

APPENDIX E: User Defined Solution Control

You can control the I/O, monitor the energies and other solution norms of interest, and shut down the problem whenever you please with user subroutine `uctrl1` in `dyn21.F`. The arguments are defined in the listing provided below. This subroutine is called each time step and does not need any control card to operate.

```

      subroutine uctrl1 (numnp,ndof,time,dt1,dt2,prtc,pltc,frci,prto,
      . plto,frco,vt,vr,at,ar,ut,ur,xmst,xmsr,irbody,rbdyn,usrhv,
      . messag,totalm,cycle,idrint,mtype,lrb,nrba,rbcor,x,rbv,nrbn,
      . nrb,xrb,yrb,zrb,axrb,ayrb,azrb,dtx,nmmat,rba,fvalnew,fvalold,
      . fvalmid,fvalnxt)
c
c*****
c|  Livermore Software Technology Corporation   (LSTC)           |
c|  -----
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c*****
c
c      user subroutine for solution control.  It is called at the
c      beginning of time step n+1.
c
c
c      note:  ls-dyna uses an internal numbering system to
c             accommodate arbitrary node numbering.  to access
c             information for user node n, address array location m,
c             m = lqf8(n,1).  to obtain user node number, n,
c             corresponding to array address m, set n = lqfinv(m,1)
c
c      arguments:
c      numnp = number of nodal points
c      ndof  = number of degrees of freedom per node
c      time  = current solution time at n+1
c      dt1   = time step size between time n-1 and n
c      dt2   = time step size between time n and n+1
c      prtc  = output interval for taurus time history data
c      pltc  = output interval for taurus state data
c      frci  = output interval for taurus interface force data
c      prto  = output time for time history file
c      plto  = output time for state data
c      frco  = output time for force data
c      vt(3,numnp) = nodal translational velocity vector
c      vr(3,numnp) = nodal rotational velocity vector.  this
c                  array is defined if and only if ndof = 6
c      at(3,numnp) = nodal translational acceleration vector
c      ar(3,numnp) = nodal rotational acceleration vector.  this
c                  array is defined if and only if ndof = 6
c      ut(3,numnp) = nodal translational displacement vector
c      ur(3,numnp) = nodal rotational displacement vector.  this
c                  array is defined if and only if ndof = 6
c      xmst(numnp) = reciprocal of nodal translational masses
c      xmsr(numnp) = reciprocal of nodal rotational masses.  this
c                  array is defined if and only if ndof = 6

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c      fvalold      = array for storing load curve values at time n
c      fvalnew      = array for setting load curve values at time n+1
c                   only load curves with 0 input points may be user
c                   defined. When the load curve is user set, the
c                   value at time n must be stored in array fvalold.
c      fvalmid      = array for predicting load curve values at time n+3/2
c      fvalnxt      = array for predicting load curve values at time n+2
c                   for some applications it is necessary to predict the
c                   load curve values at time n+2, a time that is not known,
c                   this for instance for boundary prescribed motion. In
c                   this case the load curve values at time n+3/2 need to
c                   be predicted in fvalmid, and fvalold should be set to
c                   fvalnew and fvalnew should be set to fvalnxt. See
c                   coding below.
c
c      irbody       = 0 if no rigid bodies
c      rbdyn(numnp) = flag for rigid body nodal points
c                   if deformable node then set to 1.0
c                   if rigid body node then set to 0.0
c                   defined if and only if rigid bodies are present
c                   i.e., irbody.ne.0 if no rigid bodies are
c                   present
c      usrhv(lenhv) = user defined history variables that are stored
c                   in the restart file. lenhv = 100+7*nummat where
c                   nummat is the # of materials in the problem.
c                   array usrhv is updated only in this subroutine.
c      messag       = flag for dyna3d which may be set to:
c                   'sw1.' ls-dyna3d terminates with restart file
c                   'sw3.' ls-dyna3d writes a restart file
c                   'sw4.' ls-dyna3d writes a plot state
c      totalm       = total mass in problem
c      cycle        = cycle number
c      idrint       = flag for dynamic relaxation phase
c                   .ne.0: dynamic relaxation in progress
c                   .eq.0: solution phase
c      mtype(*)     = material type for each part in the model
c      lrb(*)       = lead rigid body for each rigid body
c                   If part n is rigid and is a lead or has not
c                   been merged then lrb(n)==n.
c      nrba(*)      = starting index in nrb(*) of the nodes for each
c                   rigid body
c      rbcor(3,*)   = rigid body cg coordinates
c      x(3,*)       = Node coordinate array
c      rbv(6,*)     = rigid body cg velocity
c      nrbn(*)      = # nodes in each rigid body
c      nrb(*)       = List of all rigid body nodes
c      xrb(*)       = RB scratch array as long as longest nrbn() value
c      yrb(*)       = RB scratch array as long as longest nrbn() value
c      zrb(*)       = RB scratch array as long as longest nrbn() value
c      axrb(*)      = RB scratch array as long as longest nrbn() value
c      ayrb(*)      = RB scratch array as long as longest nrbn() value
c      azrb(*)      = RB scratch array as long as longest nrbn() value
c      dtx          = (dt1+dt2)*0.5 except at time 0 when it is = dt2
c      nmmat        = number of parts in the model
c      rba(6,*)     = rigid body cg acceleration
c      fvalnew      = array for setting load curve values at time n+1
c                   only load curves with 0 input points may be user
c                   defined. When the load curve is user set, the
c                   value at time n must be stored in array fvalold.
c      fvalold      = array for storing load curve values at time n
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c      fvalmid      = array for predicting load curve values at time n+3/2
c      fvalnxt      = array for predicting load curve values at time n+2
c                   for some applications it is necessary to predict the
c                   load curve values at time n+2, a time that is not known,
c                   this for instance for boundary prescribed motion. In
c                   this case the load curve values at time n+3/2 need to
c                   be predicted in fvalmid, and fvalold should be set to
c                   fvalnew and fvalnew should be set to fvalnxt. See
c                   coding below.
c
c      include 'ptimes.inc'
c
c      prtims(1-37)=output intervals for ascii files
c
c      ascii files:
c      ( 1)-cross section forces
c      ( 2)-rigid wall forces
c      ( 3)-nodal data
c      ( 4)-element data
c      ( 5)-global data
c      ( 6)-discrete elements
c      ( 7)-material energies
c      ( 8)-noda interface forces
c      ( 9)-resultant interface forces
c      (10)-smug animator
c      (11)-spc reaction forces
c      (12)-nodal constraint resultant forces
c      (13)-airbag statistics
c      (14)-avs database
c      (15)-nodal force groups
c      (16)-output intervals for nodal boundary conditions
c      (17)-(32) unused at this time
c      (37)-auto tiebreak damage output
c
c      prt1st(32)=output times for ascii files above. when solution time
c                   exceeds the output time a print state is dumped.
c
c      common/rbkeng/enrbdy,rbdyx,rbdy,rbdyz
c
c      total rigid body energies and momentums:
c      enrbdy = rigid body kinetic energy
c      rbdyx = rigid body x-momentum
c      rbdyy = rigid body y-momentum
c      rbdyz = rigid body z-momentum
c
c      common/swmke/swxmom,swymom,swzmom,swkeng
c
c      total stonewall energies and momentums:
c      swxmom = stonewall x-momentum
c      swymom = stonewall y-momentum
c      swzmom = stonewall z-momentum
c      swkeng = stonewall kinetic energy
c
c      common/deengs/deeng
c
c      deeng = total discrete element energy
c
c      common/bk28/summss,xke,xpe,tt,xte0,erodeke,erodeie,selie,selke,
c      . erodehg
c

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c      xpe  = total internal energy in the finite elements
c
c      common/sprengs/spreng
c
c      spreng = total spr energy
c
c      character*(*) messag
c      integer cycle
c      real*8 x
c      dimension vt(3,*),vr(3,*),at(3,*),ar(3,*),
c      . xmst(*),xmsr(*),rbdyn(*),usrhv(*),mtype(*),lrb(*),nrba(*),
c      . rbcor(3,1),x(*),rbv(6,*),nrbn(*),nrb(*),xrb(*),yrb(*),
c      . zrb(*),axrb(*),ayrb(*),azrb(*),rba(6,*),fvalnew(*),
c      . fvalold(*),fvalmid(*),fvalnxt(*)
c      real*8 ut(3,*),ur(3,*)
c
c      sample momentum and kinetic energy calculations
c
c      remove all comments in column 1 below to activate
c      i = 1
c      if (i.eq.1) return
c      return
cc
cc
cc      initialize kinetic energy, xke, and x,y,z momentums.
cc
c      xke = 2.*swkeng+2.*enrbdy
c      xm = swxmom+rbdyx
c      ym = swymom+rbdyy
c      zm = swzmom+rbdyz
cc
c      numnp2 = numnp
c      if (ndof.eq.6) then
c          numnp2 = numnp+numnp
c      endif
c      write(iotty,*)ndof
cc
cc
cc      no rigid bodies present
cc
c      if (irbody.eq.0) then
cc          note in blank comment vr follows vt.  this fact is used below.
c          do 10 n = 1,numnp2
c              xmsn = 1./xmst(n)
c              vn1 = vt(1,n)
c              vn2 = vt(2,n)
c              vn3 = vt(3,n)
c              xm = xm+xmsn*vn1
c              ym = ym+xmsn*vn2
c              zm = zm+xmsn*vn3
c              xke = xke+xmsn*(vn1*vn1+vn2*vn2+vn3*vn3)
c      10  continue
cc
cc
cc      rigid bodies present
cc
c      else
cc          nodal accerations for rigid bodies
cc
c          do 12 n = 1,nmmat
```

```

c      if (mtype(n).ne.20.or.lrb(n).ne.n) go to 12
c      lrbn = nrba(n)
c      call stvlut(rbcor(1,n),x,vt,at,ar,vr,rbv(1,n),dt2,
c      . nrbn(n),nrb(lrbn),xrb,yrb,zrb,axrb,ayrb,azrb,dtx)
c
c      rigid body nodal accelerations
c
c      if (ndof.eq.6) then
c          call rbnacc(nrbn(n),nrb(lrbn),rba(4,n),ar)
c      endif
c
c 12    continue
cc
c      do 20 n = 1,numnp
c      xmsn = 1./xmst(n)
c      vn1 = rbdyn(n)*vt(1,n)
c      vn2 = rbdyn(n)*vt(2,n)
c      vn3 = rbdyn(n)*vt(3,n)
c      xm = xm+xmsn*vn1
c      ym = ym+xmsn*vn2
c      zm = zm+xmsn*vn3
c      xke = xke+xmsn*(vn1*vn1+vn2*vn2+vn3*vn3)
c 20    continue
c      if (ndof.eq.6) then
c          do 30 n = 1,numnp
c          xmsn = 1./xmsr(n)
c          vn1 = rbdyn(n)*vr(1,n)
c          vn2 = rbdyn(n)*vr(2,n)
c          vn3 = rbdyn(n)*vr(3,n)
c          xm = xm+xmsn*vn1
c          ym = ym+xmsn*vn2
c          zm = zm+xmsn*vn3
c          xke = xke+xmsn*(vn1*vn1+vn2*vn2+vn3*vn3)
c 30    continue
c      endif
c      endif
cc
cc      total kinetic energy
c      xke=.5*xke
cc      total internal energy
c      xie = xpe+deeng+spreng
cc      total energy
c      xte = xke+xpe+deeng+spreng
cc      total x-rigid body velocity
c      xrbv = xm/totalm
cc      total y-rigid body velocity
c      yrbv = ym/totalm
cc      total z-rigid body velocity
c      zrbv = zm/totalm
c      return
c      end

```

