## Astro 410 Extra-Credit HomeWork 4

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List of included files:

- hw4-dubey.ipynb
- hw4-dubey.py (Python code not as a Jupyter Notebook)
- hw4-dubey.pdf (contains printed pdf of html file)
- hw4-dubey.html (contains html version of Jupyter Notebook for better readability)

(a lot of the inspiration for the code has been taken from the leap-frog.c file on canvas and various internet portals, although I have done my best to mold it to suit the assignment's needs by disecting it the best I can and annotating for easy understanding)

The following report attempts to approximate the behaviours of N-bodies in orbit around each other.

We begin by calculating acceleration.

```
In [1]: import numpy as np
         import matplotlib.pyplot as plt
         def getAcc( pos, mass, G, softening ):
             # positions r = [x,y,z] for all particles
                 x = pos[:,0:1]
                 y = pos[:,1:2]
                 z = pos[:,2:3]
                 # matrix that stores all pairwise particle separations: r j - r i
                 dx = x.T - x
                 dy = y.T - y
                 dz = z.T - z
                 # matrix that stores 1/r^3 for all particle pairwise particle separati
         ons
                 inv r3 = (dx^{**2} + dy^{**2} + dz^{**2} + softening^{**2})
                 inv r3[inv r3>0] = inv r3[inv r3>0]**(-1.5)
                 ax = G * (dx * inv_r3) @ mass
                 ay = G * (dy * inv_r3) @ mass
                 az = G * (dz * inv r3) @ mass
                 # pack together the acceleration components
                 a = np.hstack((ax,ay,az))
                 return a
```

Building a function to calculate KE and PE seperately for each body. Where formula for KE is:

$$\frac{1}{2}\sum_{i}mv^{2}$$

And potential energy (PE) is calculated as below:

$$rac{1}{2}\sum_{1 < i < j < N} rac{Gm_im_j}{|r_j - r_i|}$$

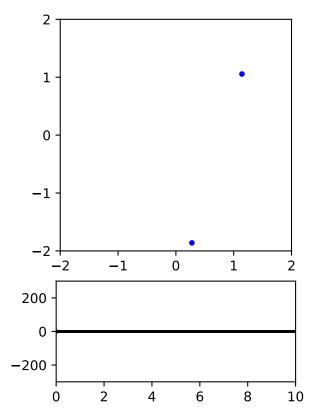
```
In [3]: def getEnergy( pos, vel, mass, G ):
            # Kinetic Energy:
            # kinetic energy of the system
            KE = 0.5 * np.sum(np.sum( mass * vel**2 ))
            # Potential Energy:
            # positions r = [x,y,z] for all particles
            x = pos[:,0:1] # Selecting Column 1 from pos matrix
            y = pos[:,1:2] # Selecting Column 2 from pos matrix
            z = pos[:,2:3] # Selecting Column 3 from pos matrix
            # matrix that stores all pairwise particle separations: r_j - r_i
            dx = x.T - x # Transpose of array X - array X
            ####################### where each position X is subtracted from every other
            dz = z.T - z # X in the matrix
            dy = y.T - y # Same with Y and Z to make a 2D matrix
            # matrix that stores 1/r for all particle pairwise particle separations
            inv r = np.sqrt(dx**2 + dy**2 + dz**2)
            inv r[inv r > 0] = 1.0 / inv r[inv r > 0]
            # sum over upper triangle, to count each interaction only once
            # potential energy of the system
            PE = G * np.sum(np.sum(np.triu(-(mass*mass.T)*inv r,1)))
            # np.triu function: Return a copy of an array with the elements below the
         k-th diagonal zeroed
            return KE, PE
```

To run the simulation, we need to first input initial parameters, for example, gravitaitonal constant G, the number of bodies N and so on as listed below. Tweaking these values allow us to view a wide variety of variations by changing one or a combination of parameters at a time. Currently, the code does not run a simulation for N bodies, but instead for 2 particles mimicking the Earth-Sun orbit. Uncommenting the lines above the specified mass will let us add more bodies in the simulation.

```
In [19]: # Simulation parameters
                 = 2
                         # Number of particles
         N
         t
                  = 2
                         # current time of the simulation
                  = 10.0 # time at which simulation ends
         tEnd
                  = 0.01 # timestep
         dt
         softening = 0.1 # softening length
                  = 1.0
                           # Newton's Gravitational Constant
         plotRealTime = True # switch on for plotting as the simulation goes along
         # Generate Initial Conditions
         np.random.seed(17)
                                     # set the random number generator seed
         # mass = 20.0*np.ones((N,1))/N # total mass of particles is 20
         mass = np.array(([19.999 999 9], [0.000\ 000\ 1]))/N # 20.0*np.ones((N,1))/N #
          total mass of particles is 20
         pos = np.random.randn(N,3) # randomly selected positions and velocities
         pos[0][0] = 0 # This is so that the first particle has x = 0
         pos[0][1] = 0 # y = 0
         pos[0][2] = 0 # and z = 0
         pos[1][0] = 1 # This is so that the first particle has x = 0
         pos[1][1] = 0 # y = 0
         pos[1][2] = 0 # and z = 0
         vel = np.random.randn(N,3)
         # Convert to Center-of-Mass frame
         vel -= np.mean(mass * vel,0) / np.mean(mass)
         # calculate initial gravitational accelerations
         acc = getAcc( pos, mass, G, softening )
         # calculate initial energy of system
         KE, PE = getEnergy( pos, vel, mass, G )
         # number of timesteps
         Nt = int(np.ceil(tEnd/dt))
         # save energies, particle orbits for plotting trails
         pos save = np.zeros((N,3,Nt+1))
         pos save[:,:,0] = pos
         KE save = np.zeros(Nt+1)
         KE_save[0] = KE
         PE save = np.zeros(Nt+1)
         PE save[0] = PE
         t all = np.arange(Nt+1)*dt
         # prep figure
         fig = plt.figure(figsize=(4,5), dpi=80)
         grid = plt.GridSpec(3, 1, wspace=0.0, hspace=0.3)
         ax1 = plt.subplot(grid[0:2,0])
         ax2 = plt.subplot(grid[2,0])
         # Simulation Main Loop
         for i in range(Nt):
             #(1/2) kick
             vel += acc * dt/2.0
```

```
# drift
   pos += vel * dt
   # update accelerations
   acc = getAcc( pos, mass, G, softening )
   #(1/2) kick
   vel += acc * dt/2.0
   # update time
   t += dt
   # get energy of system
   KE, PE = getEnergy( pos, vel, mass, G )
   # save energies, positions for plotting trail
   pos save[:,:,i+1] = pos
   KE save[i+1] = KE
   PE save[i+1] = PE
   # plot in real time
   if plotRealTime or (i == Nt-1):
       # Scatter plot following N bodies
       plt.sca(ax1)
        plt.cla()
       xx = pos_save[:,0,max(i-50,0):i+1]
       yy = pos_save[:,1,max(i-50,0):i+1]
        plt.scatter(xx,yy,s=1,color=[.7,.7,1])
       plt.scatter(pos[:,0],pos[:,1],s=10,color='blue')
        ax1.set(xlim=(-2, 2), ylim=(-2, 2))
        ax1.set_aspect('equal', 'box')
        ax1.set_xticks([-2,-1,0,1,2])
        ax1.set_yticks([-2,-1,0,1,2])
        # Scatter plot following KE, PE and Totale Energy
        plt.sca(ax2)
        plt.cla()
        plt.scatter(t all,KE save,color='red',s=1,label='KE' if i == Nt-1 else
        plt.scatter(t all,PE save,color='blue',s=1,label='PE' if i == Nt-1 els
e "")
       plt.scatter(t_all,KE_save+PE_save,color='black',s=1,label='Etot' if i
== Nt-1 else "")
        ax2.set(xlim=(0, tEnd), ylim=(-300, 300))
        ax2.set aspect(0.007)
        plt.pause(0.001)
# add Labels/Legend
plt.sca(ax2)
plt.xlabel('time')
plt.ylabel('energy')
ax2.legend(loc='upper right')
```

```
# Save figure
plt.savefig('nbody.png',dpi=240)
plt.show()
```



```
ValueError
                                           Traceback (most recent call last)
<ipython-input-19-1145540d4350> in <module>
            # plot in real time
            if plotRealTime or (i == Nt-1):
     70
---> 71
                plt.sca(ax1)
     72
                plt.cla()
                xx = pos_save[:,0,max(i-50,0):i+1]
     73
~\AppData\Local\Programs\Python\Python37\lib\site-packages\matplotlib\pyplot.
py in sca(ax)
    856
                    m.canvas.figure.sca(ax)
    857
                    return
            raise ValueError("Axes instance argument was not found in a figur
--> 858
e")
    859
    860
```

ValueError: Axes instance argument was not found in a figure