

## Reading Chapter 1

2/16/25

### 1.1 Radioactive Decay

- A typical example of nuclear isotope  $^{235}\text{U}$  (uranium nucleus that contains 143 neutrons and 92 protons).

- Process of radioactive decay is random in the following sense.

- It is useful to imagine that we have a sample containing a large number of  $^{235}\text{U}$  nuclei.

- If  $N_U(t)$  is the number of U nuclei that are present in a sample at time  $t$ .

$$\frac{dN_U}{dt} = -\frac{N_U}{\tau}$$

where  $\tau$  is the "time constant" for decay. You can then show by direct sub:

$$N_U = N_U(0)e^{-t/\tau}$$

### 1.2 A Numerical Approach

- We can now consider a simple method for solving this problem.

- Our goal is to obtain  $N_i$  as a function of  $t$ .

- Given the value of  $N_i$  at a particular value of  $t$ , we want to estimate the value at later times (initial value problem).

- We can use Taylor expansion for  $N_i$ :

$$N_i(\Delta t) = N_i(0) + \frac{dN_i}{dt} \Delta t + \frac{1}{2} \frac{d^2 N_i}{dt^2} (\Delta t)^2 + \dots$$

where  $N_i(0)$  is the value of our func. at time  $t=0$ ,  $N_i(\Delta t)$  is its value at  $t=\Delta t$ , and derivatives are evaluated at  $t=0$ .

- While programming, like handwriting, is a highly individual process, there are certain recommended practices.

- It is important that we be able to understand programs written by others, as well as by ourselves.

- The first thing you should do is think. Construct an outline of how the problem is to be solved.

- Example 1.1

- Declare necessary variables + arrays

- Initialize variables



- Do actual calculation.

- Store results.

- Note that this is the main program.

- Consider how this looks in Fortran.

C. Simulation of radioactive decay.

C Program to accompany "Computational Phys"  
by M. Gordon & H. Nakanishi

Program decay

C. declare arrays

double precision n-uranium(100), t(100)

C use subroutines to do the work

call initialize(n-uranium, t, tau, dt, n)

call calculate(n-uranium, t, tau, dt, n)

call store(n-uranium, t, n)

stop

end

Example 1.2 Pseudocode for subroutine

- Prompt for and assign  $N(0)$ ,  $\tau$ , and  $\Delta t$ .

- Set initial value of time,  $t(0)$ .

- Set numbers of time steps for calc.

- Fortran version of subroutine.

```

Subroutine initialize(nuclei, t, tc, dt, n)
  Initialize variables
  double precision nuclei(n), t(n)
  Print *, 'initial number of nuclei → '
  C read(s, *) nuclei(1)
  Print *, 'time constant → '
  read(s, *) tc
  Print *, 'time step → '
  read(s, *) dt
  Print *, 'total time → '
  read(s, *) time
  t(1) = 0
  n = min(int(time/dt), 100)
  return
end

```

### Example 1.3

- For each time step  $i$  (beginning with  $i=1$ ), calculate  $N_i$  and  $t$  at step  $i+1$ :
  - $N_i(t_{i+1}) = N_i(t_i) - (N_i(t_i)/\tau) dt$  (Euler method)
  - $t_{i+1} = t_i + \Delta t$ .
  - repeat for  $n-1$  time steps.

Subroutine calculate in Fortran:

```

Subroutine calculate(n-uranium, t, tau, dt, n)
  C Now use Euler method
  C Variable dimensioning is used for
  arrays n-uranium(n) and t(n) do i=1, n-1
    n-uranium(i+1) = n-uranium(i) - (n-uranium(i)/tau) dt
    t(i+1) = t(i) + dt

```



```
end do  
return  
end
```

- Final store writes the result to a file.

Subroutine store in Fortran

```
Subroutine store (n-uranium, t, n)  
double Precision n-uranium(n), t(n)  
Open(i, file='decay.dat')  
do i=1, n  
    write(i, 20), t(i), n-uranium(i)  
end do  
close(i)  
format(1x, 1p, 2(e15.5, 2x))  
return  
end
```

Graph:

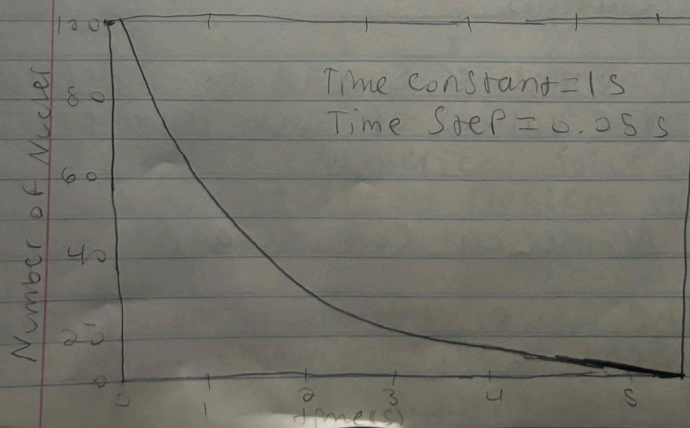


Figure 1.1: Circles indicate the numerical solution to the radioactive decay problem.

#### 1.4 Testing Your Program.

- We should not really consider it to be a working function until we are convinced that its output is correct.
- Checking a program is not always a trivial task, but there are some general guidelines.
- After a program has been debugged so that it can be run without any complaints about the syntax.
- Always check your program: gives the same answer for different "step sizes". Our decay program involved a timestep variable,  $dt$ , and most other numerical calculations.
- Final answer should be independent of the values of such parameters.

#### 1.5 Numerical Considerations

- The issue of numerical errors is central to the computational solution of any problem.



- Questions such as how to design or choose the best algorithm for a particular problem are central topics in many computer science and applied math courses.

- With our radioactive decay program errors were introduced by the approx used to estimate the solution of the diff eq.

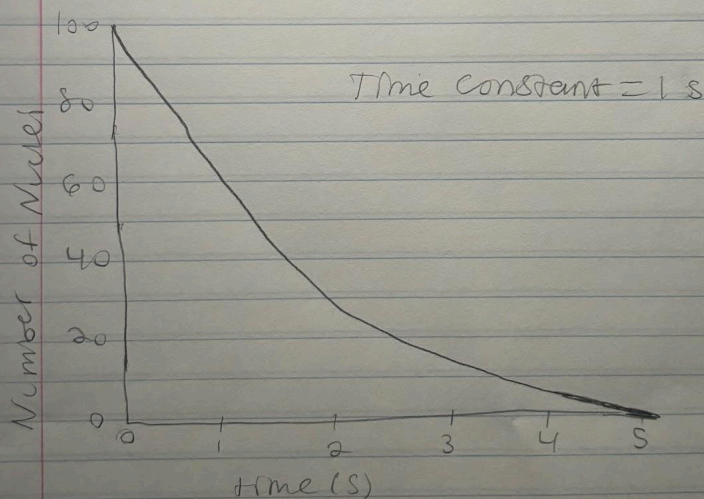


Figure 1.2: Numerical solution of the radioactive decay problem using the Euler method.