

N-Body Gravitationally Interacting System!



Background!

- We approximate a gravitationally attractive multibody system by integrating the equation of motion
- Three different scenarios are considered:
 - Solar: Central massive body, smaller planets orbiting around
 - Lunar: Moon orbiting a planet as planet orbits a star
 - Random: Particles with a random range of masses in random positions, given velocity according to a Gaussian distribution
- Each scenario is ran for a range of step sizes, after which initial/final momentum & energy are compared

```
class Particle(object):
```

```
    def __init__(self, mass, x, y, z, vx, vy, vz):  
        self.mass = mass  
        position = np.array([x,y,z])  
        velocity = np.array([vx,vy,vz])
```

```
  
        self.position = position  
        self.velocity = velocity
```

```
  
        self.position_list = [position]  
        self.velocity_list = [velocity]
```

```
  
    def combine_particles(self, particle2, particle_list, crit_radius):
```

```
  
        distance = self.get_distance(particle2)
```

```
        r = (distance[0] ** 2.0 + distance[1] ** 2.0 + distance[2] ** 2.0) ** (1.0 / 2.0)
```

```
  
        #print (distance)
```

```
        #print (r)
```

```
  
        if r < crit_radius:
```

```
            print("Combining Particles")
```

```
            # print(particle_list)
```

```
  
            M = self.get_mass() + particle2.get_mass()
```

```
            new_v = (self.get_mass() * self.get_velocity_vector() + particle2.get_mass() * particle2.get_velocity_vector()) / M
```

```
            new_p = (self.get_position_vector() + particle2.get_position_vector()) / 2.0
```

```
  
            new_particle = Particle(M, new_p[0], new_p[1], new_p[2], new_v[0], new_v[1], new_v[2])
```

```
            particle_list.append(new_particle)
```

```
  
            particle_list.remove(self)
```

```
            particle_list.remove(particle2)
```

Math!

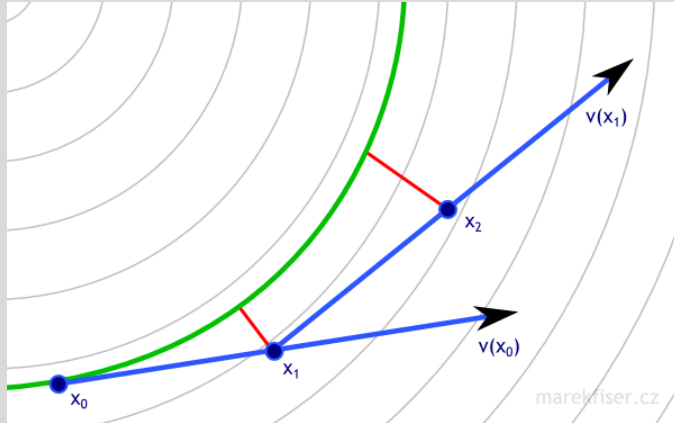
$$U = -G \frac{Mm}{r}$$

$$\vec{a} = \frac{\vec{F}}{m} = -G \frac{M}{r^2} \hat{r}$$

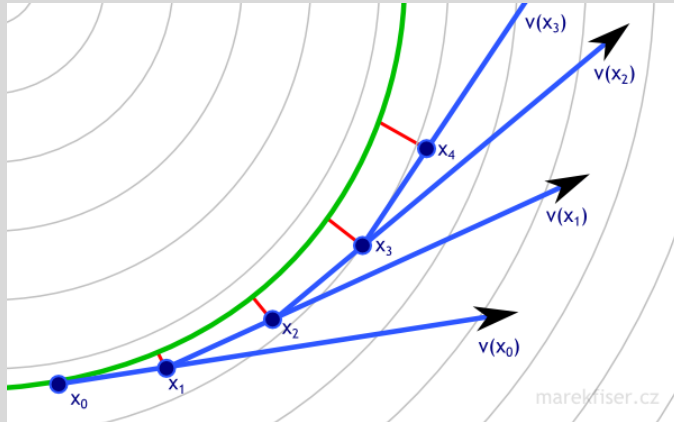
$$\vec{F} = -\nabla U = -G \frac{Mm}{r^2} \hat{r}$$

- Each vector is coded as a numpy array
- Movement is due to the net acceleration felt by all other particles
- Movement determined by integrator (RK4)

Integrator: Runge-Kutta 4th Order!



$dt = 0.5$



$dt = 0.25$

- Euler integration two different timesteps
- Vector field shown as grey lines
- Green curve represents actual path
- Blue is the Euler Method (RK1)
- Red lines shows error

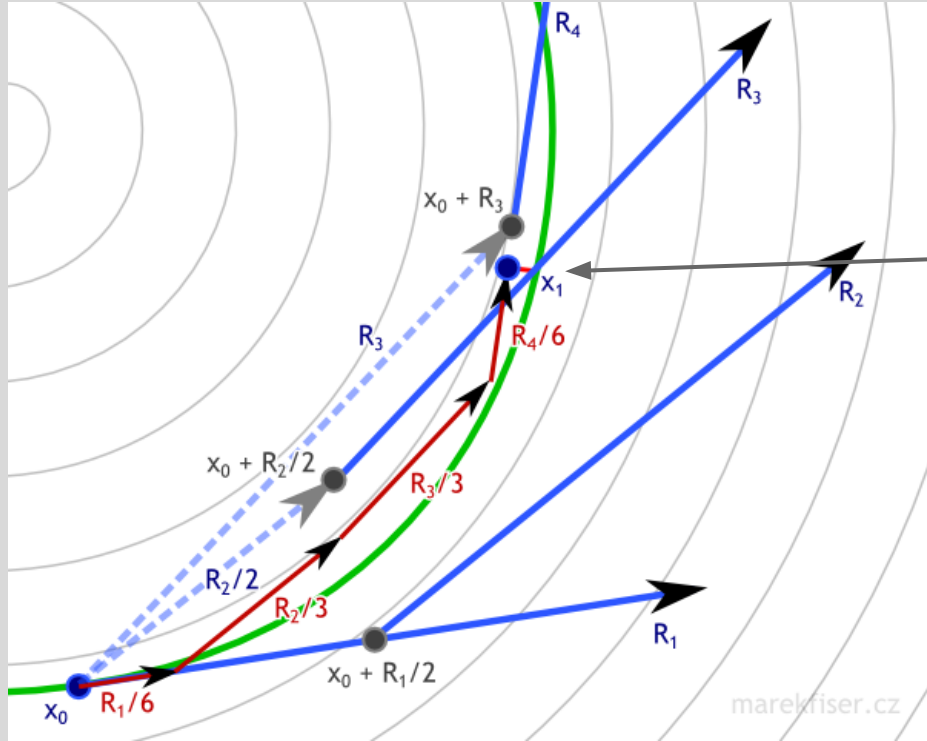
Integrator: Runge-Kutta 4th Order!

$\vec{k}_{1r_{i+1}} = \vec{v}_i$		$\vec{k}_{1v_{i+1}} = \vec{a}(\vec{r}_i)$
$\vec{k}_{2r_{i+1}} = \vec{v}_i + \vec{k}_{1v_{i+1}} \frac{h}{2}$		$\vec{k}_{2v_{i+1}} = \vec{a}(\vec{r}_i + \vec{k}_{1r_{i+1}} \frac{h}{2})$
$\vec{k}_{3r_{i+1}} = \vec{v}_i + \vec{k}_{2v_{i+1}} \frac{h}{2}$		$\vec{k}_{3v_{i+1}} = \vec{a}(\vec{r}_i + \vec{k}_{2r_{i+1}} \frac{h}{2})$
$\vec{k}_{4r_{i+1}} = \vec{v}_i + \vec{k}_{3v_{i+1}} h$		$\vec{k}_{4v_{i+1}} = \vec{a}(\vec{r}_i + \vec{k}_{3r_{i+1}} h)$

$$\vec{r}_{i+1} = \vec{r}_i + \frac{h}{6}(\vec{k}_{1r_{i+1}} + 2\vec{k}_{2r_{i+1}} + 2\vec{k}_{3r_{i+1}} + \vec{k}_{4r_{i+1}})$$

$$\vec{v}_{i+1} = \vec{v}_i + \frac{h}{6}(\vec{k}_{1v_{i+1}} + 2\vec{k}_{2v_{i+1}} + 2\vec{k}_{3v_{i+1}} + \vec{k}_{4v_{i+1}})$$

Integrator: Runge-Kutta 4th Order!



$$\vec{r}_{i+1} = \vec{r}_i + \frac{h}{6}(\vec{k}_{1r_{i+1}} + 2\vec{k}_{2r_{i+1}} + 2\vec{k}_{3r_{i+1}} + \vec{k}_{4r_{i+1}})$$
$$\vec{v}_{i+1} = \vec{v}_i + \frac{h}{6}(\vec{k}_{1v_{i+1}} + 2\vec{k}_{2v_{i+1}} + 2\vec{k}_{3v_{i+1}} + \vec{k}_{4v_{i+1}})$$

Error in measurement: little red line

These coefficients are the slope of the function at three separate points during the timestep: the beginning, the mid-point and the end.

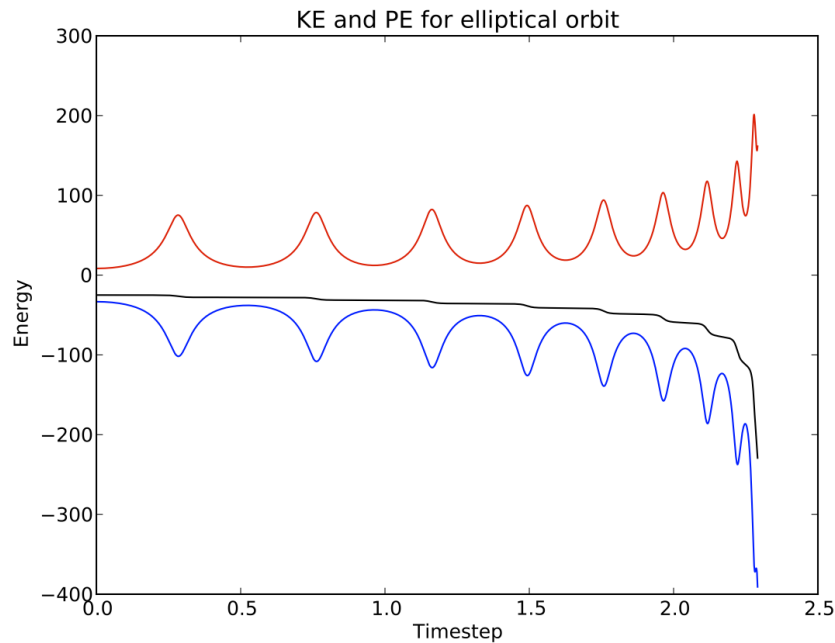
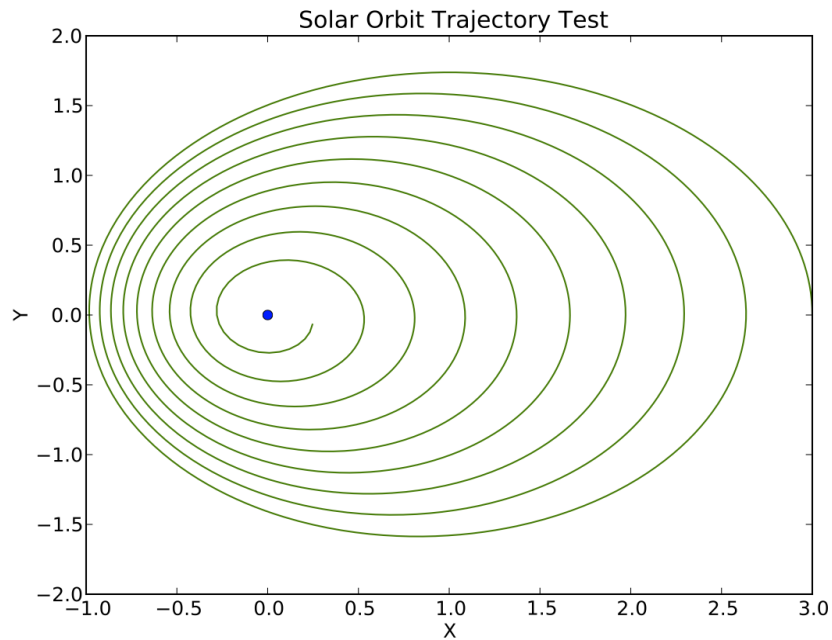
RK4 (dt = 1)

Acceleration Calculation: The Work Horse!

```
def acceleration(self, position1):  
    """Accepts particle's position; returns net acceleration"""  
    ax, ay, az = 0, 0, 0  
    for p2 in particles:  
        if p2 != self:  
            r = distance(self, p2)  
            a = G * p2.mass / r**2.0  
            a_x = a * (p2.position[0] - position1[0]) / r  
            a_y = a * (p2.position[1] - position1[1]) / r  
            a_z = a * (p2.position[2] - position1[2]) / r  
            ax += a_x  
            ay += a_y  
            az += a_z  
    return (ax, ay, az)
```


Effects of Stepsize!

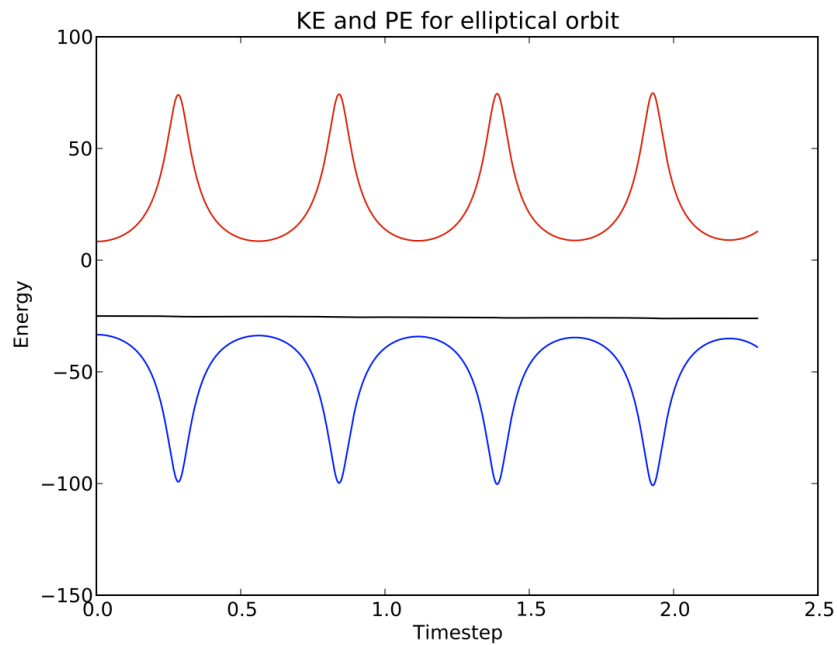
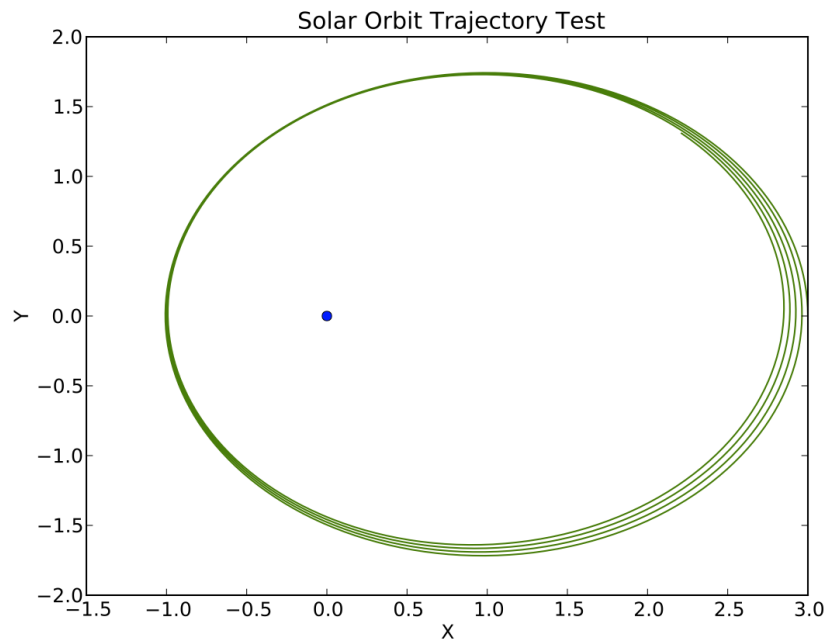
A Small Timestep is the Enemy of Energy Conservation



Solar Orbit Examples

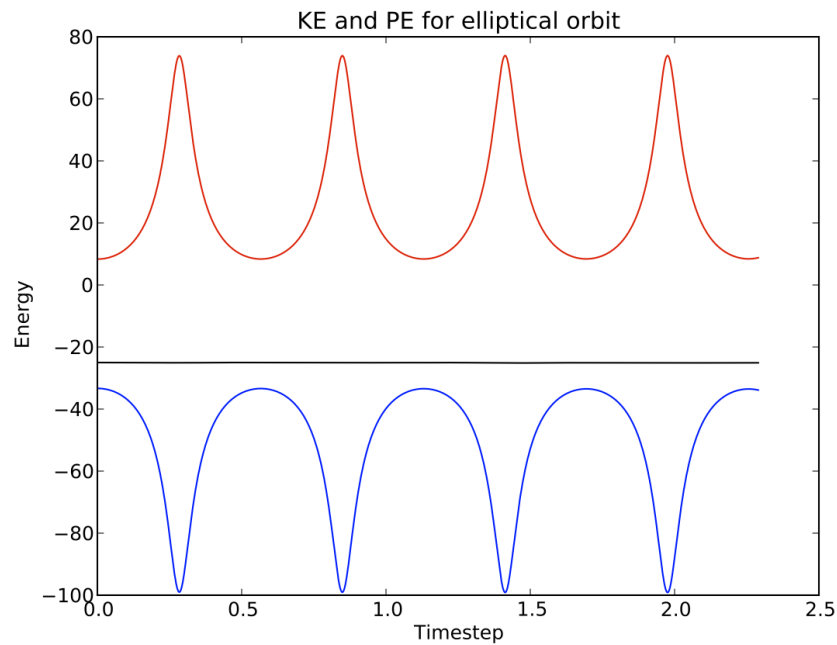
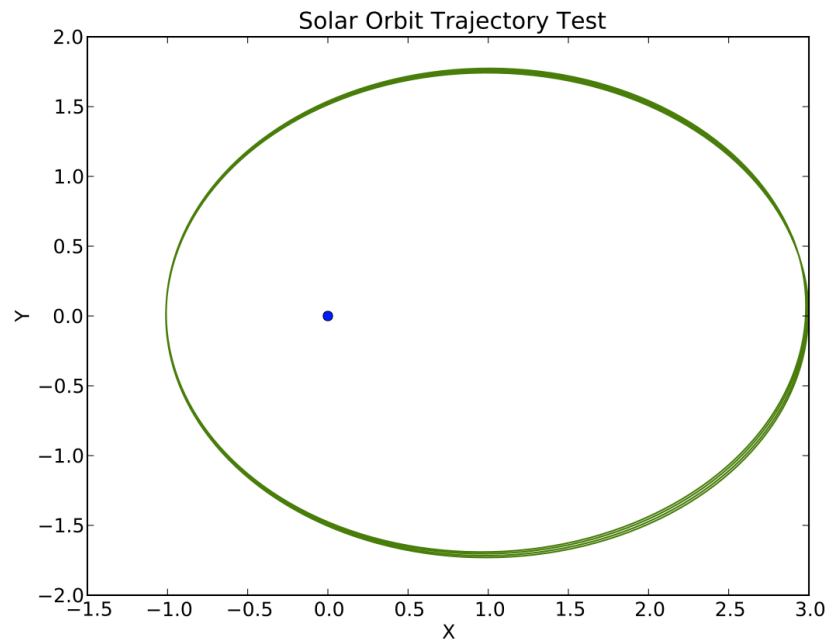
$h = 0.001$

Effects of Stepsize!



$h = 0.0001$

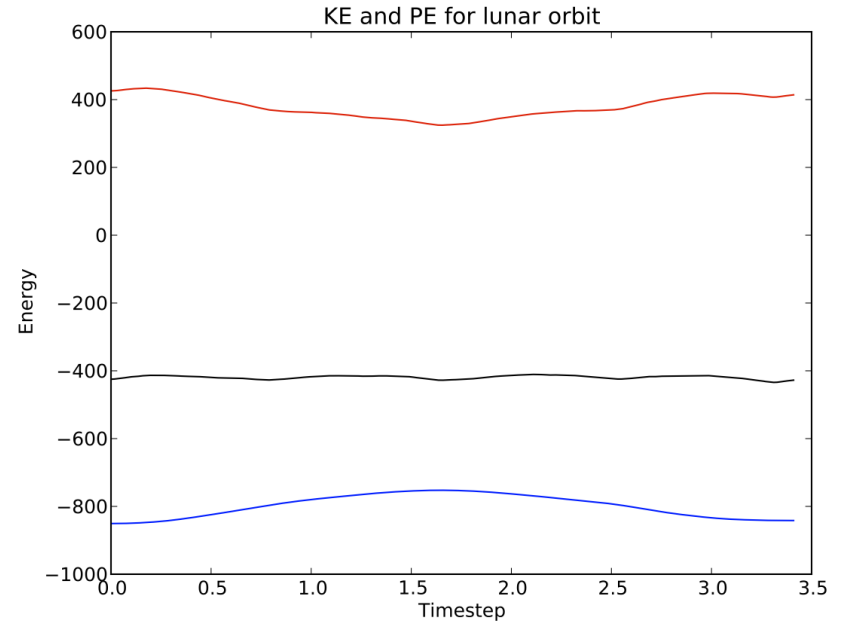
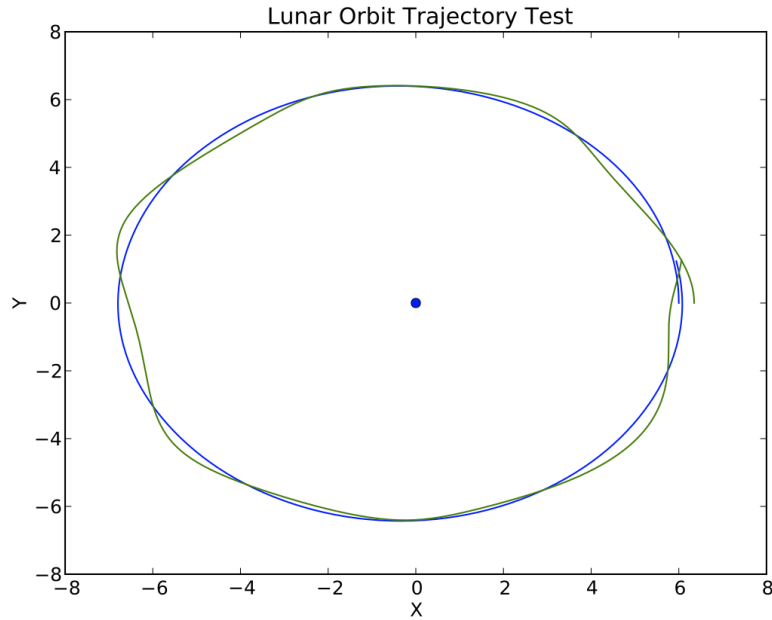
Effects of Stepsize!



$h = 0.00001$

Lunar Orbit Example

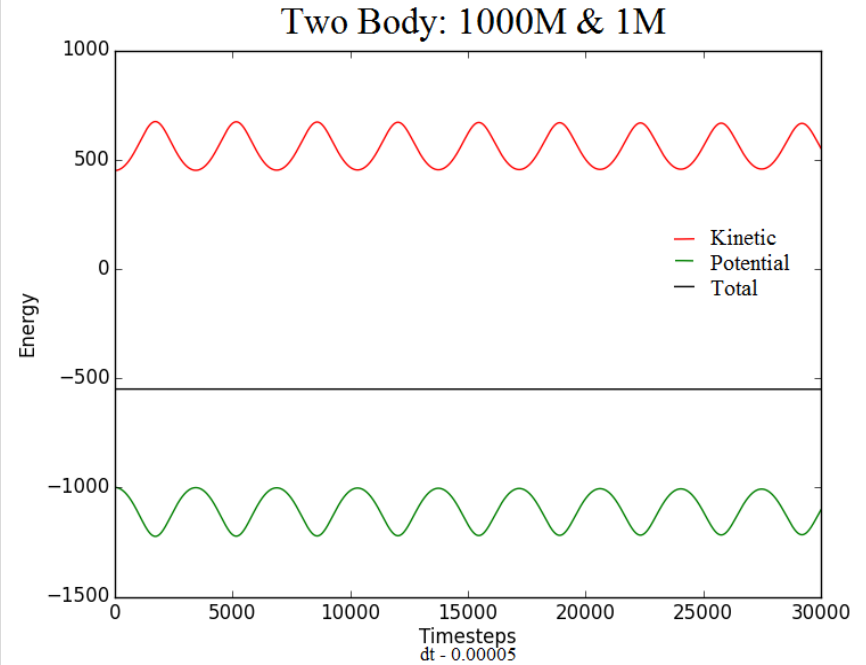
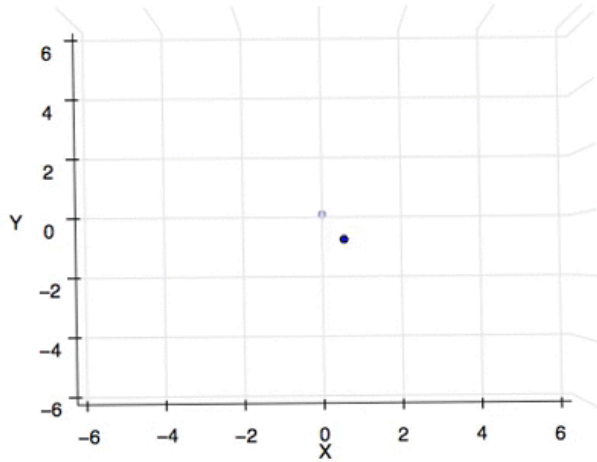
$h = 0.00001$



The lunar orbit highlights the problem with close interactions when dealing with discrete timesteps.

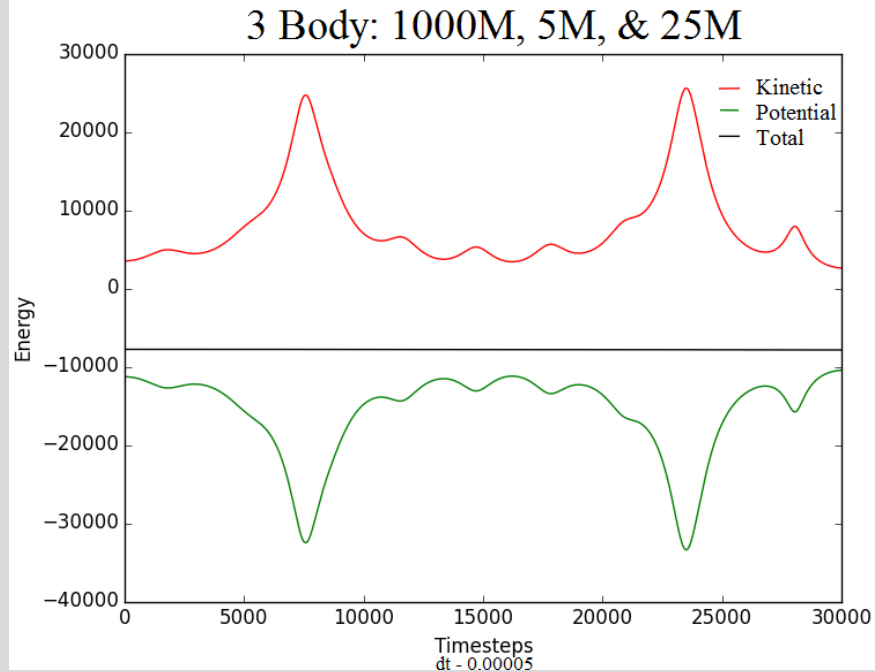
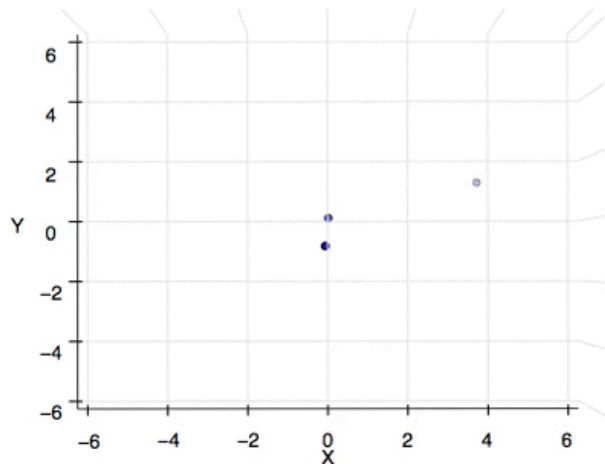
Animation!

Two Body System



Animation!

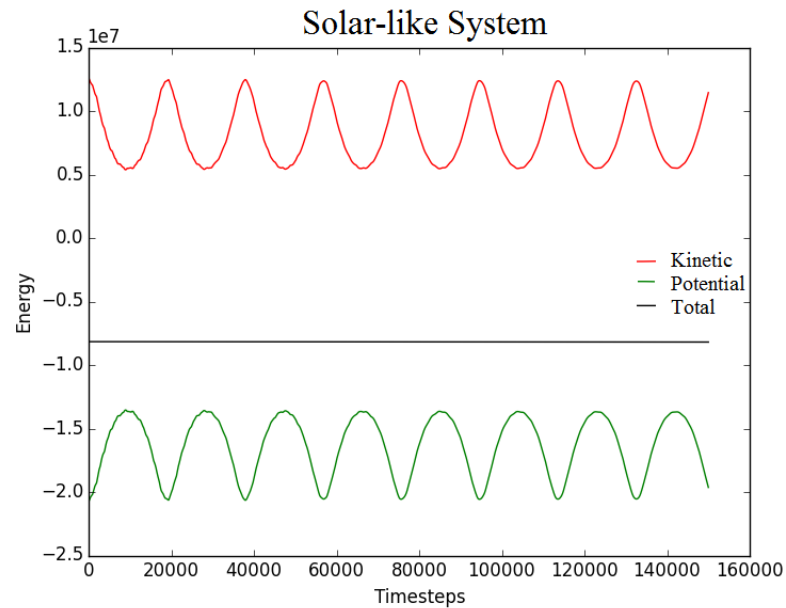
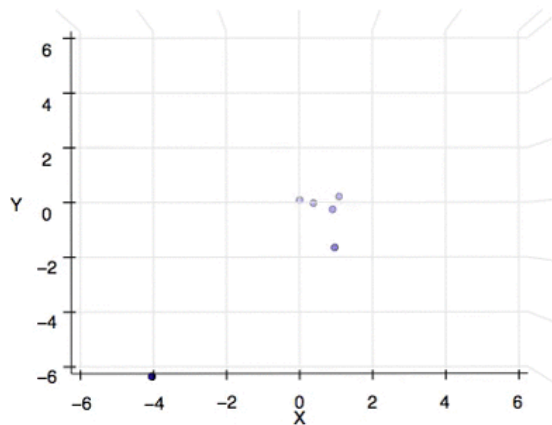
Three Body System

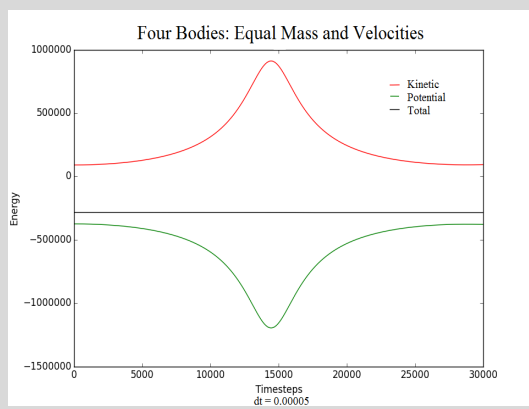


Animation!

Solar-like System (Mercury to Jupiter)

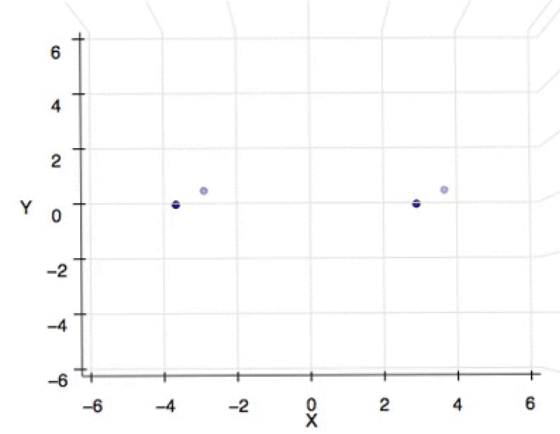
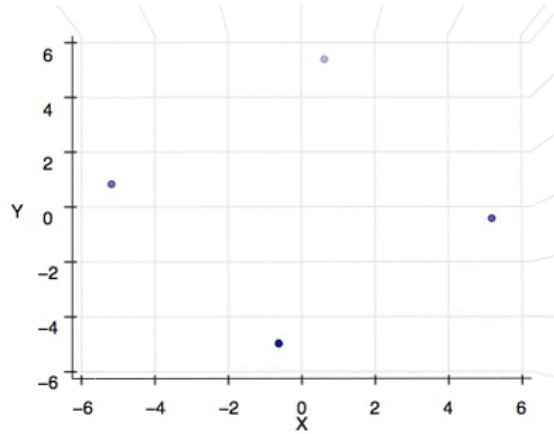
- Similar masses to inner planets plus Jupiter
- Jupiter dominates





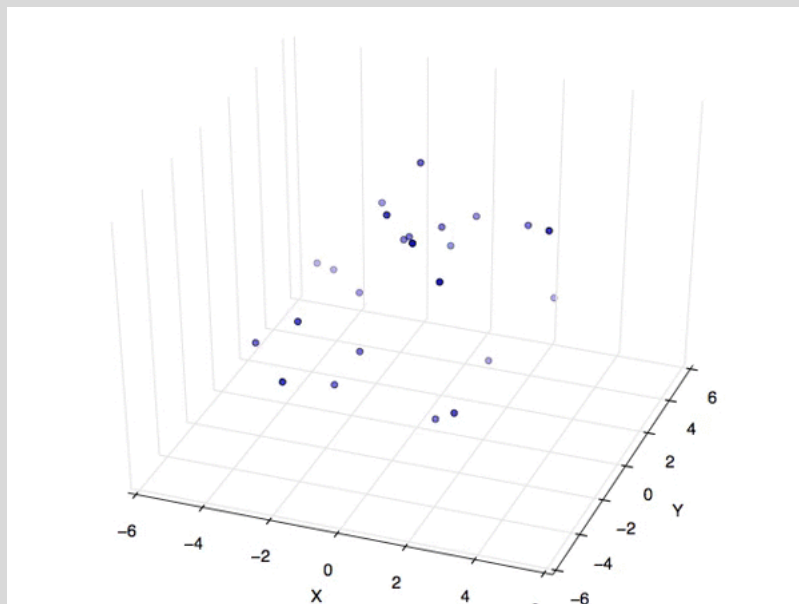
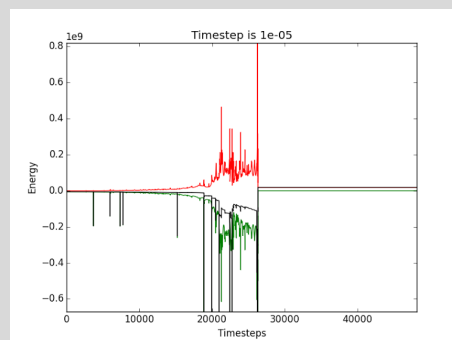
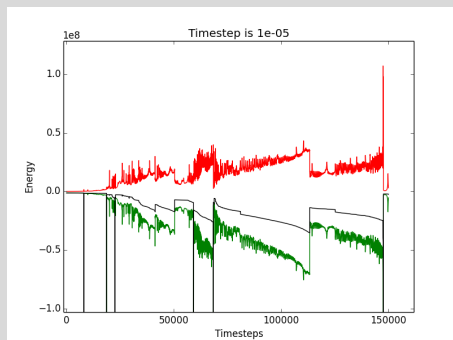
Animation!

Four Equal Mass System

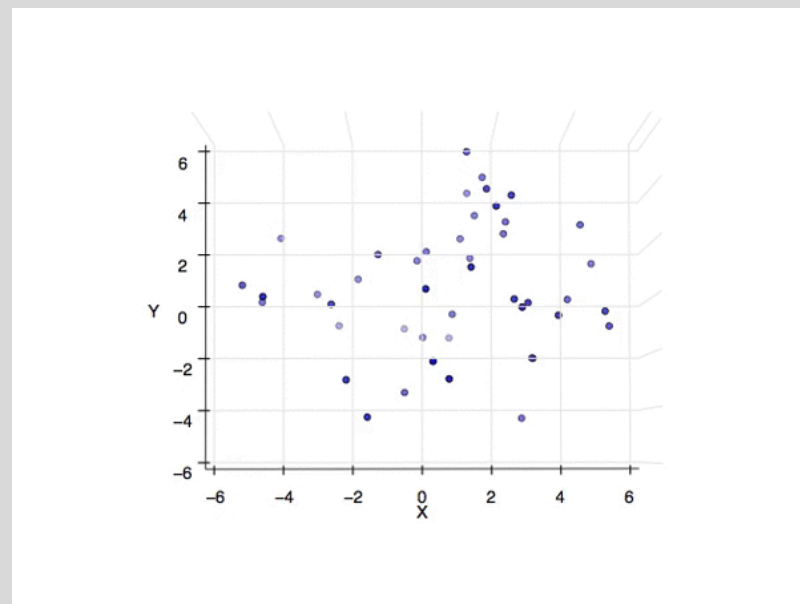


Animation!

N-body Collisional System



$N = 25$

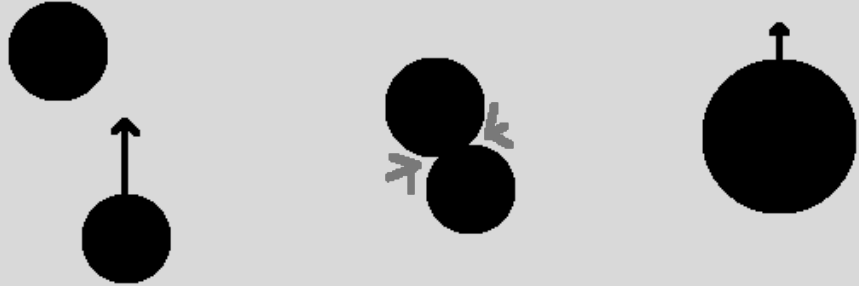


$N = 50$

Energy Conservation!

Sources of Error

- Every collision produces some amount of error (inelastic collision)
- High velocities demand high timestep
 - Dynamical timestep



→ `h = h_min * (top_velocity/current_max)`

Room for Improvement!

- Import FORTRAN for workhorse calculations
- Don't double record acceleration
- Initialize particles with a spin property that adds during collisions