

```

=====
c      Computes and reports estimate of machine epsilon.
c
c      Recall: machine epsilon is smallest positive 'eps'
c      such that
c
c              (1.0d0 + eps ) .ne. (1.0d0)
c
c      Program accepts optional argument which specifies
c      division factor: values close to 1.0 will result
c      in more accurate estimate of machine epsilon.
=====
      program          meps

      implicit        none

c-----
c      Note use of 'r8arg', available in 'libp410f.a' which
c      works exactly like 'i4arg' except that it returns
c      a real*8 value parsed from the specified command-line
c      argument
c-----
      real*8          r8arg

      real*8          default_fac
      parameter      ( default_fac = 2.0d0 )

      real*8          eps,          neweps,          fac

      fac = r8arg(1,default_fac)
      write(0,*) 'meps: using division factor: ', fac

```

```
eps      = 1.0d0
neweps   = 1.0d0
do while( .true. )
  if( 1.0d0 .eq. (1.0d0 + neweps) ) then
    write(*,*) eps
    stop
  else
    eps      = neweps
    neweps   = neweps / fac
  end if
end do

stop

end
```

```
#####  
# Output from 'meps' on SGIs (IEEE 64-bit floating point).  
#####
```

Script started on Wed Sep 20 20:15:05 2000

```
sgi1 1> make meps  
f77 -g -64 -c meps.f  
f77 -g -64 -L/usr/local/lib meps.o -lp410f -o meps
```

```
sgi1 2> meps  
meps: using division factor: 2.0000000000000000  
2.2204460492503131E-16
```

```
sgi1 3> meps 1.01  
meps: using division factor: 1.0100000000000000  
1.1104218387155329E-16
```

```
sgi1 4> meps 1.0001  
meps: using division factor: 1.0001000000000000  
1.1102645224601785E-16
```

```
#####  
# Output from 'meps' on PC Linux machine (80 bit floating pt)  
#####
```

Script started on Mon Oct 1 16:51:34 2001

lnx1 1> make meps

pgf77 -g -Msecond\_underscore -c meps.f

pgf77 -g -Msecond\_underscore -L/usr/local/PGI/lib meps.o -lp410f -o mep

Linking:

lnx1 2> meps

meps: using division factor: 2.0000000000000000  
1.0842021724855044E-019

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lnx1 3> meps 1.01

meps: using division factor: 1.0100000000000000  
5.4364534909517435E-020

FORTRAN STOP

lnx1 4> meps 1.0001

meps: using division factor: 1.0001000000000000  
5.4212146310714582E-020

FORTRAN STOP

```

=====
c   Program illustrating "catastrophic" loss of precision
c   resulting from the subtraction of two nearly equal
c   floating point values.
=====
      program          catprec

      implicit        none

      real*8          x
      parameter      ( x = 0.2d0 )

      integer         i
      real*8          h,      dsinx

      write(*,*) '      h      d(sin) approx  '//
&      'd(sin) exact  d(sin) err'
      write(*,*)

      h = 0.5d0
      do i = 1 , 16

c-----
c      Algebraically, in the limit  $h \rightarrow 0$ , dsinx should
c      approach  $\cos(x)$ , but  $\sin(x+h) \rightarrow \sin(x)$  so
c      catastrophic loss of precision occurs.
c-----

      dsinx = (sin(x+h) - sin(x)) / h
      write(*,1000) h, dsinx, cos(x), dsinx - cos(x)
1000  format(1P,E12.3,2E16.8,E12.3)
      h = 0.125d0 * h
      end do

      stop
      end

```

```
#####
# Output from 'catprec' illustrating catastrophic precision
# loss due to subtraction of nearly-equal floating point
# values.
#####
```

Script started on Mon Oct 1 16:53:38 2001

```
lnx1 1> make catprec
pgf77 -g -Msecond_underscore -c catprec.f
pgf77 -g -Msecond_underscore -L/usr/local/PGI/lib catprec.o -o catprec
Linking:
```

```
lnx1 2> catprec
      h          d(sin) approx    d(sin) exact    d(sin) err
5.000E-01  8.91096713E-01  9.80066578E-01  -8.897E-02
6.250E-02  9.73222242E-01  9.80066578E-01  -6.844E-03
7.813E-03  9.79280560E-01  9.80066578E-01  -7.860E-04
9.766E-04  9.79969416E-01  9.80066578E-01  -9.716E-05
1.221E-04  9.80054450E-01  9.80066578E-01  -1.213E-05
1.526E-05  9.80065062E-01  9.80066578E-01  -1.516E-06
1.907E-06  9.80066388E-01  9.80066578E-01  -1.895E-07
2.384E-07  9.80066554E-01  9.80066578E-01  -2.368E-08
2.980E-08  9.80066575E-01  9.80066578E-01  -2.960E-09
3.725E-09  9.80066577E-01  9.80066578E-01  -3.702E-10
4.657E-10  9.80066578E-01  9.80066578E-01  -5.373E-11
5.821E-11  9.80066578E-01  9.80066578E-01  -1.701E-10
7.276E-12  9.80066577E-01  9.80066578E-01  -8.686E-10
9.095E-13  9.80066568E-01  9.80066578E-01  -1.018E-08
1.137E-13  9.80066538E-01  9.80066578E-01  -3.998E-08
1.421E-14  9.80066299E-01  9.80066578E-01  -2.784E-07
FORTRAN STOP
```

```

=====
c      Implements matrix-matrix multiply
c
c      c = a b
c
c      where a, b and c are n x n (square) real*8 matrices.
=====
      subroutine dmmult(a,b,c,n)

          implicit      none

          integer      n
          real*8      a(n,n),    b(n,n),    c(n,n)

          integer      i,    j,    k

          do j = 1 , n
              do i = 1 , n
                  c(i,j) = 0.0d0
                  do k = 1 , n
                      c(i,j) = c(i,j) + a(i,k) * b(k,j)
                  end do
              end do
          end do

          return

      end

```

```

c=====
c   Writes a double precision matrix (two dimensional
c   array) to file 'fname'.  If 'fname' is the
c   string '-', the matrix is written to standard input.
c
c   This routine is modelled on 'dvto' previously
c   discussed in class: see ~phys410/f77/ex3/dvto.f
c=====
      subroutine dmt0(fname,a,d1,d2)
c-----
c   Arguments:
c
c       fname:  (I)   File name
c       a:      (I)   Input matrix
c       d1:     (I)   First dimension of a
c       d2:     (I)   Second dimension of a
c-----
      implicit      none
      integer      indlnb,      getu

      character*(*) fname
      integer      d1,          d2
      real*8       a(d1,d2)

      integer      uestdout
      parameter    ( uestdout = 6 )

      integer      uto,          rc
c-----
c       Parse fname: either "attach" 'uto' to stdout or
c       get a unit number using 'getu', and open the
c       file 'fname' for formatted I/O via 'uto'
c-----

```



```

if( fname .eq. '-' ) then
    uto = stdout
else
    uto = getu()
    open(uto,file=fname(1:indlnb(fname)),
&        form='formatted',iostat=rc)
    if( rc .ne. 0 ) then
        write(0,*) 'dmto: Error opening ',
&                fname(1:indlnb(fname))
        return
    end if
end if
c-----
c      Write dimensions, then array elements
c-----

write(uto,*,iostat=rc) d1, d2
if( rc .ne. 0 ) then
    write(0,*) 'dmto: Error writing dimensions'
go to 500
end if

write(uto,*,iostat=rc) a
if( rc .ne. 0 ) then
    write(0,*) 'dmto: Error reading matrix'
end if
c-----
c      Exit: Close file and return
c-----
500    continue
        close(uto)

        return
        end

```

```

c=====
c   Returns a double precision matrix (two dimensional
c   array) read from file 'fname'.  If 'fname' is the
c   string '-', the matrix is read from standard input.
c
c   The dimensions of the matrix must precede the matrix
c   elements themselves in the file.  Specifically, the
c   file should have been created using the following
c   list-directed, formatted READ statement
c   (or equivalent):
c
c       integer      d1,      d2
c       real*8       a(d1,d2)
c       integer      uout
c       write(uout,*) d1, d2
c       write(uout,*) a
c
c   This routine is modelled on 'dvfrom' previously
c   discussed in class: see ~phys410/f77/ex3/dvfrom.f
c
c   Note the use of helper routine 'dmfrom1' which
c   reads actual array values once bounds have been
c   extracted from file.
c=====
c       subroutine dmfrom(fname,a,d1,d2,asize)
c-----
c   Arguments:
c
c       fname:  (I)   File name
c       a:      (O)   Return matrix
c       d1:     (O)   First dimension of a
c       d2:     (O)   Second dimension of a
c       asize:  (I)   Maximum size (d1 * d2) of a
c-----

```

```

implicit          none

integer           indlnb,      getu

character*(*)     fname
integer           d1,          d2,          asize
real*8            a(d1,d2)

integer           ustdin
parameter        ( ustdin = 5 )

integer           ufrom,      rc

```

```

c-----
c      Parse fname: either "attach" 'ufrom' to stdin or
c      get a unit number using 'getu', and open the
c      file 'fname' for formatted I/O via 'ufrom'
c-----
      if( fname .eq. '-' ) then
          ufrom = ustdin
      else
          ufrom = getu()
          open(ufrom,file=fname(1:indlnb(fname)),
&             form='formatted',iostat=rc,status='old')
          if( rc .ne. 0 ) then
              write(0,*) 'dmfrom: Error opening ',
&                         fname(1:indlnb(fname))
              return
          end if
      end if
end if

```

```
c-----  
c      Read dimensions and abort if there is insufficient  
c      storage for the entire matrix. Note the 'go to'  
c      to the 'exit block' since we've opened a file now  
c      and should close it, even if there's an error.  
c      Also, we set the dimensions to 0 for all error  
c      conditions as a way of communicating failure to  
c      the calling routine.  
c-----
```

```
      read(ufrom,*,iostat=rc) d1, d2  
      if( rc .ne. 0 ) then  
          write(0,*) 'dmfrom: Error reading dimensions'  
          d1 = 0  
          d2 = 0  
      go to 500  
      end if  
      if( (d1 * d2) .gt. asize ) then  
          write(0,*) 'dmfrom: Insufficient storage'  
          d1 = 0  
          d2 = 0  
      go to 500  
      end if
```

```
c-----  
c      Now that dimensions have been determined call  
c      helper routine to read values  
c-----
```

```
      call dmfrom1(ufrom,a,d1,d2,rc)  
      if( rc .ne. 0 ) then  
          write(0,*) 'dmfrom: Error reading matrix'  
          d1 = 0  
          d2 = 0  
      end if
```

```

c-----
c      Exit: Close file and return
c-----
500      continue
          close(ufrom)

          return
end

c=====
c      Helper routine for dmfrom: Reads array values, returns
c      I/O status to calling routine via 'rc'
c=====
      subroutine dmfrom1(ufrom,a,d1,d2,rc)

          implicit      none

          integer      d1,      d2,      ufrom,      rc
          real*8      a(d1,d2)

          read(ufrom,*,iostat=rc) a

          return

end

```

```

=====
c      Test program for subroutine 'dmfrom', 'dmto' and
c      'dmmult' (see 'dmroutines.f')
c
c      Program expects one argument, the name of a file which
c      contains a real*8 square matrix written as described
c      in the documentation for 'dmfrom' in 'dmroutines.f'
c      Use '-' to read from stdin. Program then computes
c      square of matrix and outputs result to stdout.
=====

```

```

      program          tdm

      implicit        none

      integer         iargc

      character*256   fname

```

```

-----
c      Maximum size for input and output arrays (matrices).
-----

```

```

      integer         maxsize
      parameter      ( maxsize = 100 000 )
      real*8         a(maxsize),  asq(maxsize)
      integer        d1a,         d2a

```

```

      if( iargc() .ne. 1 ) go to 900
      call getarg(1,fname)

```

```

-----
c      Read matrix ...
-----
      call dmfrom(fname,a,d1a,d2a,maxsize)

```

```

        if( d1a .gt. 0 .and. d2a .gt. 0 ) then
            if( d1a .eq. d2a ) then
c-----
c          Compute square ...
c-----
                call dmmult(a,a,asq,d1a,d1a)
c-----
c          ... and output.
c-----
                call dmt0('-',asq,d1a,d1a)
            else
                write(0,*) 'tdm: Input array not square'
            end if
        else
            write(0,*) 'tdm: dmfrom() failed'
        end if

        stop

900 continue
        write(0,*) 'usage: tdm <file name>'
        write(0,*)
        write(0,*) '          Use ''tdm -'' to read ',
&          'from standard input'

        stop

    end

```

```
#####  
# Building 'tdm' and sample output  
#####
```

```
Script started on Mon Oct 1 16:54:49 2001
```

```
lnx1 1> pwd; ls
```

```
/home/phys410/f77/ex6
```

```
Makefile          catprec.f          dmroutines.f      meps.f            tdm.f
```

```
lnx1 2> make tdm
```

```
pgf77 -g -Msecond_underscore -c tdm.f
```

```
pgf77 -g -Msecond_underscore -c dmroutines.f
```

```
pgf77 -g -Msecond_underscore -L/usr/local/PGI/lib tdm.o dmroutines.o -l
```

```
Linking:
```

```
lnx1 3> tdm
```

```
usage: tdm <file name>
```

```
Use 'tdm -' to read from standard input
```

```
FORTRAN STOP
```

```
lnx1 4> tdm -
```

```
2 2
```

```
1 2 3 4
```

```
2 2
```

```
7.0000000000000000
```

```
10.0000000000000000
```

```
15.0000000000000000
```

```
22.0000000000000000
```

```
FORTRAN STOP
```

```
lnx1 5> tdm -
```

```
2 3
```

```
1 2 3 4 5 6
```

```
tdm: Input array not square
```

```
FORTRAN STOP
```



.IGNORE:

F77\_COMPILE = \$(F77) \$(F77FLAGS) \$(F77CFLAGS)  
F77\_LOAD = \$(F77) \$(F77FLAGS) \$(F77LFLAGS)

.f.o:

\$(F77\_COMPILE) \$\*.f

EXECUTABLES = meps catprec tdm

all: \$(EXECUTABLES)

meps: meps.o

\$(F77\_LOAD) meps.o -lp410f -o meps

catprec: catprec.o

\$(F77\_LOAD) catprec.o -o catprec

tdm: tdm.o dmroutines.o

\$(F77\_LOAD) tdm.o dmroutines.o -lp410f -o tdm

clean:

rm \*.o

rm \$(EXECUTABLES)

rm core