Exercises Lecture III:

Random numbers with non uniform distributions; simulations of simple random processes

homework: n. 1,3,4

- 1. Random numbers with non uniform distributions: Inverse Transformation Method
 - (a) With the Inverse Transformation Method we can generate random numbers according to the exponential distribution $f(z) = \lambda e^{-\lambda z}$, starting from random numbers with uniform distribution: if x is the random variable with uniform distribution in [0,1], then z = -ln(x) is distributed according to e^{-z} . Write a code implementing the algorithm. An example is given in expdev.f90.
 - (b) Check—doing a histogram—that the random variate z generated with that algorithm is actually exponentially distributed.

 (What is convenient to plot in order to check this behavior? Hint: with gnuplot you can print the log of your data (e.g., suppose you saved the values of z in column 1 and its frequency in column 2, plot with u 1:(log(\$2)) or u 1:(log10(\$2))).
 - (c) With gnuplot you can also do the fit of the histogram with an exponential function using the least-square method, with λ as fitting parameter. Check whether you get the expected value of λ . (It is convenient to make a semilog plot as suggested above and then make a least-square linear fit; the slope is λ)

Remember that with the method of the least-square fit we get for a linear regression: y = ax + b:

$$a = \frac{\overline{xy} - \overline{yy}}{(\Delta x)^2}; \qquad b = \overline{y} - a\overline{x}$$

where $(\Delta x)^2 = \overline{x^2} - \overline{x}^2$ (other definitions are trivial ...).

2. Random numbers with non uniform distributions: comparison between different algorithms

Suppose you want to generate a random variate x in (-1,1) with distribution

$$p(x) = \frac{1}{\pi} (1 - x^2)^{-1/2}.$$

Consider both methods suggested below, do the histograms and check that both methods give correct results.

- (a) From the Inverse Transformation Method: generate a random number U with uniform distribution in [0,1] and consider $x = \sin \pi (2U 1)$.
- (b) Generate two random numbers U and V with uniform distribution in [0,1]. Disregard them if $U^2 + V^2 > 1$. Otherwise consider

$$x = \frac{U^2 - V^2}{U^2 + V^2}$$

Note 1: the last method has the advantage of using only elementary operations.

Note 2: since x is also negative, pay attention to the algorithm used to make the histogram; you should notice the difference between the intrinsic functions int and nint; see also floor. From Chapman's book:

AINT(A,KIND): Real elemental function

- Returns A truncated to a whole number.

AINT(A) is the largest integer which is smaller than |A|, with the sign of A. For example, AINT(3.7) is 3.0, and AINT(-3.7) is -3.0.

- Argument A is Real; optional argument KIND is Integer

ANINT(A, KIND): Real elemental function

- Returns the nearest whole number to A.

For example, ANINT(3.7) is 4.0, and AINT(-3.7) is -4.0.

- Argument A is Real; optional argument KIND is Integer

FLOOR(A, KIND): Integer elemental function

- Returns the largest integer < or = A.

For example, FLOOR(3.7) is 3, and FLOOR(-3.7) is -4.

- Argument A is Real of any kind; optional argument KIND is Integer
- Argument KIND is only available in Fortran 95

NINT(A[,KIND])

- Integer elemental function
- Returns the nearest integer to the real value $\mbox{\mbox{\sc A}}.$
- A is Real

3. Random numbers with gaussian distribution: Box-Muller algorithm

Consider the Box-Muller algorithm to generate a random number gaussian distribution (see for instance boxmuller.f90; the gasdev subroutine used inside is similar to what you can find in "Numerical Recipes": it gives a gaussian distribution with $\sigma=1$ and average $\mu=0$). Do a histogram of the data generated, calculate numerically from the sequence the average value and the variance, check with the expected results.

4. Simulation of radioactive decay

- (a) Write a program for a numerical simulation of the radioactive decay, with a decay parameter λ in input. (See for instance decay.f90).
- (b) Use the code with "reasonable" values of the parameters (e.g., N(0) about 1000) and save N(t) in a data file. Check whether $N(t) = N(0)e^{-\lambda t}$ as expected. (Hint: As for the exercise 1, you could make use of a least-square fit by considering $\ln N(t)$ vs. t, i.e. the relationship in a semilog form in order to manage a linear fit.)
- (c) Change N(0) (100 or less; 10000 or more). What do you see?

Notice that in decay.f90 the upper bound of the inner loop (nleft) is changed within the execution of the loop; but in the execution the loop goes on up to the nleft set at the beginning of the loop; this ensures that the implementation of the algorithm is correct. See the programs checkloop.f90 and decay_checkloop.f90 in the same directory.

5. Random deviates with other distributions (Optional)

You can try t_random.f90 which uses the module random.f90 to generate random deviates with other distributions. Remember to compile first the module: g95 (or gfortran) random.f90 t_random.f90

```
print *," length of the sequence >"
 read *, n
 print *," exponential decay factor (lambda)>"
 read *, lambda
 print *," Collecting numbers generated up to 2/lambda (disregard the others)"
 print *," and normalizing the distribution in [0,+infinity[ "
 print *," Insert number of bins in the histogram>"
 read *, nbin
 delta = 2./lambda/nbin
   allocate (histo(nbin))
 histo = 0
 do i = 1,n
    call expdev(x)
    ibin = int (x/lambda/delta) + 1
    if (ibin <= nbin)histo(ibin) = histo(ibin) + 1</pre>
 open (unit=7,file="expdev.dat",status="replace",action="write")
 do ibin= 1 ,nbin
    write(unit=7,fmt=*)(ibin-0.5)*delta,histo(ibin)/float(n)/delta
  end do
contains
 subroutine expdev(x)
   REAL, intent (out) :: x
   REAL :: r
      call random_number(r)
      if(r > 0) exit
   end do
   x = -log(r)
 END subroutine expdev
end program test_expdev
! boxmuller.90
! uses the Box-Muller algorithm to generate
! a random variate with a gaussian distribution (sigma = 1)
program boxmuller
 implicit none
 real :: rnd,delta
 real, dimension(:), allocatable :: histog
 integer :: npts,i,ibin,maxbin,m
```

```
print*,' input npts, maxbin >'
 read*, npts,maxbin
 allocate(histog(-maxbin/2:maxbin/2))
 histog = 0
 delta = 10./maxbin
 do i = 1, npts
     call gasdev(rnd)
     ibin = nint(rnd/delta)
     if (abs(ibin) < maxbin/2) histog(ibin) = histog(ibin) + 1
  end do
 open(1,file='gasdev.dat',status='replace')
  do ibin = -maxbin/2, maxbin/2
     write(1,*)ibin*delta, histog(ibin)/real(npts)/delta
  end do
  close(1)
 deallocate(histog)
 stop
contains
 SUBROUTINE gasdev(rnd)
    IMPLICIT NONE
    REAL, INTENT(OUT) :: rnd
    REAL :: r2,x,y
    REAL, SAVE :: g
   LOGICAL, SAVE :: gaus_stored=.false.
    if (gaus_stored) then
       rnd=g
       gaus_stored=.false.
    else
       do
          call random_number(x)
          call random_number(y)
          x=2.*x-1.
          y=2.*y-1.
          r2=x**2+y**2
          if (r2 > 0. .and. r2 < 1.) exit
       end do
       r2=sqrt(-2.*log(r2)/r2)
       rnd=x*r2
       g=y*r2
       gaus_stored=.true.
    end if
 END SUBROUTINE gasdev
end program boxmuller
```

```
! decay.f90
! Simulation of radioactive decay
PROGRAM decay
 IMPLICIT none
 REAL, PARAMETER ::
                         lambda=0.2
 REAL :: r
 INTEGER :: i, t, nleft, start, sizer
 integer, dimension(:), allocatable :: seed
 call random_seed(sizer)
 allocate(seed(sizer))
 print *,'Here the seed has ',sizer,' components; insert them (or print "/") >'
 read(*,*)seed
 call random_seed(put=seed)
           initial values
 print *,"initial number of nuclei >"
 read *, start
 t = 1
               ! initialize time
 nleft = start ! at the beginning N(t=0)=start
 ! N(t) nuclei left at time t,
  ! that have a given probability lambda of decay
  ! in the time interval t:t+dt
OPEN(unit=7, FILE="decay.dat", status="replace",action="write")
 WRITE (unit=7,fmt=*) "# t ,
                                N(t)"
 WRITE (unit=7,fmt=*) "0 ", nleft !REAL(nleft)/start
DO
                                ! time loop
    DO i = 1, nleft
                                 ! loop on the nuclei left
       call random_number(r)
       IF (r <= lambda) THEN</pre>
          nleft = nleft - 1
                                 ! update the number of nuclei left
       ENDIF
    END DO
    1
    WRITE (unit=7,fmt=*) t , nleft ! or REAL(nleft)/start
    if (nleft == 0) exit
    t = t + 1
 END DO
  !
 close(7)
 stop
END program decay
```