

Source file: meps.f

```
c=====
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.
c=====
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
c----- real*8      default_fac
c----- parameter    ( default_fac = 2.0d0 )
c----- real*8      eps,           neweps,        fac
c
fac = r8arg(1,default_fac)
write(0,*) 'meps: using division factor: ', fac
c
eps     = 1.0d0
neweps = 1.0d0
do while( 1.0d0 .ne. (1.0d0 + neweps) )
    eps     = neweps
    neweps = neweps / fac
end do
c
write(*,*) eps
c
stop
c
```

Source file: meps-pclinux-output

```
#####
# Output from 'meps' on PC Linux machine (80 bit floating pt)
#####
lnx1 1> make meps
pgf77 -g -c meps.f
pgf77 -g -L/usr/local/PGI/lib meps.o -lp410f -o meps
lnx1 2> meps
meps: using division factor: 2.000000000000000
1.0842021724855044E-019
lnx1 3> meps 1.01
meps: using division factor: 1.010000000000000
5.4364534909517435E-020
lnx1 4> meps 1.0001
meps: using division factor: 1.000100000000000
5.4212146310714582E-020
```

Source file: meps-sgi-output

```
#####
# Output from 'meps' on SGIs (IEEE 64-bit floating point).
#####
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: 2.000000000000000
2.2204460492503131E-16
sgi1 3> meps 1.01
meps: using division factor: 1.010000000000000
1.1104218387155329E-16
sgi1 4> meps 1.0001
meps: using division factor: 1.000100000000000
1.1102645224601785E-16
```

Source file: catprec.f

```
c=====
c      Program illustrating "catastrophic" loss of precision
c      resulting from the subtraction of two nearly equal
c      floating point values.
c=====
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 -> 0, dsinx should
c      approach cos(x), but sin(x+h) -> 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
```

Source file: catprec-output

```
#####
# Output from 'catprec' illustrating catastrophic precision
# loss due to subtraction of nearly-equal floating point
# values.
#####
lnx1 1> make catprec
pgf77 -g -c catprec.f
pgf77 -g -L/usr/local/PGI/lib catprec.o -o catprec
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
```

Source file: dmroroutines.f

```
c=====
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.
c=====
subroutine dmmmult(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 dmto(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          ustdout
parameter        ( ustdout = 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 = ustdout
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-----
c----- write(uto,*,iostat=rc) d1, d2
c----- if( rc .ne. 0 ) then
c-----   write(0,*) 'dmto: Error writing dimensions'
c----- go to 500
c----- end if

c----- write(uto,*,iostat=rc) a
c----- if( rc .ne. 0 ) then
c-----   write(0,*) 'dmto: Error reading matrix'
c----- end if

c----- Exit: Close file and return
c-----
500  continue
c----- if( uto .ne. ustdout ) then
c-----   close(uto)
c----- end if

c----- return
c----- 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----- 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----- integer      d1,      d2
c----- real*8      a(d1,d2)
c----- integer      uout
c----- write(uout,*) d1, d2
c----- write(uout,*) a
c----- This routine is modelled on 'dvfrom' previously
c----- discussed in class: see ~phys410/f77/ex3/dvfrom.f
c----- Note the use of helper routine 'dmfrom1' which
c----- reads actual array values once bounds have been
c----- extracted from file.
c=====

      subroutine dmfrom(fname,a,d1,d2,asize)
c-----
c----- Arguments:
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-----
c----- implicit none
c----- integer      indlnb,      getu
c----- character(*)  fname
c----- integer      d1,          d2,          asize
c----- real*8      a(d1,d2)
c----- integer      ustdin
c----- parameter    ( ustdin = 5 )
c----- 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-----
c----- if( fname .eq. '--' ) then
c-----   ufrom = ustdin
c----- else
c-----   ufrom = getu()
c-----   open(ufrom,file=fname(1:indlnb(fname)),form='formatted',iostat=rc,status='old')
c-----   if( rc .ne. 0 ) then
c-----     write(0,*) 'dmfrom: Error opening ', fname(1:indlnb(fname))
c-----   end if
c-----   return
c----- end if
c----- end if

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-----
c----- read(ufrom,*,iostat=rc) d1, d2
c----- if( rc .ne. 0 ) then
c-----   write(0,*) 'dmfrom: Error reading dimensions'
c-----   d1 = 0
c-----   d2 = 0
c----- go to 500
c----- end if
c----- if( (d1 * d2) .gt. asize ) then
c-----   write(0,*) 'dmfrom: Insufficient storage'
c-----   d1 = 0
c-----   d2 = 0
c----- go to 500
c----- end if
c----- Now that dimensions have been determined call
c----- helper routine to read values
c-----
c----- call dmfrom1(ufrom,a,d1,d2,rc)
c----- if( rc .ne. 0 ) then
c-----   write(0,*) 'dmfrom: Error reading matrix'
c-----   d1 = 0
c-----   d2 = 0
c----- end if
c----- Exit: Close file and return
c-----
500  continue
c----- if( ufrom .ne. ustdin ) then
c-----   close(ufrom)
c----- end if

c----- return
c----- 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)
c-----
c----- implicit none
c----- integer      d1,          d2,          ufrom,      rc
c----- real*8      a(d1,d2)
c----- read(ufrom,*,iostat=rc) a
c----- return
c----- end

```

Source file: tdm.f

```
c=====
c      Test program for subroutine 'dmfrom', 'dmto' and
c      'dmmmult' (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.
c=====
```

```
program      tdm
implicit      none
integer       iargc
character*256 fname

c-----
c      Maximum size for input and output arrays (matrices).
c-----
integer      maxsize
parameter     ( maxsize = 100 000 )
real*8       a(maxsize),   asq(maxsize)
integer       dia,          d2a

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

c-----
c      Read matrix ...
c-----
call dmfrom(fname,a,dia,d2a,maxsize)
if( dia .gt. 0 .and. d2a .gt. 0 ) then
  if( dia .eq. d2a ) then
c-----
c      Compute square ...
c-----
call dmmult(a,a,asq,dia,dia)
c-----
c      ... and output.
c-----
call dmto('-',asq,dia,dia)
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
```

Source file: tdm-output

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

lnx1 1> make tdm
pgf77 -g -c tdm.f
pgf77 -g -c dmroutines.f
pgf77 -g -L/usr/local/PGI/lib tdm.o dmroutines.o -lp410f -o tdm

lnx1 2> tdm
usage: tdm <file name>

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

```
lnx1 3> tdm -
2 2
1 2 3 4
               2               2
               7.000000000000000   10.000000000000000
               22.000000000000000  15.000000000000000

lnx1 4> tdm -
2 3
1 2 3 4 5 6
tdm: Input array not square
```

Source file: Makefile

```
.IGNORE:

F77_COMPILE = $(F77) $(F77FLAGS) $(F77CFLAGS)
F77_LOAD    = $(F77) $(F77FLAGS) $(F77LFLAGS)

.f.o:
$(F77_COMPILE) $*.f

EXECUTABLES = meps catprec tdm

all: $(EXECUTABLES)

meeps: meeps.o
$(F77_LOAD) meeps.o -lp410f -o meeps

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
```