTEGER)
* = 1; WIPP Contact Handled (CH) TRU waste
* = 2; WIPP Remote Handled (RH) TRU waste
* = 3; INEL/WINCO HLW/SNF
* DIAMMOD= Outside diameter of drillstring annulus, for each
* intrusion or hit, array (REAL).
* AINTC = Borehole area, as computed by DRILL routine array
* TINT = Intrusion times of each hit, in SECONDS array (REAL)
* NNIC = No. of Nuclides in each decay chain, array (INTEGER)
* RADCHN = Names of radioisotopes in each chain, array (CHAR*8)
* NUCNAM = Names of radioisotopes used by CUTTINGS (CHARACTER*8)
* IDMTRL = Names of Generic Radioisotope Data Base (GRDB)
*           entries, array (CHARACTER*8)
* HALFY = Half-Life of each radioisotope in GRDB [years],
*           array (REAL)
* AWT = Atomic WeighT of each radioisotope in GRDB [Kg/Mole],
*           array (REAL)
* ACTCNV = ACTivity CoNVersion of each radioisotope in
*           GRDB [Ci/Kg], array (REAL)
* XMOLES = Dummy array used (REAL)
* CIMPSQ = CURIES per SQ. METER at time zero (REAL)
* CIZERO = CURIES of each radioisotope, at time zero (REAL)
* EPAREL = EPA normalization factor for each radioisotope in
*           GRDB, array (REAL)
* MAPSAV = Map of unique names of radioisotopes to be saved
*           MAPSAV(I) .GT. 0 save
*           MAPSAV(I) .EQ. 0 do not save
*           to output CAMDAT and debug/diagnostics file (CHAR).
* CIZERSV= Values of radioisotopes to be written to the
*           output CAMDAT and debug/diagnostics file (REAL).
* C      = Dynamic base array (CHARACTER*1)
* Q      = Dynamic base array (REAL)
*
* LOCAL/
*
* EXIT/
* --through subroutine call
* RADHIS = Radioisotope Inventory (Curies) History array [REAL]
* SUMCI = Sum of all hits (in Curies) for each radioisotope
*           array (REAL)
* XNORM = Normalized radioisotope relaese array (REAL)
* XINVT0 = Inventory at time=0, array [REAL]
*
*!
*          Types of data sets
* -----
*
* INPUT/
* BINAY: in units 'IDB'
* ASCII: in units 'INPUTP', 'INPUTF', 'INPSDB'
```

```
*      BINARY: in UNITS 'IBRAG'  
*  
*      OUTPUT/  
*      BINARY: in units 'ODB', 'DB_BIN'  
*      ASCII:  in units 'ICTRN', 'IHISTO', 'IPLT1', 'IPLT2', 'IPLT3',  
*                  'NHISTO', 'NOUTFL', 'NVERFY', 'DB_ASC'  
*  
*                      Files used  
*-----  
*  
* $DEFINE logical-name file-path-name  
*  
* logical file name   unit   description  
*  
* 'CUSP_INP$TXT0'      1      Input text file (REQUIRED) (Radionuclide  
*                               input, Model & Site specific properties)  
*                               or if CUSP_INP$TXT0 is not defined  
* 'CUSP_INP$BINO'       1      Input binary file (REQUIRED) (Radionuclide  
*                               input, Model & Site specific properties)  
*                               or  
* 'CUSP_INP$TXT1'       2      Input text file (REQUIRED) (Drilling  
*                               properties and to control testing inputs  
* 'CUSP_INP$BRAGCDB'    7      Input BRAGFLOW CDB file (OPTIONAL)  
* 'CUSP_OUT$DBG'        8      Output debug File  
* 'CUSP_INP$CDB'        10     Input Computational Data Base (REQUIRED)  
* 'CUSP_OUT$CDB'        11     Output CDB File (REQUIRED)  
*  
* The following files are (OPTIONAL):  
* 'CUSP_OUT$NVERFY'     9      Output verify File  
* 'CUSP_OUT$IHISTO'     20     Output Radionuclide Inventory History  
* 'CUSP_OUT$NHISTO'     21     Output Normalized History file  
* 'CUSP_OUT$PLT1'        22     Output radioisotope decay plot file (long)  
* 'CUSP_OUT$PLT2'        23     Output radioisotope decay plot file (medium)  
* 'CUSP_OUT$PLT2'        24     Output radioisotope decay plot file (short)  
* 'CUSP_OUT$ICTRN'      25     Output CCDFPERM transfer file  
*  
*  
* The following files are required for  
* CUSP$PRE_CUTTINGS:  
*  
* 'CUSP_INP$CDB'        10     Input Computational Data Base (REQUIRED)  
* 'CUSP_INP$SDB'         4      Secondary Data Base Flatfile  
*                               (REQUIRED pre_cuttings)  
* 'CUSP_INP$TXT1'        2      Input text file (REQUIRED) (Drilling  
*                               properties and to control testing inputs  
* 'CUSP_OUT$DBG'         8      Output debug File  
* 'CUSP_OUT$DB_ASC'      12     Output ascii SDB file  
* 'CUSP_OUT$DB_BIN'      13     Output binary SDB file  
*  
*****  
*!  
*****          PARBINO          *****  
*  
* PURPOSE:      reads a CUTTINGS binary file for parameters inputs  ****  
*  
* AUTHOR:       Robert A. Cole  ****  
*  
* UPDATED:      02-FEB-1996 --RAC: First Ed.  ****
```

```

* CALLED BY: PROBIN
*
* CALLS: QAABORT
*****
*!
* Subroutine partxt0
*****
*      P_A_R_a_m_e_t_e_r_s SCAN of T_e_X_T file
*      - - -           - - -
*
* PURPOSE: reads a CUTTINGS input text file for parameters inputs
*
* AUTHOR: Robert A. Cole
*
* UPDATED: 02-FEB-1996 --RAC: First Ed.
*
* CALLED BY: PROCIN
*             MAIN (CHK_OUT)
*
* CALLS: NEXTO
*        NEXT_I
*        NEXT_R
*
* ARGUMENTS:
* ENTRY/
* --through subroutine call (none)
*****
*!
* Subroutine prepro( fidb, ftxt0, fbin0, ftxt1, fodb, fbrg )
*
* PURPOSE: Interacts with DEC DCL stream to determine input &
*          outputs file and logical settings
*****
*!
* Subroutine probin( maxq,    maxc,    iq,    lq,    q,
* $           c,       komat,   outmat, iwasflg, tunits,
* $           nqarec, c(lgainfo) )
*****
* PURPOSE: Processes the CUTTINGS binary file (BIN0) information
*          & the CUTTINGS ASCII file (TXT1) information
*
* AUTHOR: Robert A. Cole
* CALLED BY: CUTTINGS {main program}
*
* CALLS: SCABINO
*        SCATXT1
*        PARBINO
*        ADDMEM2
*        DYNMEM
*        DYNMEMC
*        ZERODA
*        REABINO
*        REATXT1
*        REAEPA

```

```
*  
* ARGUMENTS:  
* ENTRY/  
* --through subroutine call  
*   MAXQ    = Total no. of REAL words allocated (INTEGER)  
*   MAXC    = Total no. of CHARACTER*1 words allocated (INTEGER)  
*   IQ      = INTEGER Dynamic Q-array (equivalenced to Q)  
*   LQ      = <LOGICAL> Dynamic Q-array (equivalenced to Q)  
*   Q       = REAL Dynamic memory base array  
*   C       = CHARACTER Dynamic memory base array  
  
* EXIT/  
* --through subroutine call  
*   MAXQ    = Total no. of REAL words allocated (INTEGER)  
*   MAXC    = Total no. of CHARACTER*1 words allocated (INTEGER)  
*   KOMAT   = CAMDAT material index to which CUTTINGS output  
*             variables will be written as material property values  
*             (INTEGER)  
*   OUTMAT  = CAMDAT material name to which CUTTINGS output  
*             variables will be written as material property values  
*             (CHARACTER*8)  
*   IWASFLG= radWASTE type FLaG (INTEGER)  
*             = 1; WIPP Contact Handled (CH) TRU waste  
*             = 2; WIPP Remote Handled (RH) TRU waste  
*             = 3; INEL/WINCO HLW/SNF waste(s)  
*   TUNITS  = Units associeated with the time of intrusions (CHAR*8)  
  
*****  
*!  
* Subroutine prosdb( maxc, maxq, iq, lq, q, c , iwasflg, qainfo )  
  
*****  
*  
* PURPOSE:      Read a flat ascii flat to extract properties &  
*               data from the INGRESS data base  
*  
* AUTHOR:       Robert A. Cole  
*  
* UPDATED:      02-FEB-1996 --RAC: First Ed.  
*  
* CALLED BY:    PROSDB used by PRE_CUSP  
*  
* CALLS:        SCATXT0  
*                 SDBQUERY  
*                 REASDB  
*                 PROCIN  
*                 CMDREAD  
*                 CHCASE  
*                 BLANK  
*                 SORT  
*                 QAABORT  
*  
*  
*  
*!  
* REABINO  
*  
* PURPOSE:      Reads a CUTTINGS binary file
```

```
* AUTHOR:      Rovert A. Cole           ****
*           ****
* UPDATED:     02-FEB-1996 --RAC: First Ed.   ****
*           ****
* CALLED BY:    PROBIN                 ****
*           ****
* CALLS:       QABORT                ****
*           ****
*****          ****
*!
*     Subroutine readbrag (diammod, poro, presgas,
*     &                      presbrn, SATGASX, nhits, tint,
*     &                      ele_n, ele_id, ele_blk, att_n,
*     &                      gridvol, prop_n, lvar_b, ele_prp,
*     &                      ele_hprp, ilog, MULTP, ierrdb)
*****          ****
*
* PURPOSE:      To extract the BRAD CDB file information
*
* Arguments to this subroutine
*
* MULTP      = Number of multiple intrusions into BRAGCDB file at
*               same intrusion time
* DIAMMOD   = the drill diameter
* PORO       <= the porosity in the repository
* PRESGAS    <= the pressure of the gas in the repository
* PRESBRN    <= the pressure of the brine in the repository
* SATGASX   <= the saturation of the brine in the repository
* NHITS      => number of intrusions
* TINT        => the intrusion times
* ELE_N      <= the element block names
* ELE_ID     <= the element block id's
* ELE_BLK    <= the number of elements per block
* ATT_N      <= the attributes names
* GRIDVOL   <= the array for the grid volumes
* PROP_N     <= the property names
* LVAR_B     <= logical property table
* ELE_PRP    <= element property values
* ELE_HPRP   <= element history property values
* IERRDB    <= the error flag returned from call to "DBI---"
*
*****          ****
*!
*****          ****
****          ****
* PURPOSE:      Extract a property value from an INGRES data base   ****
*           ****
* AUTHOR:      Robert A. Cole           ****
*           ****
* UPDATED:     02-FEB-1996 --RAC: First Ed.   ****
*           ****
* CALLED BY:    PROSDB used by PRE_CUSP   ****
*           ****
* CALLS:       SDBGETL                ****
*               QABORT                 ****
*           ****
*****          ****
*!
*****          ****
```

```
*****          SCABINO          *****
* PURPOSE:      Reads a binary parameter/data file for
*                 sizing parameters
*
* AUTHOR:       Robert A. Cole
*
* UPDATED:      02-FeB-1996 --RAC: First Ed.
*
* CALLED BY:    PROCIN
*
* CALLS:
*
*****
```

Usefull text extracted from include code

```
*!
* /BLOWPROP/ (from FORTRAN statement: include 'CUSP_BLOWPROP.INC')
*
* PURPOSE:      Provide common space for model 1 & model 2 properties
*                 & WIPP site specific properties
*
* AUTHOR:       Robert A. Cole
*
* UPDATED:      02-FEB-1996 --RAC: Added model2
*
* BRAGCDB = <LOGICAL> <default true>
*             true   <=> reads BRAGFLOW "CDB", WIPP run
*             false  <=> check out runs, test problems
*
* CHK_OUT = <LOGICAL> <default false>
*            false  <=> WIPP run
*            true   <=> calibration for experiments
*
* MODTYPE = Model type:
*           MODEL1
*           MODEL2
*
* Model 1 & 2 specific inputs:
*
* The following values are input through the code via
* the logical input file CUSP_INP$TXT0:
*
* PR_MAX = The maximum pressure allowed by model
* PR_MIN = The minimum pressure allowed by model
* PE_MAX = The maximum permeability allowed by model
* PE_MIN = The minimum permeability allowed by model
* PCUT   = Pressure (Pa) which defines the border between the
*           erosion phase and the blowout phase
* PGAS   = Pressure (Pa) which defines the border between the gas
*           erosion phase and the stuck pipe phase
* KCRIT  = Permeability which defines the phase border the blowout
*           phase and the gas-erosion/stuck pipe phase
*
* Model 2 specific inputs:
*
* DELTR  = Delta radius for derivative in stress calculation
* FGE    = Gravity    effectiveness factor
* FSE    = Strength   effectiveness factor
```

```
*      FCE      = Cementation effectiveness factor
*      CEMSIG   = Cementation Stress
*      ITER     = Number of iteration to solve for reynolds number
*      SUFTEN   = Surface Tension
*
*      Hydrogen specific constants (or Air for Experiments)
*
*      KGAS     = Ration of specific heats (unitless)
*      RGAS     = Gas Constant
*      VISC     = Hydrogen viscosity
*      CHY*    = Hydrogen sound speed (m/s), calculated from:
*                  CHY = SQRT (KGAS * RGAS * TREPO)
*
*      Repository/site specific inputs:
*
*      TREPO   = Temperature of repository (K)
*      RH01*   = Gas Density in WIPP room (kg/m**3), calcutated from:
*                  RHO1 = PG / TREPO / RGAS
*      PSUF    = Surface presure (Pa)
*      DEPTH   = Distance from repository depth to depth where casing
*                  is set (m)
*      HREPO   = Height of repository at burial time (4m)
*      RPANEL  = The equivalent radius of 1 panel (910.0m**2)
*      ROOM    = The equivalent radius of 1 room (11,640m**2)
*      RHOS    = Waste particle density (kg/m**3)
*      INPORO  = Initial repository porosity
*                  Use in CALIBRATION runs
*                  Production run uses value from BRAG CDB file
*
*      Drilling specific inputs:
*
*      L1      = Collar Length (m)
*      L2      = Drill pipe length (m)
*      COLDIA  = Collar diameter (m)
*      PIPED   = Drill pipe diameter (m)
*      FRICTF  = Friction factor (unitless)
*      FLWCNST= Percent volume of material that is carried by Drilling
*                  mud (unitless)
*
*
*      APORO   = A constant in equation to determine permeability as a
*                  function of porosity
*      NPORO   = N constant in equation to determine permeability as a
*                  function of porosity
*
*                  permeability = APORO * porosity ** NPORO
*
*****
***** ! *****
*      /BRGPAR/ ($include 'CUSP_BRGPAR.INC')
*      HEADER_B = Character description of problem, etc...
*      IMAX_B   = No. of I-lines in grid
*      JMAX_B   = No. of J-lines in grid
*      KMAX_B   = No. of K-lines in grid
*      LESSEL_B = List of Element Side sets concatenated Element List
*      LESSNL_B = List of Element Side sets concatenated Nodes associate
*                  with each element side set Node List
*      LNPSNL_B = List of Nodal Point set concatenated Nodes List
*      MAXNENO_B = Max. no. of nodes of any element
```

```
*      NDIM_B      = Number of physical DIMensions
*      NDBCAM_B    = CAMDAT versoin number
*      NELBLK_B    = Number of ELEMENT BLocks in mesh
*      NELX_B      = No. of elements along X-axis of mesh
*      NELY_B      = No. of elements along Y-axis of mesh
*      NELZ_B      = No. of elements along Z-axis of mesh
*      NQAREC_B    = Number of QA records
*      NSTEP_B     = No. of time steps in input CAMDAT (INTEGER)
*      NUMEL_B     = NUmber of ELeMents in mesh
*      NUMNOD_B    = NUMBER of NODes in mesh
*      NUMNPS_B    = No. of Nodal Point Sets
*      NUMESS_B    = No. of Element Side Sets
*      NUQATR_B    = No. of Unique ATtRIBUTE names
*      NUQPRO_B    = No. of Unique PRoperty names (old)
*      NUQPRP_B    = No. of Unique PRoPerty names (new)
*      NVAREL_B    = No. of ELEMENT type ANALYSIS variables
*      NXINFO_B    = No. of information records
*
*****  

*****!  

*****  

*      /CMM_INP/ ($include 'CUSP_CAL_INP.INC')
*      FSEMIN = Minimum FSE
*      DELFSE = Delta FSE
*      IFSE = No of increment of FSE
*      FGEMIN = Minimum FGE
*      DELFGE = Delta FGE
*      JFGE = No of increment of FGE
*      FCEMIN = Minimum FCE
*      DELFCE = Delta FCE
*      KFCE = No of increment of FCE
*      SATMIN = Minimum SAT
*      DELSAT = Delta SAT
*      ISA = No of increment of SAT
*      PGMIN = Minimum PG
*      DELPG = Delta PG
*      IPG = No of increment of PG
*      PORMIN = Minimum POR
*      DELPOR = Delta POR
*      DIA = Diameter
*      POROST = Porosity
*      IPOR = No of increment of POR
*      PARTMIN = Minimum diameter
*      PARTMAX = Maximum diameter
*      PARTMUL = Particle multiplier
*      ITMAX = Maximum number of trys for search of FGE, FSE, & FCE
*      DELERR = Delta error on convergence
*****
*!  

*****  

*      /FILEUN/ (from FORTRAN statement: include 'CUSP_FILEUN.INC')
*      ICTRN = Device no. of output CCDFPERM transfer file
*      IDB = Device no. of input CAMDAT file
*      IDBSCR = Device no. of CAMDAT scratch file
*      IHISTO = Device no. of Radioisotope Inventory History file
*      INPUTP = Device no. of input text file or binary file:
*                  Model specific parameters
*                  Site specific parameters
```

```
•          Radioisotope infomation
•
•      INPUTF = Device no. of input text file
•      INPSDB = Device no. of secondary data base SDB
•      IPLT1  = Device no. of plotting file #1 ("long-lived")
•      IPLT2  = Device no. of plotting file #2 ("medium-lived")
•      IPLT3  = Device no. of plotting file #3 ("short-lived")
•      NHISTO = Device no. of Normalized Release History file
•      NOUTFL = Device no. of output diagnostics/debug file
•      NVERIFY = Device no. of output verify file
•      ODB    = Device no. of output CAMDAT file
•      IBRAG  = Device no. of Bragflow "CDB" file
•      DB_ASC = Device no. of output SDB ascii file
•      DB_BIN = Device no. of output SDB binary file
•      ISCR3  = Device no. of scratch storage #3
*****
*!
***** /MISCDAT/ (from FORTRAN statement: include 'CUSP_MISCDAT.INC')
*     Intrusion times are in years
•      NCHAINS = No. of unique radioisotope decay chains (INTEGER)
•      NGRAD   = No. of generic radioisotopes or radionuclides (INTEGER)
•      NHITS   = No. of hits or intrusions (INTEGER)
•      NNUC    = No. of radionuclides used in CUTTINGS (INTEGER)
•      NORIAC  = No. Of Radioisotopes In All Chains (INTEGER)
•      NRADPR  = No. of RADionuclides to be PRinted (saved to
*          output CAMDAT and written to debug/diagnostics
*          file) [INTEGER]
•      SUMCUR  = Sum of all Radioisotope inventory, in Curies (REAL)
•      MULTP   = Number of multiple hits at the same intrusion time
*          Default is 1 if there is no BRAGCDB file
*****
*!
***** /RDBRAGx/ (from FORTRAN statement: include 'CUSP_RDBRAG.INC')
*
•      NULTP   = Maximum number of hits into brag calculation
*!
***** /TOOLS/ (from FORTRAN statement: include 'CUSP_TOOLS.INC')
•      DEBUG   = <LOGICAL>for debugging dynamic memory allocation
•      DBRAG   = <LOGICAL>for debugging reading bragflow CDB file
•      NOECHO  = <LOGICAL>for no echoing of input to the DBG file
•      VERIFY  = <LOGICAL> for writing a file for verifying purposes
•      PRE_CUSP = <LOGICAL> <default false>
*          false  <=> Calculation
*          true   <=> Pre_cusp, read SDB data base
•      PLODEC  = <LOGICAL>for writing long, medium, and short half-life
*          radioisotope decay history plot files.
•      WRTCTRN= <LOGICAL> for writing output CCDFPERM transfer file
•      WRTHIS  = <LOGICAL> for writing RAD WASTE-radioisotope inventory
*          history table.
•      WRTNHIS= <LOGICAL> for writing Normalized Radioisotope
*          Inventory History
```

Strucure of where how TXT0 file is read in:

SUBROUTINE PARTXT0

6720

| | Absolute line number ^^^^ |
|---|---------------------------|
| IF (CMD .NE. 'MODEL_DATA') GO TO 100 | 74 |
| IF (CMD .EQ. 'RADIONUCLIDE_DATA') GO TO 900 | 75 |
| IF (CMD .EQ. 'EOF') | GO TO 900 |
| IF (CMD .EQ. 'PR_MAX') THEN | 81 |
| IF (CMD .EQ. 'PR_MIN') THEN | 85 |
| IF (CMD .EQ. 'PE_MAX') THEN | 92 |
| IF (CMD .EQ. 'PE_MIN') THEN | 99 |
| IF (CMD .EQ. 'PCUT') THEN | 106 |
| IF (CMD .EQ. 'PGAS') THEN | 113 |
| IF (CMD .EQ. 'KCRIT') THEN | 120 |
| IF (CMD .EQ. 'DELTR') THEN | 127 |
| IF (CMD .EQ. 'FGE') THEN | 134 |
| IF (CMD .EQ. 'FSE') THEN | 141 |
| IF (CMD .EQ. 'FCE') THEN | 148 |
| IF (CMD .EQ. 'CEMENT') THEN | 155 |
| IF (CMD .EQ. 'ITER') THEN | 162 |
| IF (CMD .EQ. 'SUFLEN') THEN | 169 |
| IF (CMD .EQ. 'KGAS') THEN | 176 |
| IF (CMD .EQ. 'RGAS') THEN | 183 |
| IF (CMD .EQ. 'VISC') THEN | 190 |
| IF (CMD .EQ. 'TREPO') THEN | 197 |
| IF (CMD .EQ. 'PSUF') THEN | 204 |
| IF (CMD .EQ. 'DEPTH') THEN | 211 |
| IF (CMD .EQ. 'HREPO') THEN | 218 |
| IF (CMD .EQ. 'RPANEL') THEN | 225 |
| IF (CMD .EQ. 'ROOM') THEN | 232 |
| IF (CMD .EQ. 'RHOS') THEN | 239 |
| IF (CMD .EQ. 'L1') THEN | 246 |
| IF (CMD .EQ. 'L2') THEN | 253 |
| IF (CMD .EQ. 'COLDIA') THEN | 260 |
| IF (CMD .EQ. 'PIPED') THEN | 267 |
| IF (CMD .EQ. 'ROUGH') THEN | 274 |
| IF (CMD .EQ. 'FLWCNST') THEN | 281 |
| IF (CMD .EQ. 'APORO') THEN | 288 |
| IF (CMD .EQ. 'NPORO') THEN | 295 |
| IF (CMD .EQ. 'MODEL1') THEN | 302 |
| IF (CMD .EQ. 'MODEL2') THEN | 309 |
| | 317 |

Relative line number ^^^^

| | Absolute line number ^^^^ |
|---|---------------------------|
| SUBROUTINE REATXT0(NNIC, RADCHN, IDMTRL, NUCNAM, | 9513 |
| IF (CMD .NE. 'RADIONUCLIDE_DATA') GO TO 100 | 122 |
| IF (CMD .EQ. 'EOF') GO TO 500 | 129 |
| IF (CMD(1:5) .EQ. 'CHAIN') THEN | 131 |
| IF (CMD .EQ. 'SAVE') THEN | 162 |
| IF (CMD .EQ. '<TABLE_INPUTS') THEN | 182 |
| IF (CMD .EQ. 'END_TABLES') GO TO 110 | 188 |
| IF (CMD(1:1) .EQ. '**') THEN | 237 |
| IF (CMD .EQ. 'AWT') THEN | 254 |
| IF (CMD .EQ. 'HALFS') THEN | 261 |
| IF (CMD .EQ. 'ACTCNV') THEN | 268 |
| IF (CMD .EQ. 'EPAREL') THEN | 275 |
| IF (CMD .EQ. 'INVEN') THEN | 282 |

Relative line number ^^^^

| | |
|---|---------------------------|
| SUBROUTINE SCATXT0(IWASFLG, NAMELB, KOMAT, OUTMAT, IDMTRL, | 11038 |
| | Absolute line number ^^^^ |
| IF (CMD .NE. 'RADIONUCLIDE_DATA') GO TO 100 | 133 |
| IF (CMD .EQ. 'EOF') GO TO 900 | 139 |
| IF (CMD(1:5) .EQ. 'CHAIN') THEN | 144 |
| IF (CMD(1:4) .EQ. 'SAVE') THEN | 177 |
| IF (CMD(1:6) .EQ. 'INV_AR') THEN | 203 |
| IF (CMD(1:6) .EQ. 'RHW_AR') THEN | 221 |
| IF (CMD(1:3) .EQ. 'WUF') THEN | 239 |
| IF (CMD(1:7) .EQ. 'OUT_MAT') THEN | 257 |
| IF (CMD(1:6) .EQ. 'RADWAS') THEN | 295 |
| IF (CMD(1:3) .EQ. 'CON' .OR. | 301 |
| & CMD(1:2) .EQ. 'CH') THEN | ^ |
| IF (CMD(1:3) .EQ. 'REM' .OR. | 306 |
| & CMD(1:2) .EQ. 'RH') THEN | ^ |
| IF (CMD(1:3) .EQ. 'INE') THEN | 311 |
| IF (CMD(1:8) .EQ. 'REP_NAME') THEN | 325 |
| IF (CMD(1:8) .EQ. 'REP_GEOL') THEN | 346 |
| IF (CMD(1:6) .EQ. 'DISPOS') THEN | 365 |
| IF (CMD(1:4) .EQ. 'PACK') THEN | 385 |
| IF (CMD .EQ. '<TABLE_INPUTS') THEN | 405 |
| IF (CMD .EQ. 'END_TABLES') GO TO 120 | 413 |
| IF (CMD(1:1) .EQ. '*') THEN | 415 |
| | Relative line number ^^^^ |

Structure of where how TXT1 file is read in:

| | |
|---|---------------------------|
| SUBROUTINE REATXT1(IASPRP, NAMELB, NMATPR, XMATPR, TINT, | 10016 |
| | Absolute line number ^^^^ |
| IF (CMD .EQ. 'EOF') GO TO 666 | 108 |
| IF (CMD .EQ. 'UNITS') THEN | 113 |
| IF (CMD .EQ. 'SECONDS') THEN | 119 |
| IF (CMD .EQ. 'TINTR') THEN | 135 |
| IF (CMD .EQ. 'DIAMMOD') THEN | 165 |
| IF (CMD .EQ. 'TCLOUT') THEN | 194 |
| IF (CMD .EQ. 'PARTDIA') THEN | 222 |
| IF (CMD .EQ. 'CEMENT') THEN | 250 |
| IF (CMD .EQ. 'FGE') THEN | 278 |
| IF (CMD .EQ. 'FSE') THEN | 306 |
| IF (CMD .EQ. 'FCE') THEN | 334 |
| IF (CMD .EQ. 'POROSITY') THEN | 362 |
| IF (CMD .EQ. 'PRESSURE') THEN | 390 |
| IF (CMD .EQ. 'SATBRIN') THEN | 418 |
| | Relative line number ^^^^ |

| | |
|--|---------------------------|
| SUBROUTINE SCATXT1(IWASFLG, IASPRP, XMATPR, NAMELB, NMATPR, | 11604 |
| | Absolute line number ^^^^ |
| IF (CMD .EQ. 'EOF') GO TO 90 | 115 |
| IF (CMD(1:5) .EQ. 'NHITS') THEN | 117 |
| IF (CMD .EQ. 'EOF') GO TO 999 | 162 |
| IF (CMD(1:5) .EQ. 'TINTR') THEN | 165 |
| IF (CMD .EQ. 'DIAMMOD') THEN | 194 |

| | |
|--|-----|
| IF (CMD .EQ. 'TCLOUD') THEN | 223 |
| IF (CMD .EQ. 'PARTDIA') THEN | 252 |
| IF (CMD .EQ. 'DNSFLUID') THEN | 281 |
| IF (CMD .EQ. 'DOMEGA') THEN | 307 |
| IF (CMD .EQ. 'ABSROUGH') THEN | 334 |
| IF (CMD .EQ. 'TAUFAIL') THEN | 361 |
| IF (CMD .EQ. 'VISCO') THEN | 388 |
| IF (CMD .EQ. 'YLDSTRSS') THEN | 415 |
| IF (CMD(1:5) .EQ. 'INTR_'.AND. BRAGCDB) THEN | 442 |
| IF (CMD .EQ. 'NO_GRDVOL') THEN | 477 |
| IF (CMD .EQ. 'DIABRAG') THEN | 509 |
| IF (CMD .EQ. 'GRIDVOL') THEN | 530 |
| IF (CMD .EQ. 'PORO_NM') THEN | 551 |
| IF (CMD .EQ. 'PRGS_NM') THEN | 572 |
| IF (CMD .EQ. 'PRBR_NM') THEN | 593 |
| IF (CMD .EQ. 'SATG_NM') THEN | 614 |
| IF (CMD .EQ. 'POROSITY') THEN | 635 |
| IF (CMD .EQ. 'PRESS') THEN | 668 |
| IF (CMD .EQ. 'SATBRIN') THEN | 697 |

Relative line number ^^^^

END OF APPENDIX L