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=====
c   History: sode.f
c
c   Driver routine which integrates ODEs defining
c   model for deuteron.
c
c   See class notes and Arfken, Math. Methods for
c   Physicists, 2nd Edition, section 9.1.2
c   for more details.
=====

      program          deut

      implicit        none

      character*4     cdnm
      parameter      ( cdnm = 'deut' )

      integer         iargc,          indlnb,          i4arg
      real*8          r8arg

      real*8          r8_never
      parameter      ( r8_never = -1.0d-60 )

c-----
c   Order of system.
c-----

      integer         neq
      parameter      ( neq = 2 )

c-----
c   Storage for solution at requested output radii.
c-----

      integer         maxout
      parameter      ( maxout = 10 000 )

      real*8          y0(neq)
      real*8          vxout(maxout), vyout(maxout,neq),
&                   work(maxout)
      integer         nxout,          ixout,          nxout_succ

```

```

integer      ieq

logical      ltrace
parameter   ( ltrace = .true. )

integer      maxdump
parameter   ( maxdump = 50 )

logical      lsodatrace
parameter   ( lsodatrace = .false. )

logical      ivs_ok

```

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c-----
c   LSODA Variables.
c-----

```

```

external     fcn,      jac

real*8       y(neq)
real*8       tbgn,     tend
integer      itol
real*8       rtol,     atol
integer      itask,    istate,  iopt
integer      lrw

parameter    ( lrw = 22 + neq * 16 )
real*8       rwork(lrw)

integer      liw
parameter    ( liw = 20 + neq )
integer      iwork(liw)
integer      jt

real*8       tol
real*8       default_tol
parameter    ( default_tol = 1.0d-8 )

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c-----
c   Common communication with routine 'fcn' in 'fcn.f' ...
c-----

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include          'fcn.inc'

c-----
c   Parse command line arguments.  Deviation from
c   'sode'.  Usage is
c
c   deut <x0> <E> [<tol>]
c-----

      if( iargc() .lt. 3 ) go to 900

      x0      = r8arg(1,r8_never)
      E      = r8arg(2,r8_never)
      tol    = r8arg(3,default_tol)
      if( x0 .eq. r8_never .or. E .eq. r8_never )
&      go to 900

c-----
c   Both boundary conditions are fixed in this case.
c-----

      y0(1) = 0.0d0
      y0(2) = 1.0d0

c-----
c   Echo command line arguments if "local tracing"
c   enabled ...
c-----

      if( ltrace ) then
          write(0,*) 'x0: ', x0
          write(0,*) 'E: ', E
          write(0,*) 'tol: ', tol
      end if

c-----
c   Get output radii from standard input ...
c-----

      call dvfrom('-',vxout,nxout,maxout)
      if( nxout .le. 0 ) then
          write(0,*)
          write(0,*) cdnm//': No output radii read from standard input'
          write(0,*) cdnm//': Use (e.g.) dvmesh or dvgmesh to '//

```

```

&          'generate output radii.'
  write(0,*)
  write(0,*) cdnm//': Sample usages:'
  write(0,*)
  write(0,*) cdnm//': dvmesh  0.0 1.0 101'
  write(0,*) cdnm//': dvgmesh 0.0 1.0 101 10'
  write(0,*)
  stop
end if
if( ltrace ) then
  if( nxout .le. maxdump ) then
    call dvdump(vxout,nxout,cdnm//': output radii',0)
  else
    call dvdmp1(vxout,work,
&          int(1.0d0 * nxout / maxdump + 0.5d0),
&          nxout,cdnm//': selected output radii',0)
  end if
  write(0,*)
  write(0,*) cdnm//': Initial time: ', vxout(1)
  write(0,*)
end if

```

```

c-----
c   Set LSODA parameters ...
c-----

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```

  itol  = 1
  rtol  = tol
  atol  = tol
  itask = 1
  iopt  = 0
  jt    = 2

```

```

c-----
c   Initialize the solution ...
c-----

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```

  do ieq = 1 , neq
    y(ieq)      = y0(ieq)
    vyout(1,ieq) = y0(ieq)
  end do

```

```

c-----
c   Do the integration ...
c-----
      do ixout = 2 , nxout
          istate = 1
c-----
c   Need these temporaries since lsoda overwrites
c   tend ...
c-----

      tbgn = vxout(ixout-1)
      tend = vxout(ixout)
      call lsoda(fcn,neq,y,tbgn,tend,
&               itol,rtol,atol,itask,
&               istate,iopt,rwork,lrw,iwork,liw,jac,jt)
      if( lsodatrace ) then
          write(0,1000) cdnm, ixout, nxout, vxout(ixout),
&               vxout(ixout+1)
1000      format(' ',a,': Step ',i4,' t = ',1pe10.3,
&               ' .. ',1pe10.3)
          write(0,*) cdnm//': lsoda reurns ', istate
      end if

      if( istate .lt. 0 ) then
          write(0,1500) cdnm, istate, ixout, nxout,
&               vxout(ixout-1), vxout(ixout)
1500      format(/' ',a,': Error return ',i2,
&               ' from integrator LSODA.'/
&               '           At output time ',i5,' of ',i5/
&               '           Interval ',1pe11.3,' .. ',
&               1pe11.3/)
          nxout_succ = ixout - 1
          go to 500
      end if
      do ieq = 1 , neq
          vyout(ixout,ieq) = y(ieq)
      end do
  end do
  nxout_succ = nxout

```

```

500  continue

      do ixout = 1 , nxout_succ
        write(*,2000)  vxout(ixout),
&                    ( vyout(ixout,ieq) , ieq = 1 , neq )
c-----
c      NOTE: This format will not dump all values on a
c      single line if neq > 10 !!
c-----
2000  format(1P,10E25.16)
      end do

      stop

900  continue
      write(0,*) 'usage: '//cdnm//
&          ' <x0> <E> [<tol>]'
      write(0,*) ' '
      write(0,*) '      Output radii read from standard input'
      stop

      end

```

```
c-----  
c      Driver routine which integrates ODEs defining  
c      model for deuteron.  
c  
c      See class notes and Arfken, Math. Methods for  
c      Physicists, 2nd Edition, section 9.1.2  
c      for more details.  
c-----  
c      neq = 2  
c-----
```

```
subroutine fcn(neq,x,y,yprime)  
  implicit none  
  
  include 'fcn.inc'  
  
  integer neq  
  real*8 x, y(neq), yprime(neq)  
  
  real*8 u, w  
  
  u = y(1)  
  w = y(2)  
  yprime(1) = w  
  
  if( x .le. x0 ) then  
    yprime(2) = (-1.0d0 - E) * y(1)  
  else  
    yprime(2) = -E * y(1)  
  end if  
  
  return  
end
```

```
c-----  
c   Application specific common block for communication  
c   with derivative evaluating routine 'fcn'.  
c  
c   x0:   Range of square potential well  
c   E:   Energy (sought eigenvalue)  
c-----
```

```
      real*8  
&      x0,  
&      E  
      common / com_fcn /  
&      x0,  
&      E
```



```
c-----  
c      Integrates static equations of motion for spherically  
c      symmetric, general-relativistic boson star.  
c  
c      See class notes and Colpi et al, Phys Rev Lett,  
c      vol 57, 2485--2488 for more details  
c-----  
c      neq = 4  
c-----
```

```
subroutine fcn(neq,x,y,yprime)  
  implicit none  
  
  include 'fcn.inc'  
  
  integer neq  
  real*8 x, y(neq), yprime(neq)  
  
  real*8 A, M, B, sigma, xi  
  
  M = y(1)  
  B = y(2)  
  sigma = y(3)  
  xi = y(4)
```

```

if( x .eq. 0.0d0 ) then
  A = 1.0d0
  yprime(1) = 0.0d0
  yprime(2) = 0.0d0
  yprime(3) = 0.0d0
  yprime(4) = ( (Omegasq / B - 1.0d0) * sigma -
&                Lambda * sigma**3 ) / 3.0d0
else
  A = 1.0d0 / (1.0d0 - 2.0d0 * M / x)
  yprime(1) = x**2 * (
&    0.5d0 * (Omegasq / B + 1.0d0) * sigma**2 +
&    0.25d0 * Lambda * sigma**4 +
&    0.5d0 * xi**2 / A
&    )
  yprime(2) = A * B * x * (
&    (1.0d0 - 1.0d0 / A) / (x**2) +
&    (Omegasq / B - 1.0d0) * sigma**2 -
&    0.5d0 * Lambda * sigma**4 +
&    xi**2 / A
&    )
  yprime(3) = xi
  yprime(4) = xi * (
&    -2.0d0 / x - 0.5d0 * yprime(2) / B +
&    A * (yprime(1) - M / x) / x
&    ) - A * (
&    (Omegasq / B - 1.0d0) * sigma -
&    Lambda * sigma**3 )
end if

return
end

```

```
c-----  
c   Application specific common block for communication  
c   with derivative evaluating routine 'fcn'.  
c  
c   Lambda:   Dimensionless  $\phi^4$  coupling constant  
c   Omega:    Eigen-frequency  
c   Omegasq:  Square of eigen-frequency  
c-----
```

```
      real*8  
      &          Lambda,  
      &          Omega,  
      &          Omegasq  
      common / com_fcn /  
      &          Lambda,  
      &          Omega,  
      &          Omegasq
```