Investigation of the motion of a double pendulum

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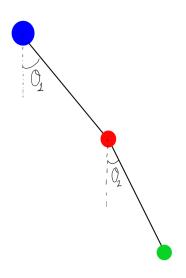
A computational method was used to obtain the trajectories of two masses of a double pendulum, which is a dynamical system with chaotic behavior. Different initial parameters, mostly different angles, were used. Program calculated 400 frames per simulation of double pendulum with given specific initial conditions. Chaotic motion had been expected to be observed with large initial angles. And on the contrary, linear alike motion had been expected to be observed with small initial angles. Chaotic motion was observed, but linear motion was not as significant as expected due to some numerical errors.

PACS numbers:

INTRODUCTION

Chaotic systems are of dynamical systems that are highly sensitive to initial conditions. That means a small change in initial parameters of the system will result a great difference in final status of the system. Since the final status of the system can differ from others within a small change of initial parameter, it is a common misconception that it is impossible to predict the behavior of the system. However some chaotic systems, like double pendulum, can be deterministic. This means that if its initial conditions are known, and can integrate it forward in time with infinite precision, we could predict its motion. For further information about chaotic motion and its applications in real life, referenced as [5] website can be visited. It shows papers on the applications of chaotic motion which can be helpful for further applications. Also Akerlof's experiment on double pendulum [7] gives rigorous explanations for those who want advanced explanations.

THEORY



In the diagram blue circle is fixed. Red is attached to blue and green is attached to red. Both red and green circles can move. Now call the position of red circle as (x_r, y_r) , position of green circle as (x_g, y_g) , distance between blue and red circle as l_1 and distance between red and green as l_2 . Their position can also be written as;

$$x_r = l_1 sin(\theta_1) \tag{1}$$

$$y_r = -l_1 cos(\theta_1) \tag{2}$$

$$x_q = l_1 sin(\theta_1) + l_2 sin(\theta_2) \tag{3}$$

$$y_q = -l_1 cos(\theta_1) - l_2 cos(\theta_2) \tag{4}$$

To find the instantaneous velocities we take the derivatives of the equations 1-4 with respect to time.

$$\frac{dx_r}{dt} = l_1 cos(\theta_1) \frac{d\theta_1}{dt} \tag{5}$$

$$\frac{dy_r}{dt} = l_1 sin(\theta_1) \frac{d\theta_1}{dt} \tag{6}$$

$$\frac{dx_g}{dt} = l_1 cos(\theta_1) \frac{d\theta_1}{dt} + l_2 cos(\theta_2) \frac{d\theta_2}{dt}$$
 (7)

$$\frac{dy_g}{dt} = l_1 sin(\theta_1) \frac{d\theta_1}{dt} + l_2 sin(\theta_2) \frac{d\theta_2}{dt}$$
 (8)

To solve the Lagrange's Equation, it is necessary to find Lagrangian quantity which is the difference of kinetic and potential energies of the system.

$$\mathcal{L} = E_K - E_P = T - V \tag{9}$$

$$V = m_r g y_r + m_g g y_g$$

 $m_r = \text{mass of the red circle}, m_g = \text{mass of the green}$

We can insert equation 2 and 4 to Potential equation:

$$V = -(m_r + m_g)gl_1cos(\theta_1) - m_ggl_2cos(\theta_2)$$
 (10)

Kinetic energy(T):

$$T = \frac{m_r V_1^2 + m_g V_2^2}{2}$$

$$T = \frac{m_r((\frac{dx_r}{dt})^2 + (\frac{dy_r}{dt})^2) + m_g((\frac{dx_g}{dt})^2 + (\frac{dy_g}{dt})^2)}{2}$$

Substitute equations 5-8 into the kinetic energy equation,

$$T = \frac{m_r[(l_1 cos(\theta_1) \frac{d\theta_1}{dt})^2 + (l_1 sin(\theta_1) \frac{d\theta_1}{dt})^2]}{2}$$

$$+$$

$$\frac{m_g[(l_1 cos(\theta_1) \frac{d\theta_1}{dt} + l_2 cos(\theta_2) \frac{d\theta_2}{dt})^2]}{2} +$$

$$\frac{(l_1 sin(\theta_1) \frac{d\theta_1}{dt} + l_2 sin(\theta_2) \frac{d\theta_2}{dt})^2}{2}$$

Then we can substitute equation 10 and 11 into equation 9.

For the last step we can apply Lagrangian's Equations for θ_1 and θ_2 to find angular speeds and angular accelerations

$$\frac{d}{dt} \left(\frac{\partial \mathcal{L}}{\partial \left(\frac{d\theta}{dt} \right)} \right) - \frac{\partial \mathcal{L}}{\partial \theta} = 0$$

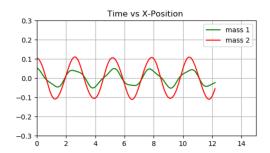
METHOD

Python program will perform continuous calculations to find the positions of the masses. To find the the angular accelaration and angular speeds of the masses, lagrangian of the system will be solved. And the equations that will be obtained from the lagrangian equation will be used to find the angles for given times. Python's built in differential equation solver will be used. User will give initial conditions, such as initial angles, masses of the objects and lengths of the rods. However masses of the rods will be neglected. Program then simulate the motion of the pendulum, plot the graphes of total energy of the system, x and y positions ys time.

THE COMPUTATIONAL RESULTS & INTERPRETATION

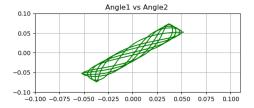
Initially, masses of the objects were set to 1 kg each and the lenths of the rods were set to 1 meters. The gravitational acceleration was set to $9.8m/s^2$. Then, for the first simulation, small angles were assigned as parameters. $\theta_1 = 3^{\circ}$ and $\theta_2 = 3^{\circ}$.

Double pendulum started to oscillate with small angles. Since the angle was small system behaved like a linear double spring. (Since double spring is out of the scope of this experiment, no detail was given. However double spring can be observed in the website with reference number [4]). The x and y positions of the masses were,





From the plot of time vs x-position, mass 2 oscilated like a simple pendulum. However, first mass oscilated strangely. The differential equation solver module could be the cause of the strange oscillations . Or it could be normal. When θ_1 vs θ_2 graph plotted,

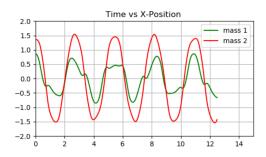


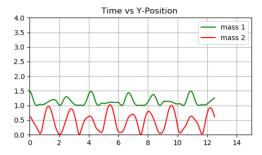
Both angles are in radians

it traced out the familiar form of a Lissajous curve. This small motion is structered well and does not look like a chaotic system. This is due to the small angle approximation. When θ_1 and θ_2 are small enough their sin and cos functions become θ and 0 respectively. Also, since their

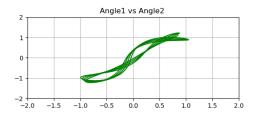
momenta are small, the product of their momenta will be very small. Thus their momenta and product of momentas can be ignored. When all these considered, nonlinear systems of equations become linear. And motion of the pendulum becomes predictable.

When we enlarged the initial angles to $\theta_1 = 50^{\circ}$ and $\theta_2 = 70^{\circ}$, the motion of the masses were





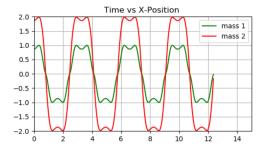
From the plot of time vs x-position, mass 2 oscilated like a simple pendulum. However, first mass oscilated strangely even more than the the oscilation of the small angle. In here, small angle approximation can still be used. However, the more the angles grow the more the error becomes. Also its Lissajous curve were plotted as,

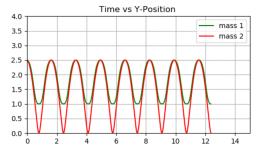


Both angles are in radians

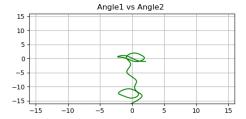
Its Lissajous curve is not as rectengular as the one with small inital angles. Its curve is stretched and bent. The curve still looks structered however it is a complex nonlinear shape.

When the initial angles were enlarged even more to $\theta_1 = 120^{\circ}$ and $\theta_2 = 90^{\circ}$, the motion of the masses were become as,





Both masses oscilated strangely. This can be seen from the x positions of masses in different times. Also its Lissajous curve was plotted as

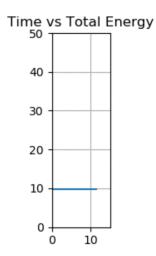


Both angles are in radians

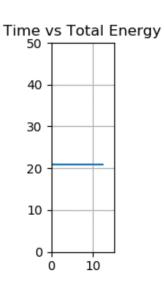
From its Lissajous curve, it can be seen that its structure is nonlinear shape.

ERRORS

For the program, pythons scipi module imported and its built in differential solver was used. This module approximated the solution better than "euler-cromer" and "leap-frog" algorithms. Although advanced error calculations were not performed, energy vs time plots were plotted. If any fluctuation in the graph were spotted, it could have been understood that the calculations have errors.



Total Energy vs. Time of the pendulum with initial angles of $\theta_1=0.5^\circ$, $\theta_2=0.5^\circ$



Total Energy vs. Time of the pendulum with initial angles of $\theta_1=60^\circ$, $\theta_2=30^\circ$

Time vs Total Energy 40 30 20 10 0 10

Total Energy vs. Time of the pendulum with initial angles of $\theta_1=60^\circ$, $\theta_2=120^\circ$

However there weren't any fluctuations spotted in any of the simulations. So it can be concluded that error was small.

APPENDIX - THE COMPUTER CODE

58 59

60

```
61
1 from numpy import sin, cos
2 import numpy as np
                                                        62
  import matplotlib.pyplot as plt
  from matplotlib.pylab import *
                                                        63
  import scipy.integrate as integrate
  import matplotlib.animation as animation
                                                        64
  class DoublePendulum:
8
                                                        65
       def __init__(self,
                     init_state = [0.0, 0.0, 0.0]
10
                                                        66
       0.01.
                                                        67
                    L1 = 1.0,
                                                        68
                    L2 = 1.0,
                    M1 = 1.0,
13
                                                        69
                    M2 = 1.0,
14
                    G = 9.8,
16
                     origin =(0,2):
           self.init_state = np.asarray(init_state,
17
                                                        71
        dtype='float')
           self.params=(L1, L2, M1, M2, G)
           self.origin = origin
19
           self.time_elapsed=0
20
                                                        74
           self.state = self.init_state*np.pi/180
21
                                                        75
22
                                                        76
       def position(self):
23
           (L1, L2, M1, M2, G) = self.params
24
                                                        77
25
                                                        78
           x = np.cumsum([self.origin[0],
26
                                                        79
                           L1 * sin(self.state[0])
27
28
                           L2 * sin(self.state[2])])
           y = np.cumsum([self.origin[1],
29
                           -L1 * cos(self.state[0]),
30
                                                        82
                           -L2 * cos(self.state[2])
31
      ])
           return (x,y)
32
33
       def energy (self):
34
           (L1, L2, M1, M2, G) = self.params
35
           x = np.cumsum([L1 * sin(self.state[0])
36
                           L2 * sin(self.state[2])])
37
38
           y = np.cumsum([-L1 * cos(self.state[0]),
                           -L2 * cos(self.state[2])
39
      1)
40
           vx = np.cumsum([L1 * self.state[1]*cos(
41
       self.state[0]),
                            L2 * self.state[3] * cos
       (self.state[2])])
           vy = np.cumsum([L1 * self.state[1] * sin
      (self.state[0]),
                            L2 * self.state[3] * sin
44
       (self.state[2])])
45
           U = G * (M1 * (2 + y[0]) + M2 * (2+y[1])
46
           K = 0.5 * (M1 * np.dot(vx, vx) + M2 * np
47
       . dot(vy, vy))
48
           return U + K
49
50
51
       def dstate_dt(self, state, t):
           (L1, L2, M1, M2, G) = self.params
53
           dydx = np.zeros_like(state)
54
           dydx[0] = state[1]
55
           dydx[2] = state[3]
56
57
```

```
cos_delta = cos(state[2] - state[0])
              sin_delta = sin(state[2] - state[0])
              den1 = (M1 + M2) * L1 - M2 * L1 *
         cos_delta * cos_delta
              dydx[1] = (M2 * L1 * state[1] * state[1]
          * sin_delta * cos_delta
                           + M2 * G * sin(state[2]) *
         cos_delta
                           + M2 * L2 * state [3] * state
         [3] * sin_delta
                            - (M1 + M2) * G * sin(state)
         [0]))/den1
              \begin{array}{l} den2 \, = \, (L2 \, / \, \, L1) \, * \, den1 \\ dydx \, [3] \, = \, (- \, \, M2 \, * \, \, L2 \, * \, \, state \, [3] \, * \, \, state \end{array}
         [3] * sin_delta * cos_delta
                           + (M1 + M2) * G * sin(state)
         [0]) * cos_delta
                            - (M1 + M2) * L1 * state[1] *
          state[1] * sin_delta
                            - (M1 + M2) * G * sin(state)
         [2])) / den2
              return dvdx
         def step(self, dt):
             self.state = integrate.odeint(self.
         dstate_dt, self.state, [0, dt])[1]
              self.time_elapsed += dt
 pendulum = DoublePendulum ([90.0, 0.0, 90.0,
        0.0])
dt = 1./30
84 # Setup figure and subplots
so fig = figure (num = 0, figsize = (12, 8))#, dpi =
         100)
fig.suptitle("Double Pendulum", fontsize=12) ax01 = subplot2grid((4, 4), (0, 0), colspan=2,
        rowspan=2, autoscale_on=False, aspect =
        equal', xlim = (-2,2), ylim = (0,4)
88 ax01.grid()
89 ax01. set_title ("Simulation")
line, = ax01.plot([],[], 'o-', lw = 2)

time_text = ax01.text(0.02, 0.95, '', transform=
        ax01.transAxes)
energy_text = ax01.text(0.02, 0.90, ', ', ')
        transform = ax01.transAxes)
95 # Data Placeholders
en1=zeros(0)
_{97} t=zeros(0)
98 \text{ xpos}1 = \text{zeros}(0)
99 xpos2 = zeros(0)
ypos1 = zeros(0)
ypos2 = zeros(0)
104 \text{ x}_{-}\text{max} = 5
105 \text{ ax}02 = \text{subplot}2\text{grid}((5, 4), (0, 2), \text{colspan} = 2, \\ \text{rowspan} = 2, \text{ aspect} = 'equal', \text{xlim} =
         (0,15), ylim = (0,50))
106 ax02.grid()
ax02.set_title("Time vs Total Energy")
energygraph, = ax02.plot(t,en1)
```

```
ax03 = subplot2grid((5, 4), (3, 0), colspan = 2,
        rowspan = 2, xlim = (0,15), ylim = (-2,2))
   ax03.grid()
111
   ax03.set_title("Time vs X-Position")
112
   xpos1graph, = ax03.plot(t,xpos1, "g-", label="
       mass 1")
   xpos2graph, = ax03.plot(t, xpos2, "r-", label="
       mass 2")
   ax04 = subplot2grid((5, 4), (3, 2), colspan = 2,
        rowspan = 2, xlim = (0,15), ylim = (0,4))
   ax04.grid()
117
   ax04.set_title("Time vs Y-Position")
   ypos1graph , = ax04.plot(t,ypos1, "g-", label="
       mass 1")
   ypos2graph,
                = ax04.plot(t,ypos2, "r-", label="
       mass 2")
122 # Setting Lagends
   ax03.legend([xpos1graph,xpos2graph], [xpos1graph
123
        . get_label(), xpos2graph.get_label()])
   ax04.legend([ypos1graph,ypos2graph], [ypos1graph
        . get_label(), ypos2graph.get_label()])
126 # Data Update
128
   time1 = 0.0
129
130
   def init():
        line.set_data([], [])
        time_text.set_text(
        energy_text.set_text('')
135
136
        return line, time_text, energy_text
137
138
139
   def animate(i):
140
        {\color{red} {\bf global} \ \ pendulum} \ , \ dt \ , \ time1 \ , \ en1 \ , \ t \ , \ x\_max \ ,
141
       xpos1, xpos2, ypos1, ypos2
142
       pendulum.step(dt)
143
        line.set_data(*pendulum.position())
144
        time_text.set_text('time = %.1f', % pendulum.
145
        time_elapsed)
        energy_text.set_text('energy = \%.3 f' %
146
       pendulum.energy())
       en1 = append(en1, pendulum.energy())
148
        t = append(t, time1)
149
        time1 += dt
       energygraph.set_data(array(t), array(en1))
        xpos1 = append(xpos1, pendulum.position()
154
        [0][1]
        xpos2 = append(xpos2, pendulum.position()
        [0][2])
        xpos1graph.set_data(t, xpos1)
       xpos2graph.set_data(t, xpos2)
158
```

ypos1 = append(ypos1, pendulum.position()

```
[1][1])
       ypos2 = append(ypos2, pendulum.position()
       [1][2]
161
       ypos1graph.set_data(t, ypos1)
       ypos2graph.set_data(t, ypos2)
162
163
       \#if time1 >= x_max - 1.00:
164
           #energygraph.axes.set_xlim(time1 - x_max
        + 1.0, time1 + 1.0)
           \#xpos1graph.axes.set\_xlim(time1 - x\_max)
         1.0, time1 + 1.0)
167
       return line , time_text , energy_text ,
       energygraph, xpos1graph, xpos2graph,
       ypos1graph, ypos2graph
169
170
   from time import time
171
172
173
t0 = time()
animate (0)
t1 = time()
   interval = 1000 * dt - (t1 - t0)
178
ani = animation.FuncAnimation(fig, animate,
       frames = 400,
                                   interval =
180
       interval, blit = False,
181
                                   init_func=init)
182
#ani.save("p90_p90.gif", writer="imagemagick",
       fps = 30)
185 plt.show()
```

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