Phy 329: Computational Physics Homework 3 Due: Monday, September 30, 3:00pm

The following assignment involves writing and testing several simple Fortran 77 programs. Follow the instructions carefully and do all compilation and execution on `einstein`*. All files required by the assignment must reside in the correct places on your SGI account for the homework to be considered complete.*

Problem 1: In directory `~/hw3/a1` on your SGI account create Fortran source files `pgm1.f` and `pgm2.f` which are to be compiled into executables `pgm1` and `pgm2` with behaviour as described below. Use `real*8` arithmetic and "free-format" I/O (i.e. `read(*,*)` and `write(*,*)`) here and throughout the homework.

- `pgm1`: reads x, y pairs (2 `real*8` numbers per line) from standard input, computes and writes $f(x, y)$ to standard output, where $f(x, y)$ is defined as follows:

$$f(x,y) = \sin(3xy)\left(\frac{1.0 + e^{-\frac{x^2+y^2}{2}}}{1.0 - e^{-\frac{x^2+y^2}{2}}}\right)$$

- `pgm2`: reads x, y, z triples (3 `real*8` numbers per line) from standard input, computes and writes $g(x, y, z)$ to standard output, where $g(x, y, z)$ is defined as follows:

$$g(x,y,z) = \begin{cases} \sqrt{x^2+y^2+z^2} & \text{if } x < 0 \\ \frac{\pi}{4} & \text{if } x = 0 \\ \cos(x)\sin(y)\tan(z) & \text{if } x > 0 \end{cases}$$

Minimal commenting of these programs is acceptable, but be sure to carefully test them. You may find it helpful to use `Maple` for testing purposes. (Recall that the `Maple` procedure `evalf()` evaluates an expression using floating-point arithmetic.) Your programs should detect and ignore invalid input as did the `mysum` example discussed in class. Should you wish to print messages if invalid input is encountered, be sure to direct the messages to standard error. I will evaluate your programs using my own input.

Problem 2: In directory `~/hw3/a2` create source files `dvvfrom.f` and `dvvto.f` which define Fortran `subroutines` with the following headers:

```
subroutine dvvfrom(fname,v1,v2,n,maxn)
   implicit        none
   character*(*)   fname
   integer         n,           maxn
   real*8          v1(maxn),    v2(maxn)


subroutine dvvto(fname,v1,v2,n)
   implicit        none
   character*(*)   fname
   integer         n
   real*8          v1(n),       v2(n)
```

These routines are to behave completely analogously to the routines `dvfrom` and `dvto` covered in the class notes except that they are to read and write *pairs* of `real*8` vectors (one-dimensional arrays) from and to a file with name `fname`. Thus the format of the input and output is:

```
v1(1)    v2(1)
v1(2)    v2(2)
v1(3)    v2(3)
   .
   .
   .
```

If `fname .eq. '-'`, the routines should use standard input or standard output as appropriate. In the same directory, write test programs `tdvvfrom.f` and `tdvvto.f` and create executables `tdvvfrom` and `tdvvto` which test your two routines. The arguments to `tdvvfrom` and `tdvvto` should be identical to the programs `tdvfrom` and `tdvto`, respectively, which were covered in class. Specifically, the usage of the test programs should be:

```
% tdvvfrom <file name>
% tdvvto   <file name> <n>
```

`tdvvfrom` should report on standard error the length of the vectors which were read (i.e. how many pairs were read); `tdvvto` need not report anything under normal execution. Both testing programs should be adequately documented and must perform argument processing—this includes detection of insufficient or invalid arguments, in which case a usage message should be printed to standard error. Use the built-in routines `iargc` and `getarg` as well as `i4arg` from the `p329f` library for argument parsing. You should also use the integer function `indlnb` from the `p329f` library to ensure that you don't create file names with trailing spaces. I will test your test programs, as well as your `dvvfrom` and `dvvto` routines themselves, with my own input and testing programs.

Problem 3: In directory `~/hw3/a3` create a source file `rw2d.f` and executable `rw2d` which simulates a random walk of a particle on a two dimensional integer lattice. Specifically, if at step n the current position of the particle performing the walk is $(x, y) = (i, j)$, where i and j are integers, then, with equal probability (i.e. $p = 1/4$), the position at step $n+1$ is either $(i+1, j)$, $(i-1, j)$, $(i, j+1)$ or $(i, j-1)$. `rw2d` should accept a single integer argument which specifies the number of steps to perform in the simulation. Start the particle at $(0, 0)$ and use `real*8` values for the particle coordinates, even though the coordinates are restricted to integer values. After each step, including initialization, the program should write to standard output the step number, the current coordinates of the particle, and the particle's current distance, $d = \sqrt{x^2 + y^2}$, from its initial position. Specifically, use a `write` statement like

```
write(*,*) step, x, y, d
```

Document and test your program thoroughly, then run it for 10000 steps and save the output in the file `out10000`. Use `gnuplot` or `maple` to produce a plot of d versus step number from the data in `out10000`. Save a postscript version of the plot in `distance.ps`. Make another plot (postscript version `distance2.ps`) which graphs d^2 as a function of step number. Note than on the SGIs, the built in `real*8` function `rand` can be used to generate a random number between 0 and 1—it is up to you to figure out how to use `rand` to generate the particular randomness required by the simulation. Be sure that you declare the function properly in the main program, and note that since it has no arguments, it should be invoked with an empty argument list:

```
real*8      rand

write(*,*) 'This is a pseudo-random number ', rand()
```

Problem 4: *The Chaos Game*
Let

$$V_i \equiv (\tilde{x}_i, \tilde{y}_i) \qquad i = 1, 2, 3$$

be the vertices of an equilateral triangle with side length 1 lying in the xy plane. The orientation and position of the triangle in the plane is not important but, for concreteness, you may wish to "center" the triangle at the origin $(x, y) = (0, 0)$. The game is initialized by choosing, *at random*, some point in the xy plane, such that the point lies outside of the triangle. Call this randomly chosen location the *active point*. One then *randomly* chooses one of the three triangle vertices, V_i, draws an imaginary line between the vertex and the active point, bisects the line and makes the point of bisection the active point. This last sequence of randomly choosing a vertex, bisecting the line between it and the active point and updating the active point with the point of bisection is then repeated some specified number of times. If one plots all the active points which are generated, an interesting pattern appears after a sufficiently large number of iterations.

In directory ~/hw3/a4 on your SGI account write a well-documented Fortran program called `chaos2d` which implements this algorithm. As with the previous problem, you will probably find use for the `real*8` random number function, `rand`. The program should take a single integer argument which specifies the number of iterations to perform. Again, I suggest that you use the `i4arg` function discussed in class and available in the `p329f` library to parse the argument. The program should write to standard output the x and y coordinates of all of the active points which are generated (including the original randomly chosen one). Specifically, use a `write` statement like the following

```
write(*,*) xactive, yactive
```

in your iteration loop, so that the coordinate pairs appear one per line on standard output. You may optionally output (also to standard output) the (x, y) coordinates of the three triangle vertices V_i. If you choose to do so, output them *once only*, presumably at the beginning of the game. Finally, although the initial active point is to be chosen randomly, ensure that its distance from the center of the triangle does not exceed 2.5 (recall that the sides of the triangle are of length 1.0). Run your program for 10, 100, 1000 and 10000 iterations and redirect the output to files `out10, out100, out1000,` and `out10000` respectively. Use `gnuplot` or `maple` to plot the points defined by each of the output files and save the plots as postscript files `out10.ps, out100.ps, out1000.ps,` and `out10000.ps`