

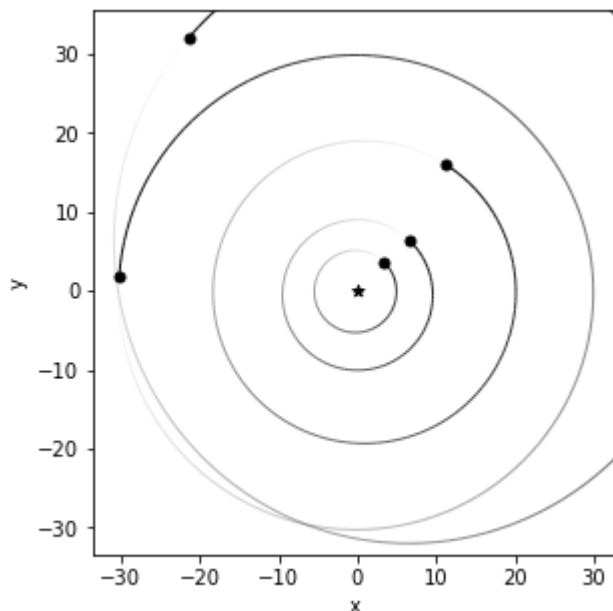
# Integrations with REBOUND

```
In [1]: import numpy as np
import matplotlib.pyplot as plt
import rebound
```

```
In [2]: sim = rebound.Simulation()
rebound.data.add_outer_solar_system(sim)
```

```
In [3]: rebound.OrbitPlot(sim)
```

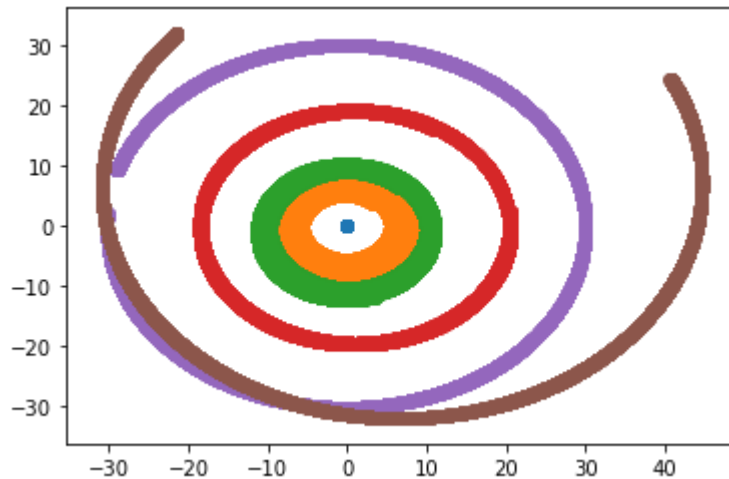
```
Out[3]: (<Figure size 360x360 with 1 Axes>,
<matplotlib.axes._subplots.AxesSubplot at 0x7fe4f223e2d0>)
```



```
In [4]: def eulerstep(sim):
sim.integrator = "None"
sim.step() # Calculates the accelerations
for i in range(sim.N):
    sim.particles[i].x += sim.dt * sim.particles[i].vx
    sim.particles[i].y += sim.dt * sim.particles[i].vy
    sim.particles[i].z += sim.dt * sim.particles[i].vz
    sim.particles[i].vx += sim.dt * sim.particles[i].ax
    sim.particles[i].vy += sim.dt * sim.particles[i].ay
    sim.particles[i].vz += sim.dt * sim.particles[i].az
```

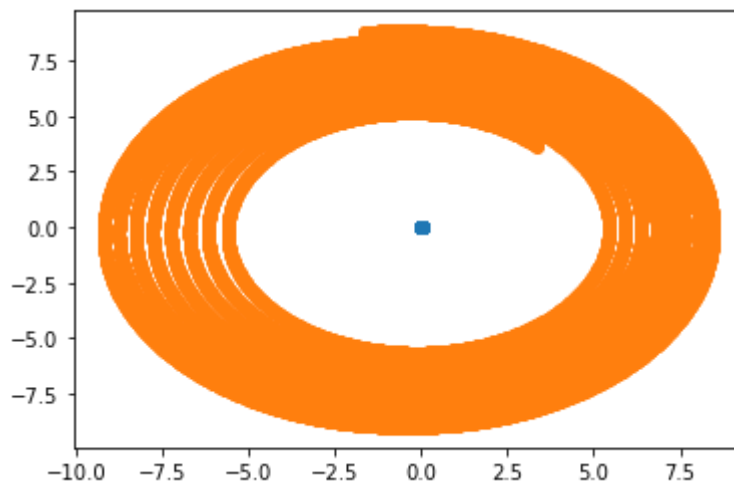
```
In [5]: sim.dt = 1e-1
Nsteps = int(1000/sim.dt)
xyz = np.zeros((Nsteps, sim.N,3))
for i in range(Nsteps):
    eulerstep(sim)
    for j in range(sim.N):
        xyz[i][j] = sim.particles[j].xyz
```

```
In [6]: for i in range(sim.N):
        plt.scatter(xyz[:,i,0], xyz[:,i,1])
```



```
In [7]: plt.scatter(xyz[:,0,0], xyz[:,0,1])
        plt.scatter(xyz[:,1,0], xyz[:,1,1])
        # Euler method is really bad, and you can see that below because the
        # orbit is varying too much
```

Out[7]: <matplotlib.collections.PathCollection at 0x7fe4f04f5990>

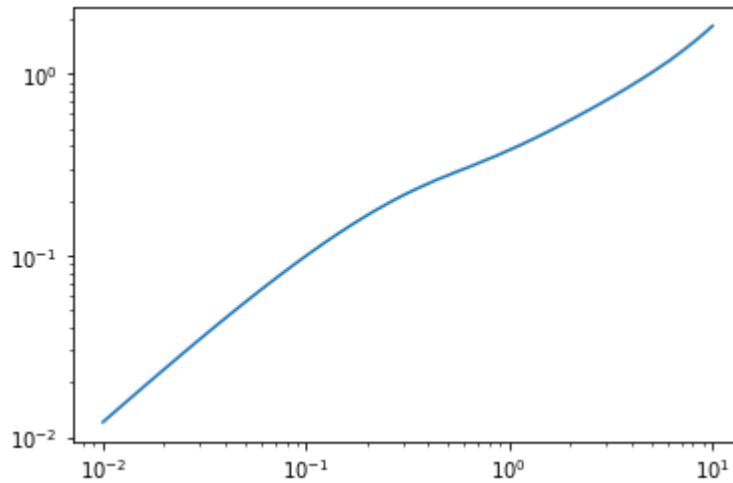


```
In [8]: def measureError(dt, step):
        sim = rebound.Simulation()
        rebound.data.add_outer_solar_system(sim)
        sim.dt = dt
        Nsteps = int(100/sim.dt)
        Ei = sim.calculate_energy()
        for i in range(Nsteps):
            step(sim)
        Ef = sim.calculate_energy()
        return np.abs((Ef-Ei)/Ei)
```

```
In [9]: dts = np.logspace(-2,1,100)
        errors_euler = np.zeros(len(dts))
        for i in range(len(dts)):
            errors_euler[i] = measureError(dts[i], eulerstep)
```

```
In [10]: plt.xscale("log")
plt.yscale("log")
plt.plot(dts, errors_euler)
```

```
Out[10]: [<matplotlib.lines.Line2D at 0x7fe4f0463950>]
```

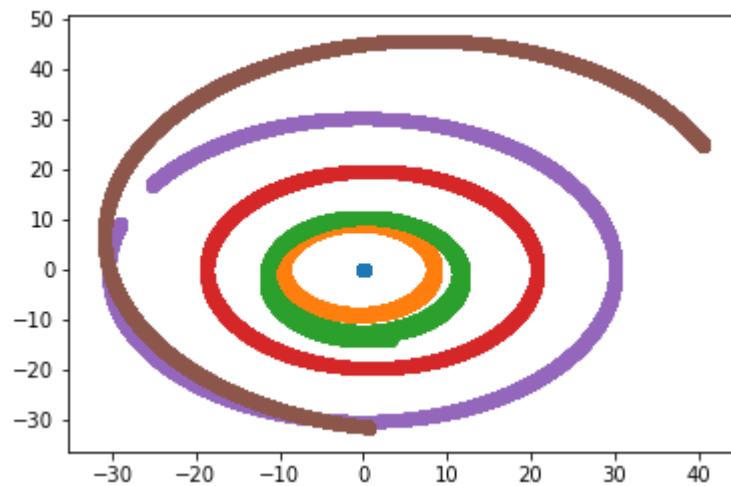


## Task 1

```
In [11]: def leapfrog(sim):
    sim.integrator = "None"
    # sim.step() # Calculates the accelerations
    for i in range(sim.N):
        sim.particles[i].x += 0.5*sim.dt * sim.particles[i].vx
        sim.particles[i].y += 0.5*sim.dt * sim.particles[i].vy
        sim.particles[i].z += 0.5*sim.dt * sim.particles[i].vz
    sim.step()
    for i in range(sim.N):
        sim.particles[i].vx += sim.dt * sim.particles[i].ax
        sim.particles[i].vy += sim.dt * sim.particles[i].ay
        sim.particles[i].vz += sim.dt * sim.particles[i].az
    for i in range(sim.N):
        sim.particles[i].x += 0.5*sim.dt * sim.particles[i].vx
        sim.particles[i].y += 0.5*sim.dt * sim.particles[i].vy
        sim.particles[i].z += 0.5*sim.dt * sim.particles[i].vz
```

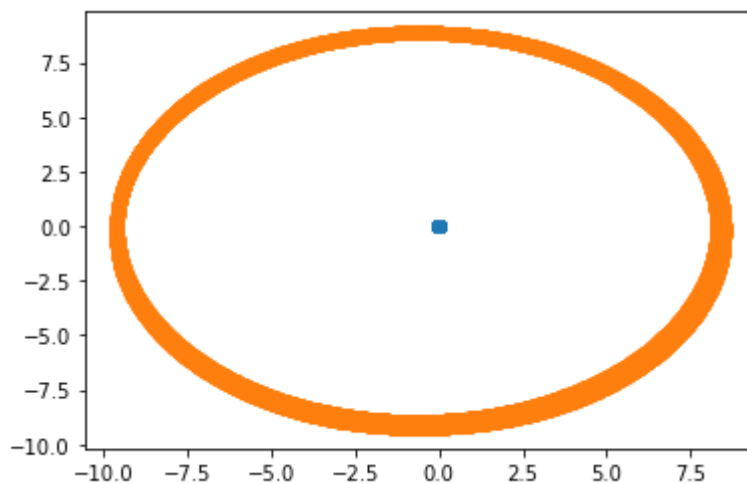
```
In [12]: # Checking out leapfrog
sim.dt = 1e-1
Nsteps = int(1000/sim.dt)
xyz = np.zeros((Nsteps, sim.N,3))
for i in range(Nsteps):
    leapfrog(sim)
    for j in range(sim.N):
        xyz[i][j] = sim.particles[j].xyz
```

```
In [13]: for i in range(sim.N):  
         plt.scatter(xyz[:,i,0], xyz[:,i,1])
```



```
In [14]: # Checking out the Jupiter orbit  
plt.scatter(xyz[:,0,0], xyz[:,0,1])  
plt.scatter(xyz[:,1,0], xyz[:,1,1])
```

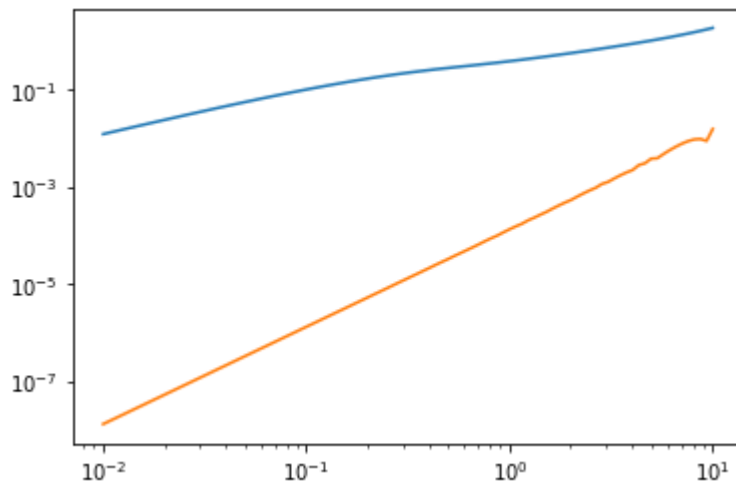
Out[14]: <matplotlib.collections.PathCollection at 0x7fe4d91281d0>



```
In [15]: dts = np.logspace(-2,1,100)  
errors_leapfrog = np.zeros(len(dts))  
for i in range(len(dts)):  
    errors_leapfrog[i] = measureError(dts[i], leapfrog)
```

```
In [16]: plt.xscale("log")
plt.yscale("log")
plt.plot(dts, errors_euler)
plt.plot(dts, errors_leapfrog)
```

```
Out[16]: [<matplotlib.lines.Line2D at 0x7fe4d93c1b50>]
```

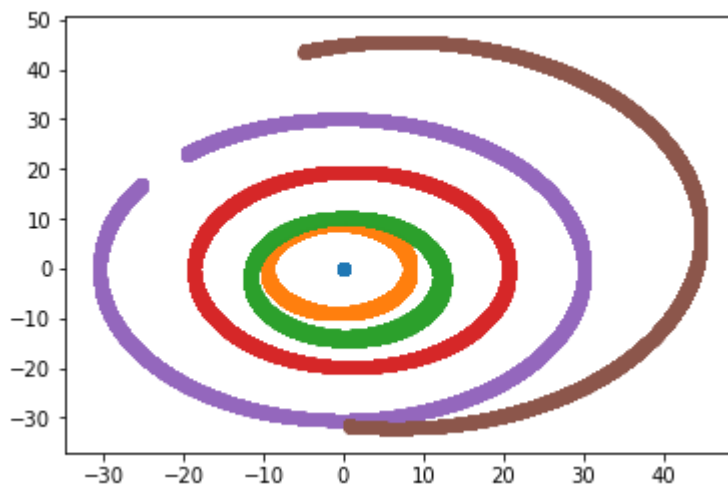


## Task 2

```
In [17]: def whfaststep(sim):
sim.integrator = "WHfast"
sim.step()
```

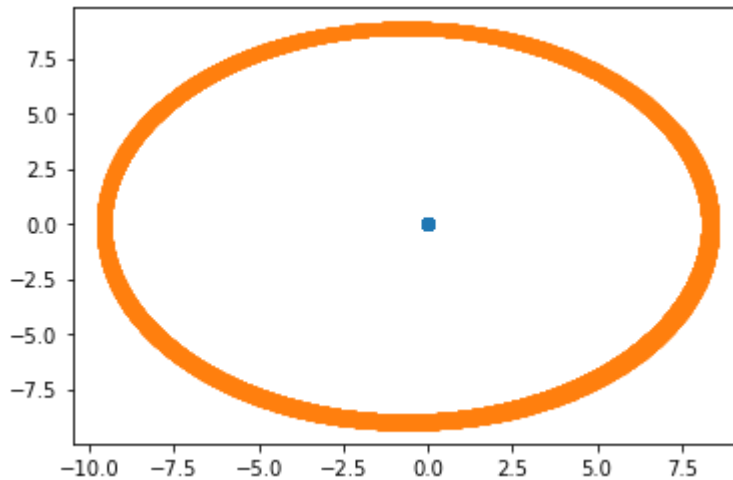
```
In [18]: # Checking out WHfast
sim.dt = 1e-1
Nsteps = int(1000/sim.dt)
xyz = np.zeros((Nsteps, sim.N, 3))
for i in range(Nsteps):
    whfaststep(sim)
    for j in range(sim.N):
        xyz[i][j] = sim.particles[j].xyz
```

```
In [19]: for i in range(sim.N):
plt.scatter(xyz[:,i,0], xyz[:,i,1])
```



```
In [20]: # Checking out the Jupiter orbit
plt.scatter(xyz[:,0,0], xyz[:,0,1])
plt.scatter(xyz[:,1,0], xyz[:,1,1])
```

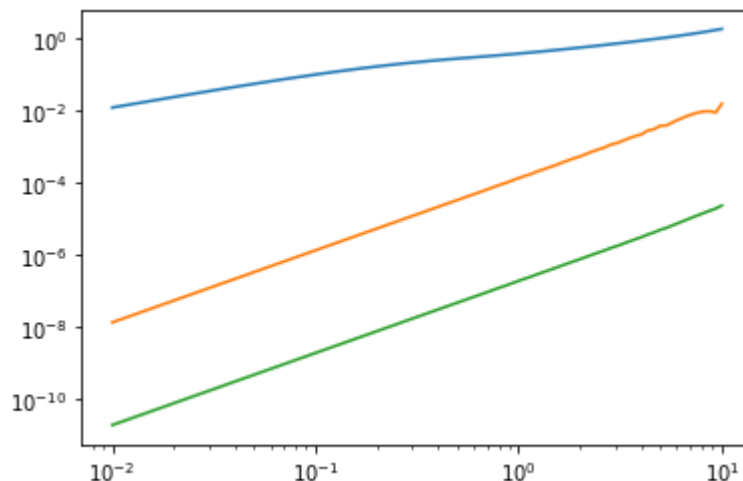
Out[20]: <matplotlib.collections.PathCollection at 0x7fe4f0489710>



```
In [21]: dts = np.logspace(-2,1,100)
errors_whfaststep = np.zeros(len(dts))
for i in range(len(dts)):
    errors_whfaststep[i] = measureError(dts[i], whfaststep)
```

```
In [22]: plt.xscale("log")
plt.yscale("log")
plt.plot(dts, errors_euler)
plt.plot(dts, errors_leapfrog)
plt.plot(dts, errors_whfaststep)
```

Out[22]: [<matplotlib.lines.Line2D at 0x7fe4d8e9ea90>]



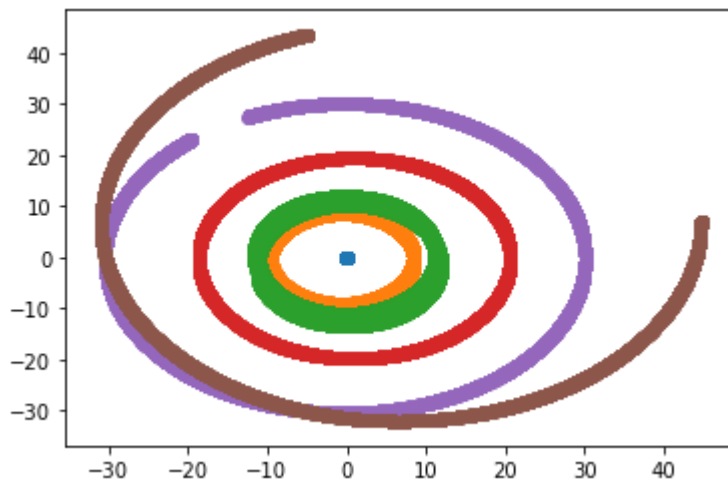
The error in the WHFaststep is significantly lower, and it is of the same order as the leapfrog method, because the slopes are the same.

## Bonus Task 1

```
In [23]: def whckl(sim):
          sim.integrator = "WHCKL"
          sim.step()
```

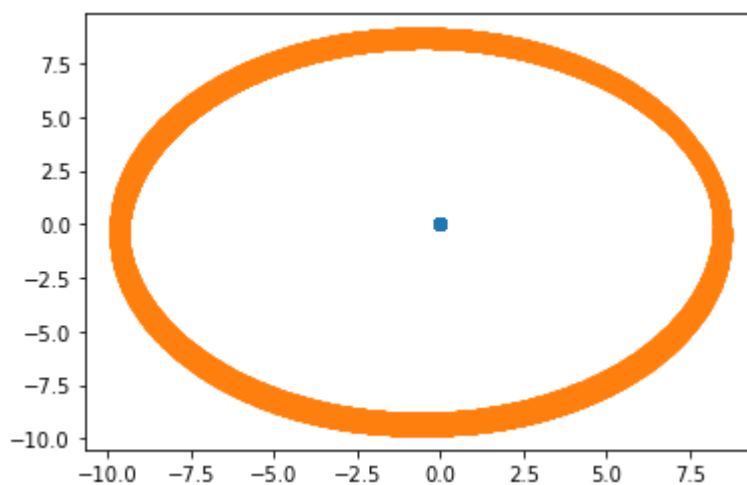
```
In [24]: # Checking out WHCKL
          sim.dt = 1e-1
          Nsteps = int(1000/sim.dt)
          xyz = np.zeros((Nsteps, sim.N,3))
          for i in range(Nsteps):
              whckl(sim)
              for j in range(sim.N):
                  xyz[i][j] = sim.particles[j].xyz
```

```
In [25]: for i in range(sim.N):
          plt.scatter(xyz[:,i,0], xyz[:,i,1])
```



```
In [26]: # Checking out the Jupiter orbit
          plt.scatter(xyz[:,0,0], xyz[:,0,1])
          plt.scatter(xyz[:,1,0], xyz[:,1,1])
```

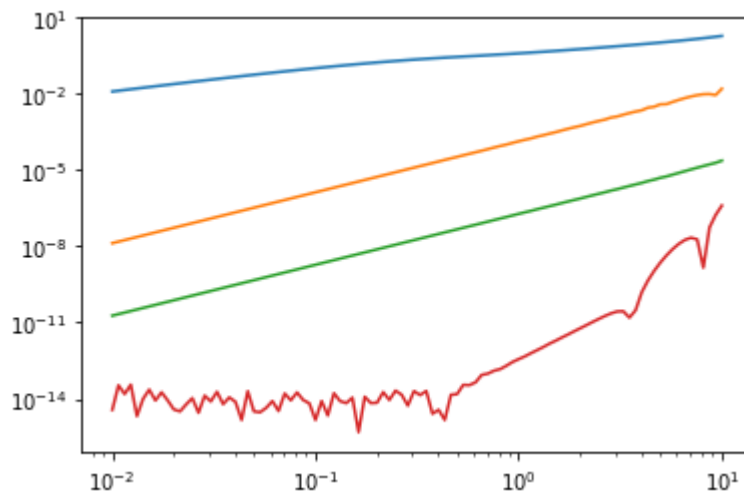
Out[26]: <matplotlib.collections.PathCollection at 0x7fe4d91c41d0>



```
In [27]: dts = np.logspace(-2,1,100)
          errors_whckl = np.zeros(len(dts))
          for i in range(len(dts)):
              errors_whckl[i] = measureError(dts[i], whckl)
```

```
In [28]: plt.xscale("log")
plt.yscale("log")
plt.plot(dts, errors_euler)
plt.plot(dts, errors_leapfrog)
plt.plot(dts, errors_whfaststep)
plt.plot(dts, errors_whckl)
```

Out[28]: [<matplotlib.lines.Line2D at 0x7fe4d8da62d0>]



The WHCKL method has a different order as the errors increase significantly as the size of the timestep is increased.

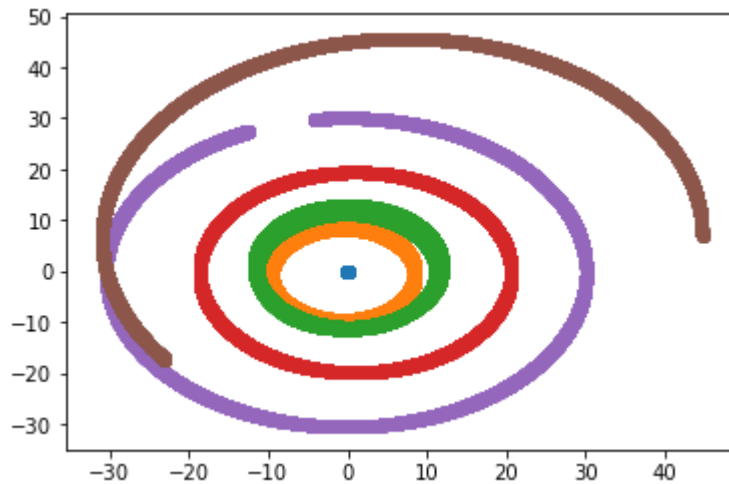
## Bonus Task 2

```
In [29]: def whfaststep_c(sim):
sim.integrator = "WHfast"
sim.ri_whfast.corrector = 17
sim.step()
```

```
In [30]: # Checking out WH with symplectic correctors
sim.dt = 1e-1
Nsteps = int(1000/sim.dt)
xyz = np.zeros((Nsteps, sim.N, 3))
for i in range(Nsteps):
    whfaststep_c(sim)
    for j in range(sim.N):
        xyz[i][j] = sim.particles[j].xyz
```

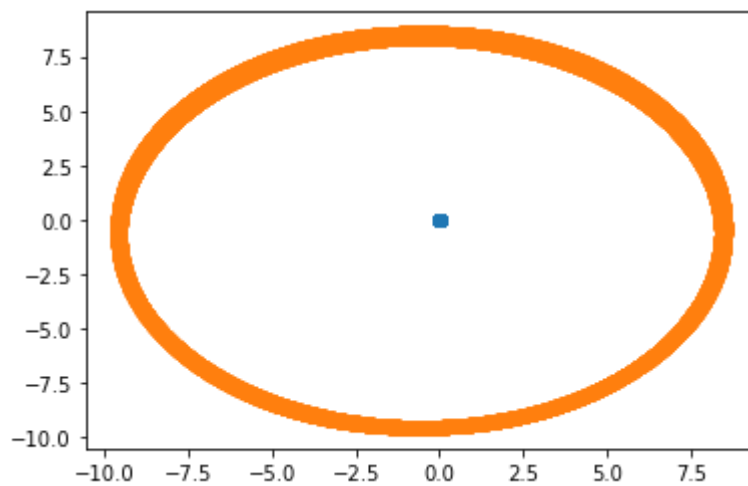


```
In [31]: for i in range(sim.N):  
         plt.scatter(xyz[:,i,0], xyz[:,i,1])
```



```
In [32]: # Checking out the Jupiter orbit  
plt.scatter(xyz[:,0,0], xyz[:,0,1])  
plt.scatter(xyz[:,1,0], xyz[:,1,1])
```

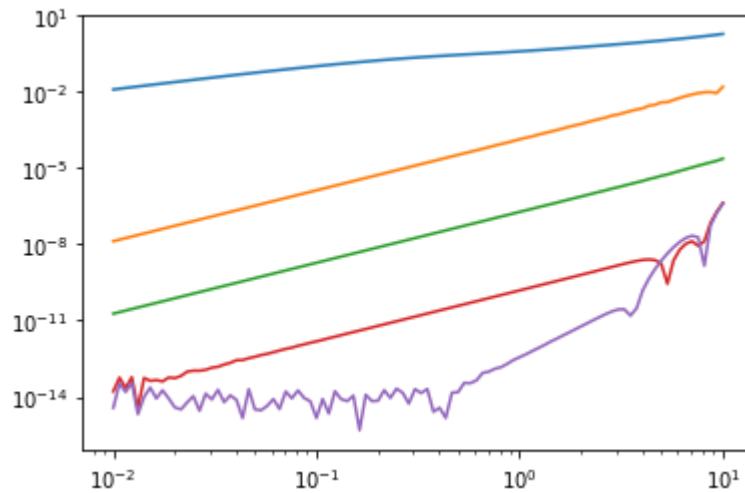
Out[32]: <matplotlib.collections.PathCollection at 0x7fe4d93c8fd0>



```
In [33]: dts = np.logspace(-2,1,100)  
errors_whfaststep_c = np.zeros(len(dts))  
for i in range(len(dts)):  
    errors_whfaststep_c[i] = measureError(dts[i], whfaststep_c)
```

```
In [34]: plt.xscale("log")
plt.yscale("log")
plt.plot(dts, errors_euler)
plt.plot(dts, errors_leapfrog)
plt.plot(dts, errors_whfaststep)
plt.plot(dts, errors_whfaststep_c)
plt.plot(dts, errors_whckl)
```

Out[34]: [<matplotlib.lines.Line2D at 0x7fe4d8c533d0>]



In [ ]: