Programming Assignment 1

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Due Wed 2/5/25

Problem 1: Finding Prime Numbers

Method 1

```
In [4]: import time

t0 = time.time()

primes = [2]
for i in range(3, 100001):
    isPrime = True
    for f in range(2, i):
        if (i%f == 0):
            isPrime = False
    if isPrime:
        primes.append(i)

tf = time.time()
print("Runtime: {:.1f} s".format(tf-t0))
```

Runtime: 198.8 s

Method 2

- Skip evens greater than 2
- Only check divisibility of primes less than i
- Only check factors up to \sqrt{i}
- Break once you find the first factor

```
In [25]: import math
import time

t0 = time.time()

primes2 = [2]
for i in range(3, 100001):
    isPrime = True
    # Skip this i if even
    if (i%2 == 0):
```

```
continue

# Check i's divisibility by primes less than it (contents of primes2), u
for p in primes2:
    if (p > math.sqrt(i)):
        break
    if (i%p == 0):
        isPrime = False
        break

# Add to list
if isPrime:
    primes2.append(i)

tf = time.time()
print("Runtime: {:.1f} s".format(tf-t0))
```

Runtime: 0.1 s

Method 1 takes ~10^4 times longer to run.

Problem 2: Nuclear Binding Energy

Part A

```
In [16]: A = int(input("Enter A: "))
Z = int(input("Enter Z: "))
print(f"A={A}, Z={Z}")

a1, a2, a3, a4 = 15.8, 18.3, 0.714, 23.2
if A%2!=0:
    a5 = 0
elif A%2==0 and Z%2==0:
    a5 = 12
else:
    a5 = -12

B = a1*A - a2*A**(2/3) - a3*Z**2 / A**(1/3) - a4*(A-2*Z)**2 / A + a5 / A**(1
print(f"B={B}")
A=58, Z=28
```

Part B

B=497.5620206224374

```
In [7]: print(f"Binding energy per nucleon: " + str(B/A))
```

Binding energy per nucleon: 8.578655527973059

Part C

```
In [22]: Z = int(input("Enter Z: "))
         print(f"Z={Z}")
         a1, a2, a3, a4 = 15.8, 18.3, 0.714, 23.2
         Bper_max = 0
         Aper max = 0
         for A in range(Z, 3*Z+1):
             # Need to recalc a5
             if A%2!=0:
                 a5 = 0
             elif A%2==0 and Z%2==0:
                 a5 = 12
             else:
                 a5 = -12
             # Calculate binding energy B
             B = a1*A - a2*A**(2/3) - a3*Z**2 / A**(1/3) - a4*(A-2*Z)**2 / A + a5 / A
             # ... per nucleon
             Bper = B/A
             # Store Bper and A if these yielded the highest B
             if Bper>Bper max:
                 Bper_max = Bper
                 Aper_max = A
         print(f"Max binding energy per nucleon (Bper): {Bper_max}
                                                                        Mass (A): {Aper
        Z = 26
```

Part D

```
In [28]: a1, a2, a3, a4 = 15.8, 18.3, 0.714, 23.2
         most_stable_B, most_stable_Z = 0, 0
         for Z in range(1, 101):
              Bper_max = 0
             Aper_max = 0
              for A in range(Z, 3*Z+1):
                 # Need to recalc a5
                 if A%2!=0:
                      a5 = 0
                 elif A%2==0 and Z%2==0:
                      a5 = 12
                 else:
                      a5 = -12
                 # Calculate binding energy B
                 B = a1*A - a2*A**(2/3) - a3*Z**2 / A**(1/3) - a4*(A-2*Z)**2 / A + a5
                 # ... per nucleon
                  Bper = B/A
```

Max binding energy per nucleon (Bper): 8.701432576808985

Mass (A): 58

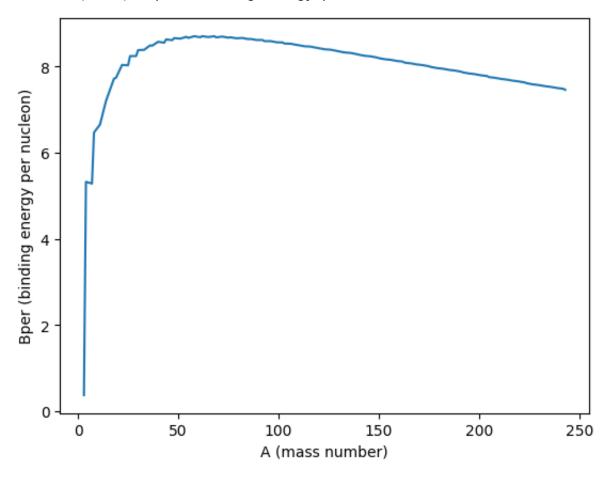
Most stable Z: 28 Binding energy per nucleon: 8.70245768367189

Part E

```
In [40]: # For every element Z,
             # find the most stable nucleus (largest Bper)
             # store the Bper and A
         # Graph Bper vs A
         import matplotlib.pyplot as plt
         # from importlib import reload
         # plt=reload(plt)
         Bper_max_list = []
         Aper_max_list = []
         a1, a2, a3, a4 = 15.8, 18.3, 0.714, 23.2
         for Z in range (1, 100):
             Bper_max = 0
             Aper max = 0
             for A in range(Z, 3*Z+1):
                 # Need to recalc a5
                 if A%2!=0:
                     a5 = 0
                 elif A%2==0 and Z%2==0:
                     a5 = 12
                 else:
                     a5 = -12
                 # Calculate binding energy B
                 B = a1*A - a2*A**(2/3) - a3*Z**2 / A**(1/3) - a4*(A-2*Z)**2 / A + a5
                 # ... per nucleon
                 Bper = B/A
                 # Update Bper_max and Aper_max if this A yielded the highest Bper
                 if Bper>Bper max:
                     Bper_max = Bper
                     Aper max = A
             # Store this Z's Bper_max and Aper_max in the lists
             Bper max list.append(Bper max)
             Aper_max_list.append(Aper_max)
```

```
In [41]: plt.plot(Aper_max_list, Bper_max_list)
  plt.xlabel("A (mass number)")
  plt.ylabel("Bper (binding energy per nucleon)")
```

Out[41]: Text(0, 0.5, 'Bper (binding energy per nucleon)')



Elements that **release** energy during **fusion**: Light elements -- has a mass number (A) less than about 50

Elements that **release** energy during **fission**: Heavy elements -- has a mass number (A) greater than about 50

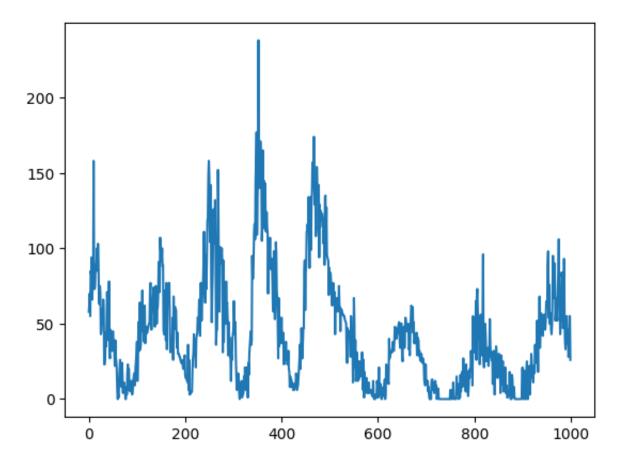
More energy released in *fusion of light nuclei* or *fission of heavy nuclei*: fusion of light nuclei, because the graph is much steeper in the left region (light elements)

Problem 3: Running Avg of Dataset

Part A

```
import numpy as np
import matplotlib.pyplot as plt
sunspots_data = np.loadtxt(fname="sunspots.txt", delimiter=" ")
t, n = sunspots_data[:1000, 0], sunspots_data[:1000, 1]
plt.plot(t,n)
```

Out[10]: [<matplotlib.lines.Line2D at 0x123ba2570>]

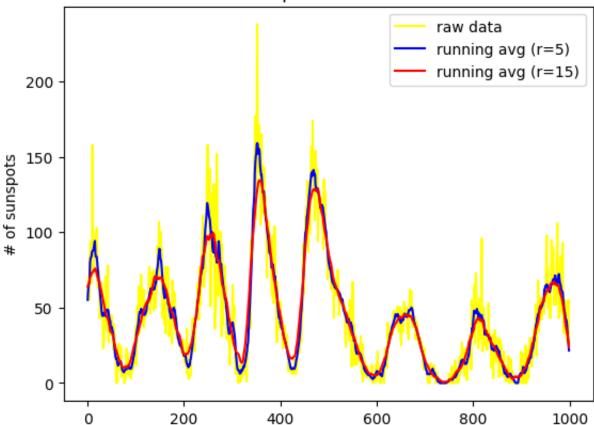


Part B

```
In [37]: # create function to find running avg data for different r values
         def running_avg(r):
             nravgs = []
             for k in range(len(n)):
                 summation = 0
                 for m in range(-r, r+1):
                      # if we run out of data towards the end for the summation, just
                      try:
                          summation += n[k+m]
                      except:
                          continue
                 Yk = 1/(2*r+1) * summation
                 nravgs.append(Yk)
             return nravgs
         # use function to find running avg data for different r values
         nravgs_r5 = running_avg(r=5)
         nravgs_r15 = running_avg(r=15)
         # plot
         fig, ax = plt.subplots()
         ax.plot(t,n, color="yellow", label="raw data")
         ax.plot(t,nravgs_r5, color="blue", label="running avg (r=5)")
```

```
ax.plot(t,nravgs_r15, color="red", label="running avg (r=15)")
ax.legend()
ax.set_ylabel("# of sunspots")
ax.set_title("Sunspots Over Time")
plt.show()
```





Increasing r smooths out the curve more.

Problem 4 (optional): Mandelbrot Set

Black/White Binary

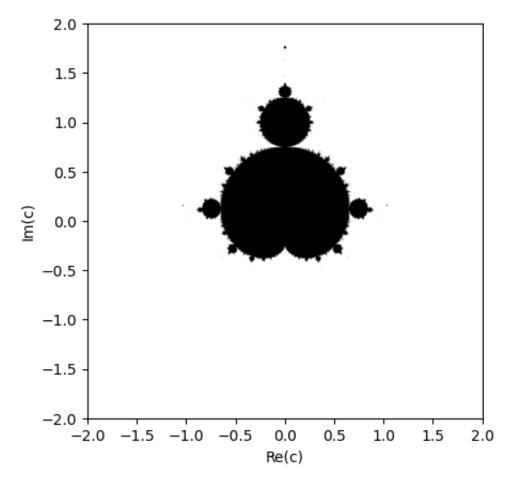
- Create grid/set of c's
- Create array mandelbrot_map
- · For each c:
 - Start with z=0, and do the following 100 (max_iter) times

$$\circ z + = z^2 + c$$

- $\circ \,\,$ If |z|>2 at any point, set mandelbrot(x,y) to 0
- Else keep mandelbrot_map(x,y) to 1
- Plot mandelbrot_map

```
In [13]: import numpy as np
         import matplotlib.pyplot as plt
         N = 1000
                         # resolution
         max_iter = 100 # iters to check
         Re, Im = np.linspace(-2,2,N), np.linspace(-2,2,N)
         m_m = np.ones((N,N))
         \# c = Re + Im*i
         for x in range(N):
             for y in range(N):
                 c = Re[x] + Im[y]*1j
                 z = 0
                 for iters in range(max_iter):
                     z = z**2 + c
                     if abs(z) > 2:
                         m_map[x, y] = 0
                         break
         plt.imshow(m_map, cmap="binary", extent=[-2,2,-2,2])
         plt.xlabel("Re(c)")
         plt.ylabel("Im(c)")
```

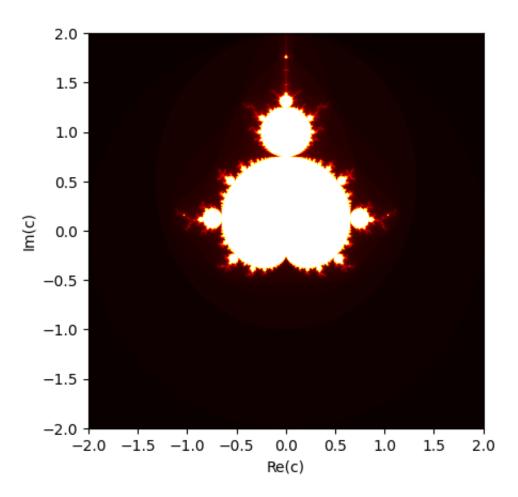
Out[13]: Text(0, 0.5, 'Im(c)')



Gradient based on Iterations

- · Create grid/set of c's
- Create array mandelbrot_map
- For each c:
 - Start with z=0, and do the following 100 (max_iter) times
 - $\circ z + = z^2 + c$
 - \circ If |z| > 2, log number of iters in m_map
 - If z never >2, log max_iters in m_map
- Plot mandelbrot_map

```
In [ ]: import numpy as np
        import matplotlib.pyplot as plt
        N = 1000
                   # resolution
        max iter = 100 # iters to check
        L = 2 # "side length" of view window, centered at 0,0
        def find_iters(c):
            z = 0
            for iters in range(max iter):
                z = z**2 + c
                if abs(z) > 2:
                    return iters
            return max_iter
        Re, Im = np.linspace(-L,L,N), np.linspace(-L,L,N)
        m_map = np.zeros((N,N))
        \# c = Re + Im*i
        for x in range(N):
            for y in range(N):
                c = Re[x] + Im[y]*1j
                c_iters = find_iters(c)
                m_map[x, y] = c_iters
        plt.imshow(m_map, cmap="hot", extent=[-L,L,-L,L])
        plt.xlabel("Re(c)")
        plt.ylabel("Im(c)")
        plt.show()
```



Gradient based on log(iters)

```
In [30]:
         import numpy as np
         import matplotlib.pyplot as plt
         import math
                         # resolution
         N = 1000
         max_iter = 100 # iters to check
         L = 2
                 # "side length" of view window, centered at 0,0
         def find_iters(c):
             z = 0
             for iters in range(max_iter):
                 z = z**2 + c
                 if abs(z) > 2:
                      return iters
             return max_iter
         Re, Im = np.linspace(-L,L,N), np.linspace(-L,L,N)
         m_m = np.zeros((N,N))
         \# c = Re + Im*i
         for x in range(N):
             for y in range(N):
                 c = Re[x] + Im[y]*1j
```

```
c_iters = find_iters(c)
    m_map[x, y] = math.log(c_iters) if c_iters>0 else 0

plt.imshow(m_map, cmap="hot", extent=[-L,L,-L,L])
plt.xlabel("Re(c)")
plt.ylabel("Im(c)")
plt.show()
```

