

Numerical Simulation of the 2-Body Problem

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1 Numerical Simulation of the 2-Body Problem

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

1.1 i)

```
[ ]: def euler(w0,s0,h,e):
    #the function will take the initial velocity
    # initial position, the whished step size
    # and the desired upper bondary for the integral

    G = 1 #grativational constant
    m = 1 #mass of both bodies
    N = int(e/h) #time steps

    s = np.zeros((N,2))
    w = np.zeros((N,2))
    LRL = np.zeros((N,2))
    s[0] = s0
    w[0] = w0
    #leap-frog algorithm
    for i in range(N-1):
        s[i+1] = s[i] + h*w[i]
        w[i+1] = w[i]-h*s[i]/np.linalg.norm(s[i])**3
        #calculatin LRL for every time step
        LRL[i] = s[i]*(np.dot(w[i],w[i])) - w[i]*(np.dot(s[i],w[i])) - s[i]/np.
        ↪linalg.norm(s[i])

    #calculating eccentricity with LRL
    epsilon = np.mean([np.linalg.norm(x) for x in LRL])
    print(np.linalg.norm(LRL[1]))
    #calculatin eccentricity geometrically
    a = max(s[:,0])- min(s[:,0])
```

```

b = max(s[:,1])-min(s[:,1])
print(a)
print(b)
epsilon2 = np.sqrt(abs(1-a**2/b**2))
print('epislon calculated geometrically {:.3}'.format(epsilon2))
#plot
fig1, ax1 = plt.subplots()
t = np.linspace(0,e,N)
ax1.plot(t,s[:,1], label= r'\omega$_0$ = {} $\quad s$_0$ = {},$\quad$ $';
↪ $ dt = {}'.format(w0,s0,h))
plt.title('Motion of 2-body problem body problem with $\epsilon$ = {:.3}$'.
↪ format(epsilon))
plt.xlabel(' time t')
plt.ylabel('relative position $s_1$')
plt.legend(loc = 8)
fig2, ax2 = plt.subplots()
ax2.plot(s[:,0], s[:,1], label= r'\omega$_0$ = {} $\quad s$_0$ = {}
↪ {},$\quad$ $'; $ dt = {}'.format(w0,s0,h))
plt.title('Orbit of the 2-body problem with $\epsilon$ = {:.3}$'.
↪ format(epsilon))
plt.xlabel('relative position $s_1$')
plt.ylabel('relative position $s_2$')
plt.legend(loc = 8)
return s,w

```

```

[ ]: w1 = [0.0,1]
s1=[1.0,0.0]
euler(w1,s1,0.001,30)

```

```

1.5000003749539096e-06
2.1054507526942885
2.0993082373604404
epislon calculated geometrically :0.0766

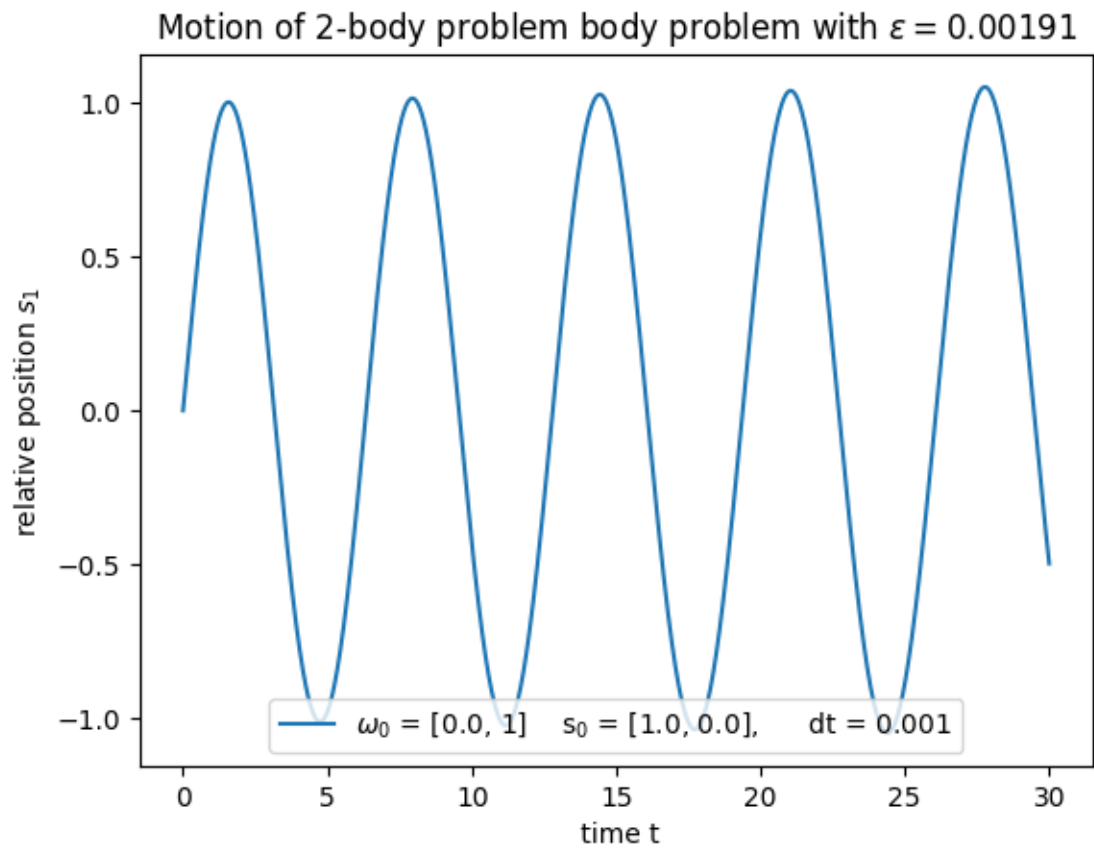
```

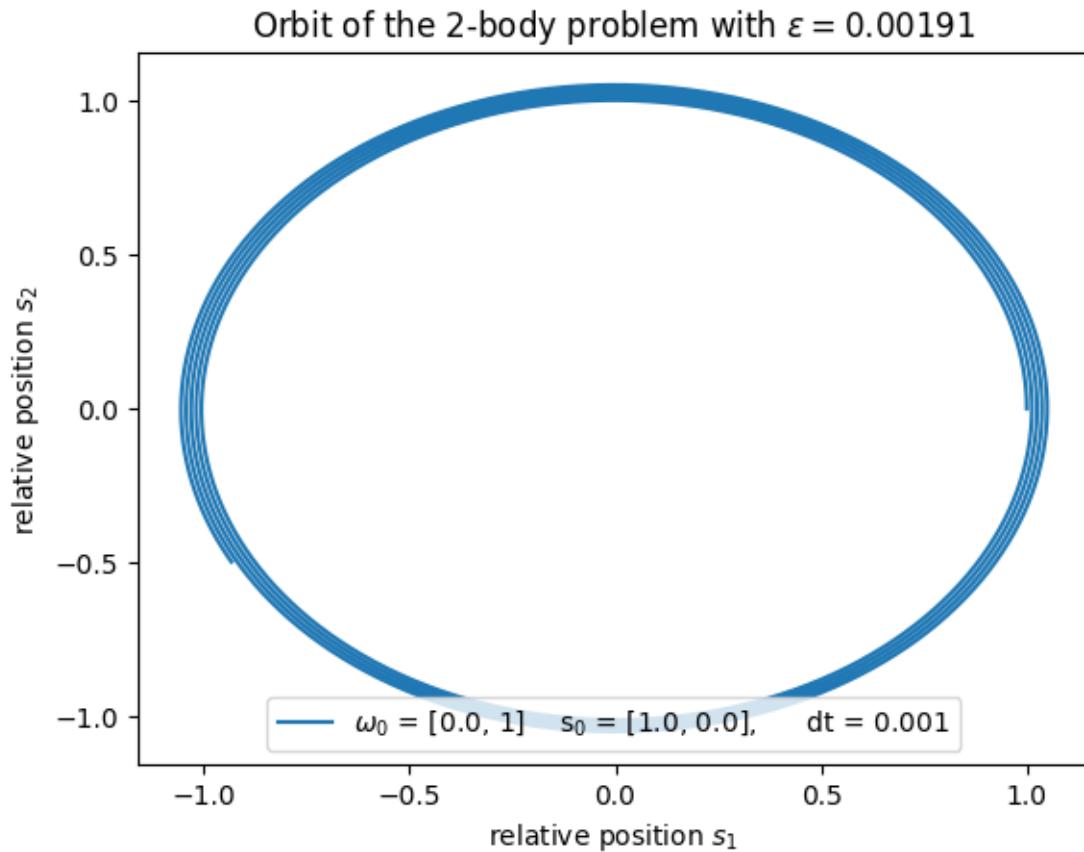
```

[ ]: (array([[ 1.          ,  0.          ],
             [ 1.          ,  0.001       ],
             [ 0.999999   ,  0.002       ],
             ...,
             [-0.93439238, -0.49522218],
             [-0.93393954, -0.49608233],
             [-0.93348591, -0.49694207]]),
array([[ 0.          ,  1.          ],
       [-0.001       ,  1.          ],
       [-0.002       ,  0.999999   ],
       ...,
       [ 0.4528374   , -0.86015207],
       [ 0.45362748  , -0.85973333],

```

[0.45441717, -0.85931387]]))





```
[ ]: def leap_frog(w0,s0,h,e):
    #the function will take the initial velocity
    # initial position, the whished step size
    # and the desired upper bondary for the integral
    G = 1 #grativational constant
    m = 1 #mass of both bodies
    N = int(e/h) #time steps

    s = np.zeros((N,2))
    w = np.zeros((N,2))
    LRL = np.zeros((N,2))

    w[0] = w0
    s[0] = s0

    #First the programm will compute the s_1/2 step using the
    #euler forward methode
    a = 0.5*h/np.linalg.norm(s0)**3
    wn = w[0] - h/2*s[0]/np.linalg.norm(s[0])
```

```

s[0] = s0 + h*wn

#leap-frog algorithm
for i in range(N-1):
    w[i + 1] = w[i]-h/2*s[i]/np.linalg.norm(s[i])**3
    s[i + 1] = s[i] + w[i]*h
    #calculatin LRL for every time step
    LRL[i] = s[i]*(np.dot(w[i],w[i])) - w[i]*(np.dot(s[i],w[i])) - s[i]/np.
↪linalg.norm(s[i])

#calculating eccentricisity with LRL
epsilon = np.mean([np.linalg.norm(x) for x in LRL])
#calculatin eccentricisity geometrically
a = max(s[:,0])-min(s[:,0])
b = max(s[:,1])-min(s[:,1])
print(a)
print(b)
epsilon2 = np.sqrt(abs(1-b**2/a**2))
print('epislron calculated geometrically {:.3}'.format(epsilon2))
#plot
fig1, ax1 = plt.subplots()
t = np.linspace(0,e,N)
ax1.plot(t,s[:,1], label= r'\omega$_0$ = {} $\quad s$_0$ = {},$\quad$ $ \backslash$;
↪$ dt = {}'.format(w0,s0,h))
plt.title('Motion of 2 body problem with $\epsilon$ = {:.3}'.
↪format(epsilon))
plt.xlabel(' time t')
plt.ylabel('relative position $s_1$')
plt.legend(loc = 8)
fig2, ax2 = plt.subplots()
ax2.plot(s[:,0], s[:,1], label= r'\omega$_0$ = {} $\quad s$_0$ = {}
↪$, $\quad$ $ \backslash$; $ dt = {}'.format(s0,w0,h))
plt.title('Orbit of the 2-body problem with $\epsilon$ = {:.3}'.
↪format(epsilon))
plt.xlabel('relative position $s_1$')
plt.ylabel('relative position $s_2$')
plt.legend(loc = 8)

```

```

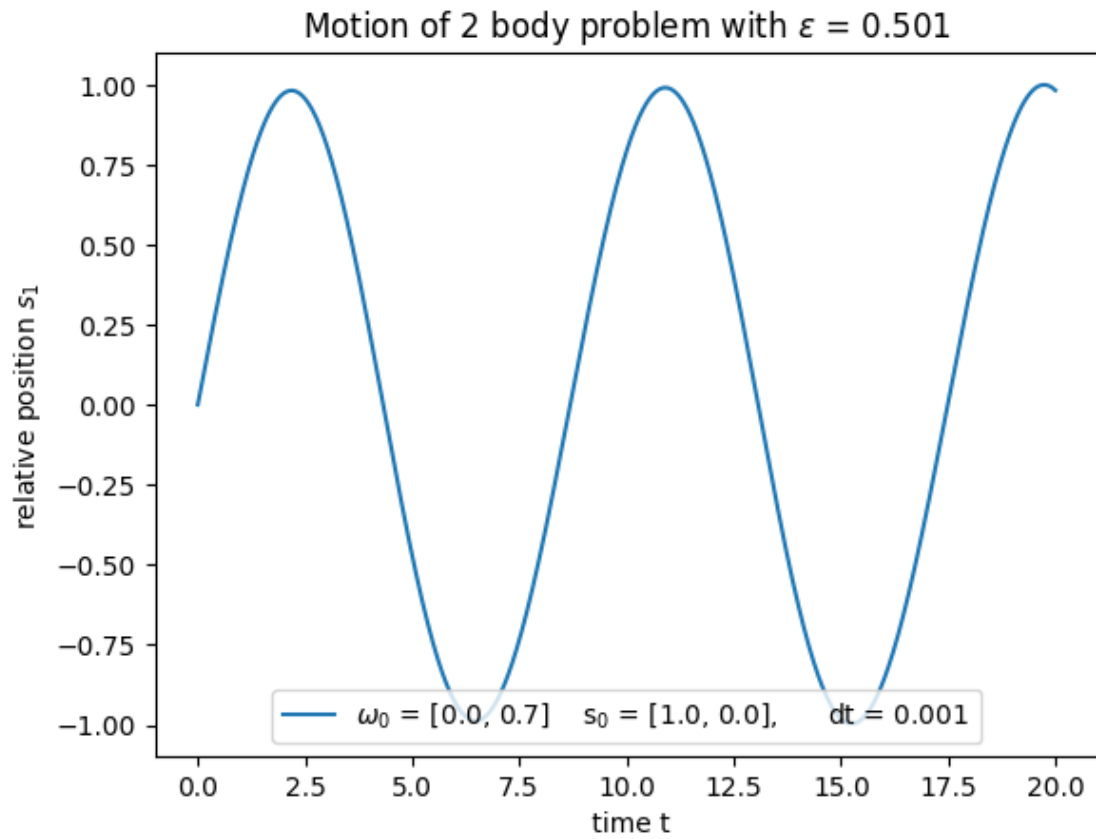
[ ]: w2 = [0.0,.70]
s2=[1.0,0.0]
leap_frog(w2,s2,0.001,20)

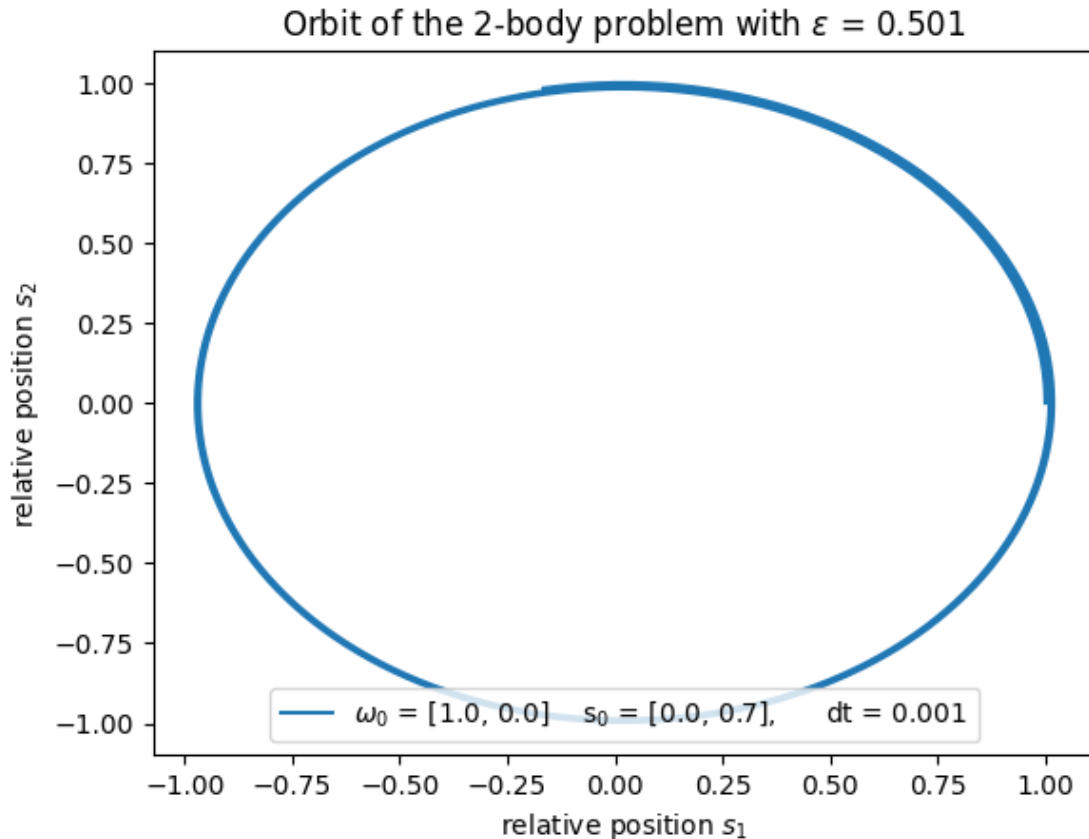
```

```

1.992265652834932
1.9961493469569298
epislron calculated geometrically :0.0625

```





2 ii)

After changing the parameters of the Euler Forward Method it has been established, that due to the errors of this method it is not possible to get a perfect circle with radius 1. However with $\omega_0 = (0, 1.0)$ the program delivers a good approximation. In the other hand, for the leap-frog algorithm to deliver a good approximation the initial velocity has to be set to $\omega_0 = (0, 0.7)$

2.1 iii)

```
[ ]: w1 = [0.0,1]
      s1=[1.0,0.0]
      euler(w1,s1,0.01,30)
```

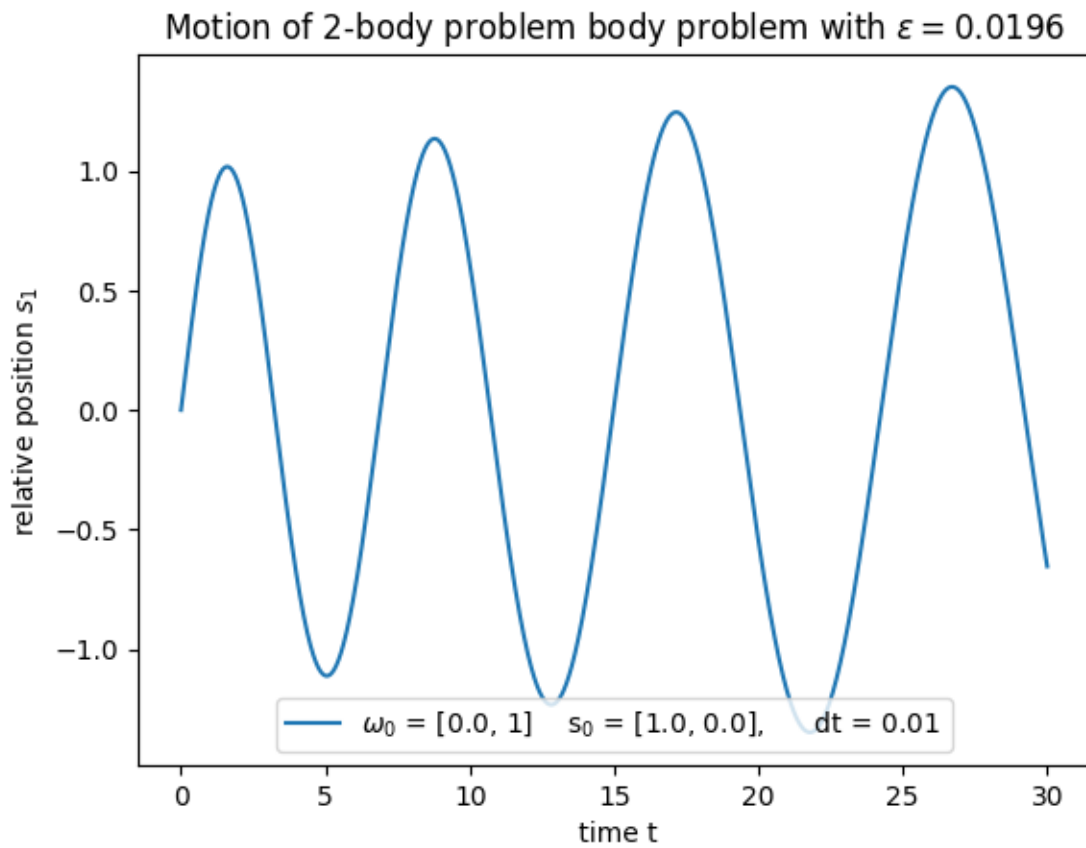
0.00015000374993734167

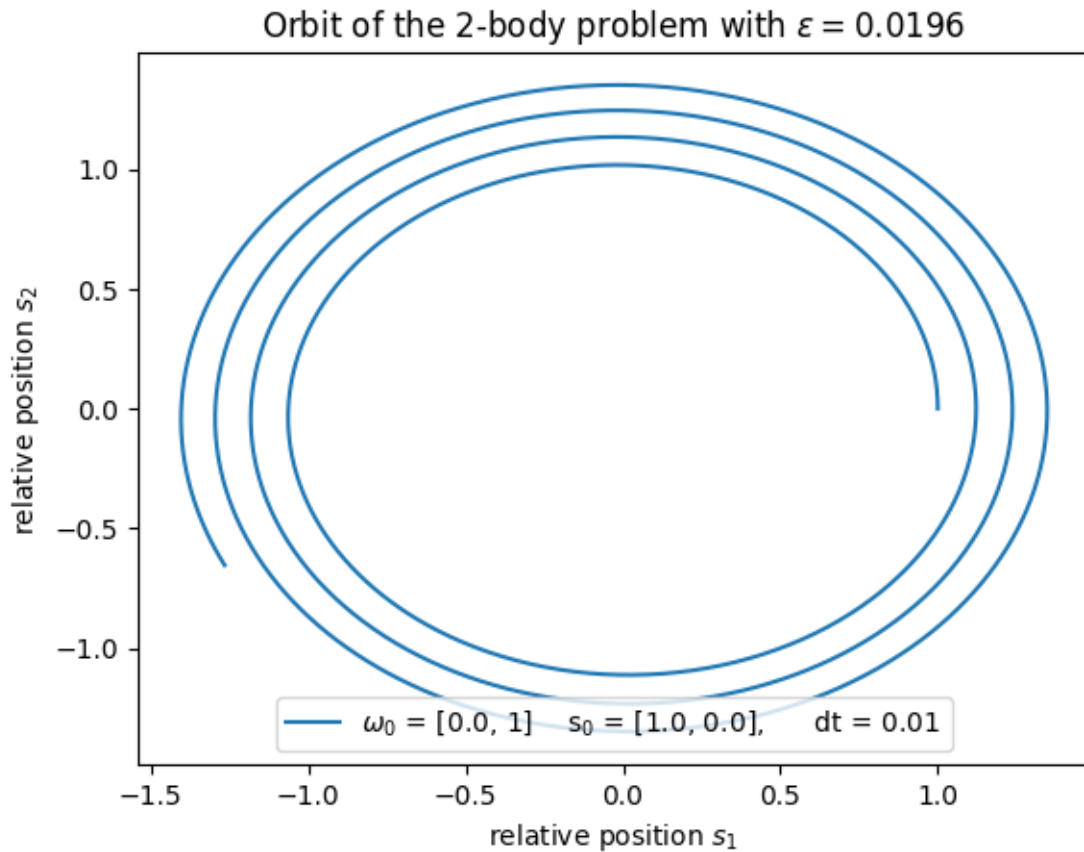
2.7516943302449928

2.6964787836941415

epsilon calculated geometrically :0.203

```
[ ]: (array([[ 1.          ,  0.          ],
            [ 1.          ,  0.01         ],
            [ 0.9999       ,  0.02         ],
            ...,
            [-1.27251341, -0.63774383],
            [-1.26897081, -0.64530039],
            [-1.26538408, -0.65283482]]),
array([[ 0.          ,  1.          ],
       [-0.01         ,  1.          ],
       [-0.0199985    ,  0.99990001],
       ...,
       [ 0.35426036, -0.75565525],
       [ 0.35867302, -0.75344376],
       [ 0.36307115, -0.75120722]]))
```



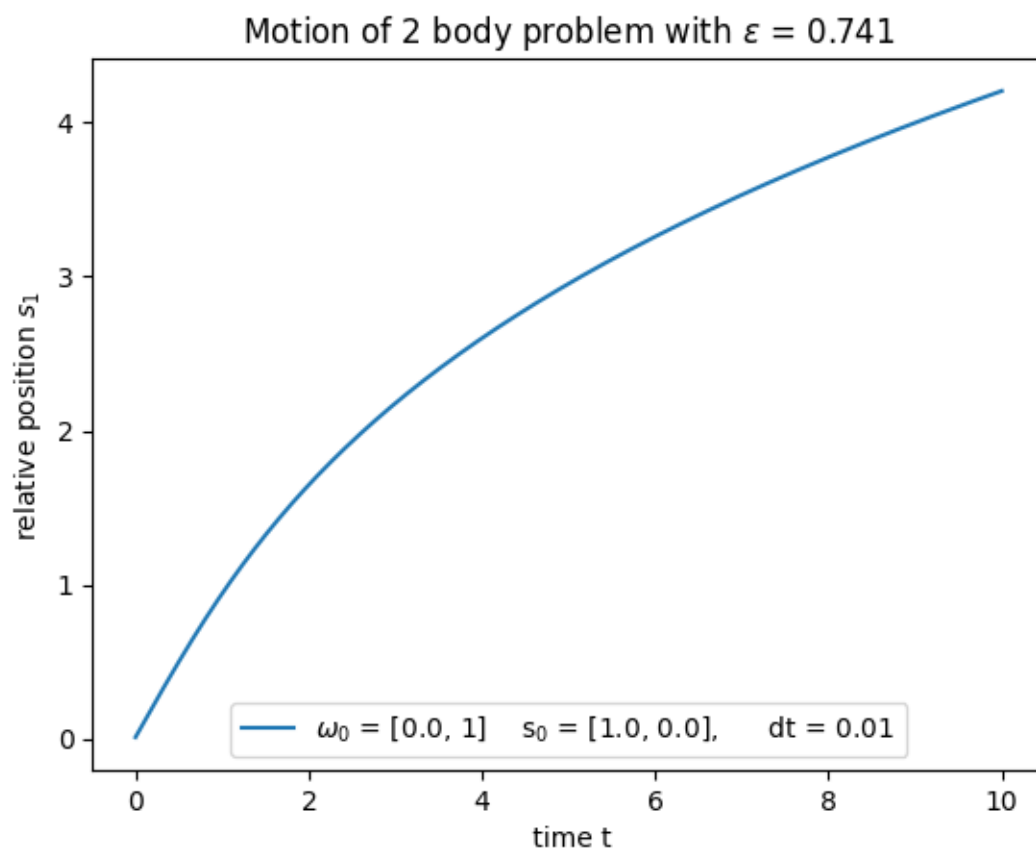


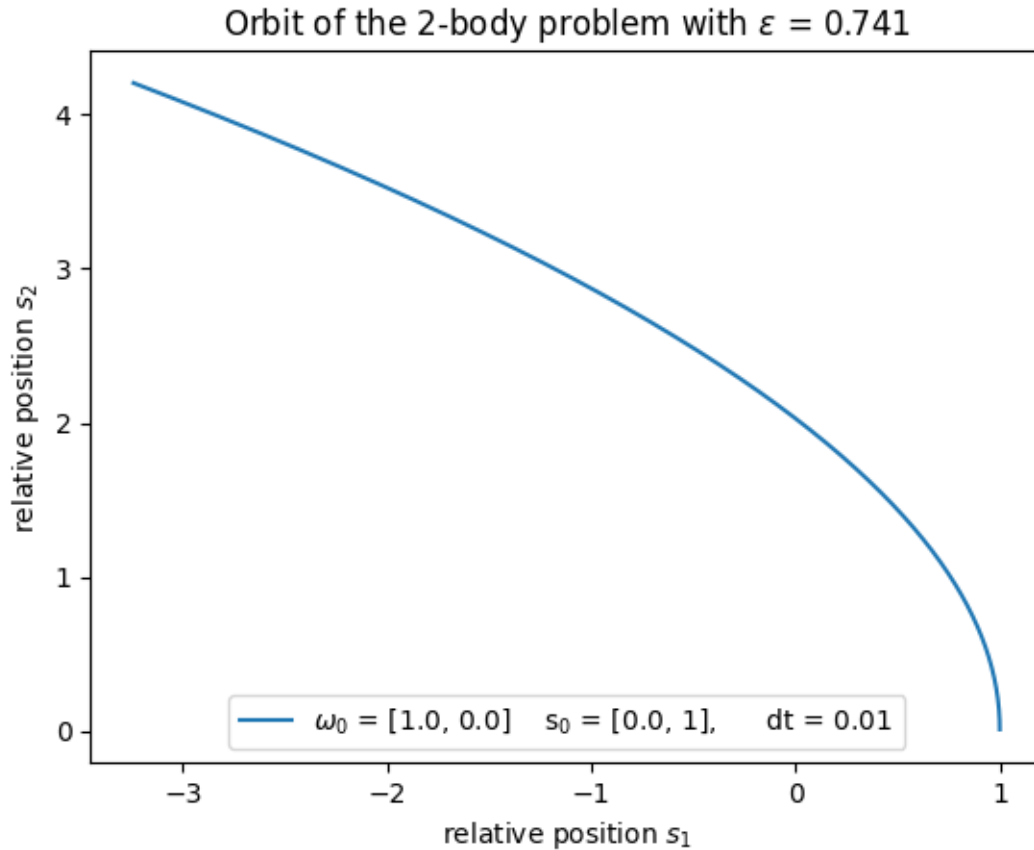
```
[ ]: leap_frog(w1,s1,0.01,10)
      #for the leap_frog algorithm the increase in the distance is lower about 0.01%
```

4.241240714984339

4.1951887488190645

epilson calculated geometrically :0.147





If the time step is increased the exactitude of both methods decrease greatly. The Euler Method plots a piral trajectory, meanwhile the leap-frog algorithm shows a hyperbola.

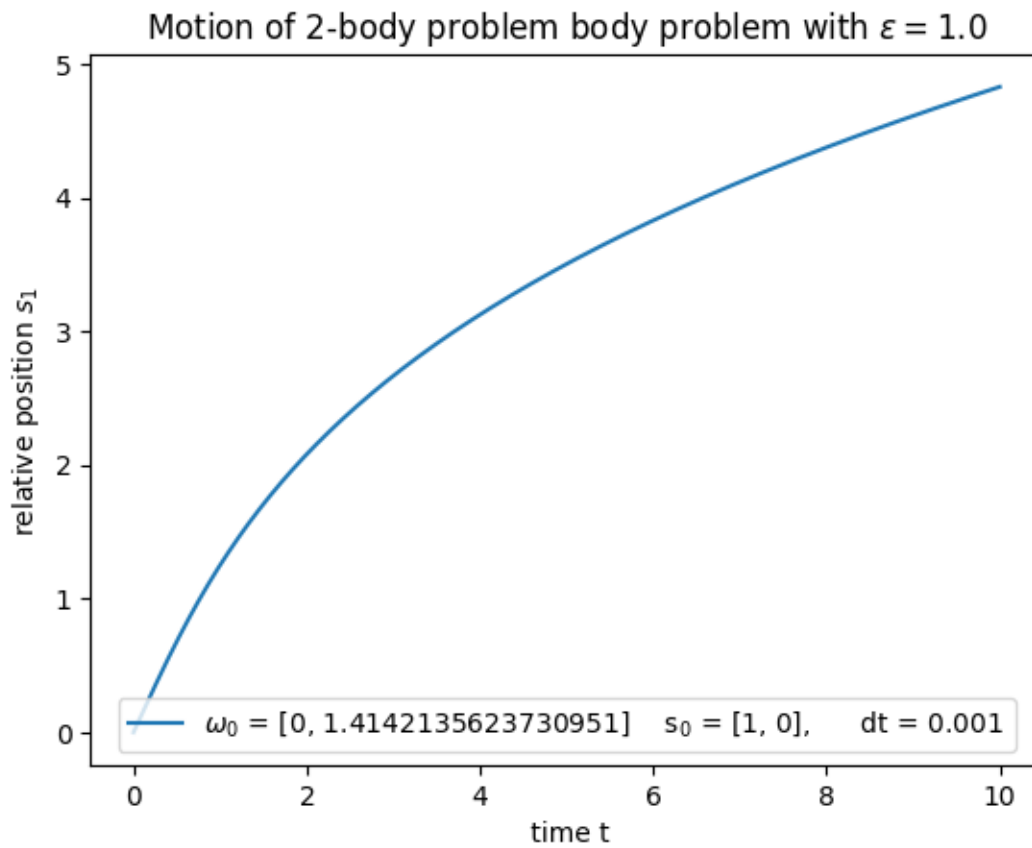
2.2 v)

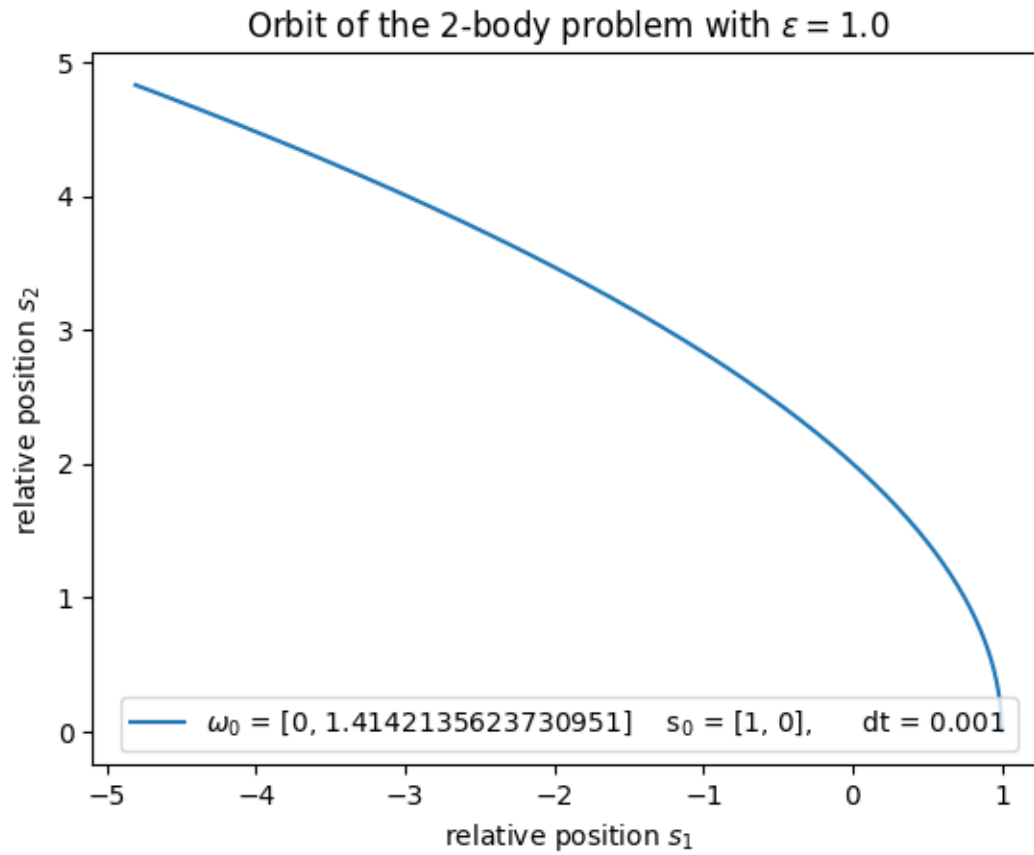
```
[ ]: euler([0,np.sqrt(2)],[1,0], 0.001,10)
1.0000029999985007
5.808259981521099
4.827925218332602
epsilon calculated geometrically :0.669
[ ]: (array([[ 1.00000000e+00,  0.00000000e+00],
             [ 1.00000000e+00,  1.41421356e-03],
             [ 9.99999000e-01,  2.82842712e-03],
             ...,
             [-4.80725699e+00,  4.82750700e+00],
             [-4.80775849e+00,  4.82771612e+00],
             [-4.80825998e+00,  4.82792522e+00]]),
      array([[ 0.00000000e+00,  1.41421356e+00],
```

```

[-1.00000000e-03,  1.41421356e+00],
[-1.99999700e-03,  1.41421215e+00],
...,
[-5.01504064e-01,  2.09114958e-01],
[-5.01488862e-01,  2.09099691e-01],
[-5.01473661e-01,  2.09084427e-01]]))

```



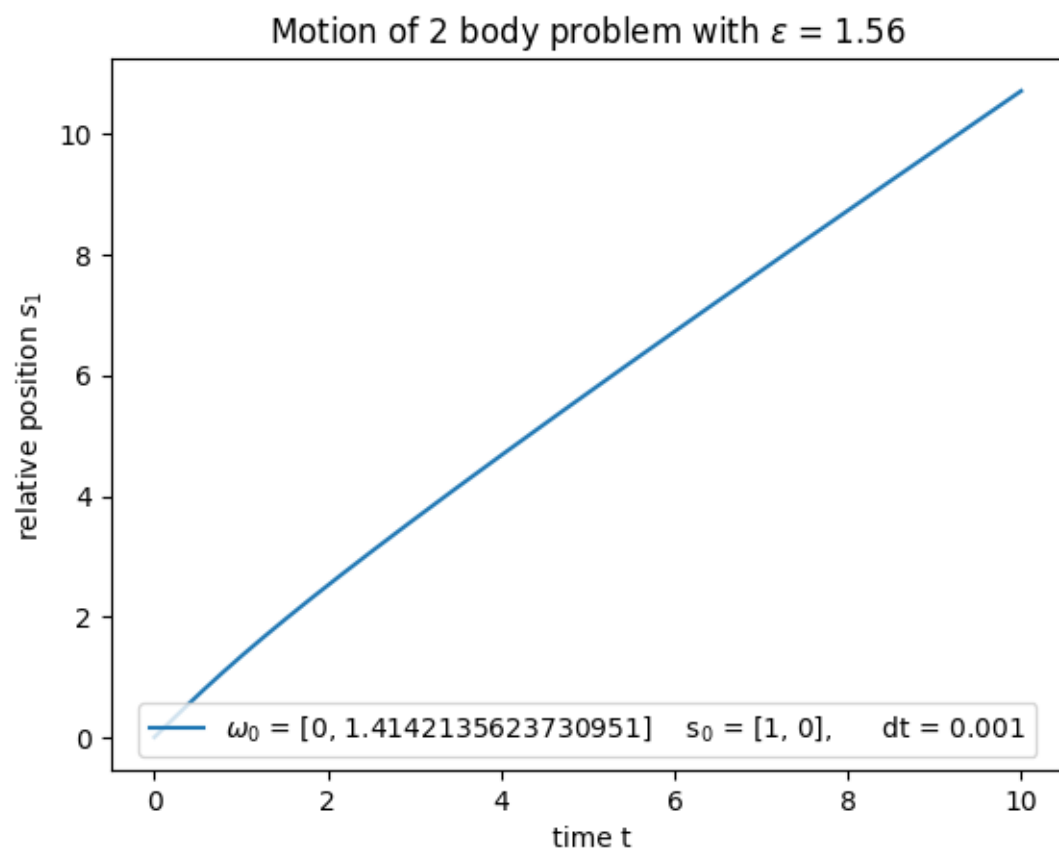


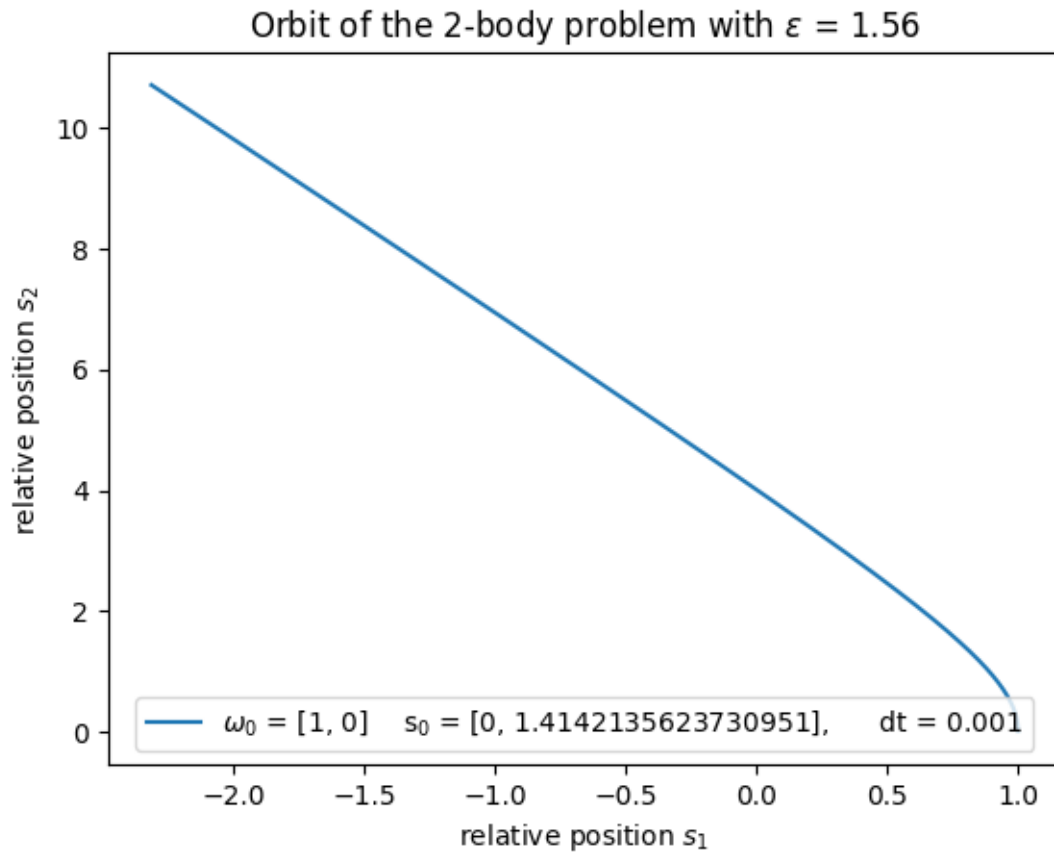
```
[ ]: leap_frog([0,np.sqrt(2)],[1,0], 0.001,10)
```

```
3.3165115257055158
```

```
10.712525087721502
```

```
epsilon calculated geometrically :3.07
```





```
[ ]: euler([0,1/np.sqrt(2)], [1,0], 0.0001,10)
```

```
0.49999999250000005
```

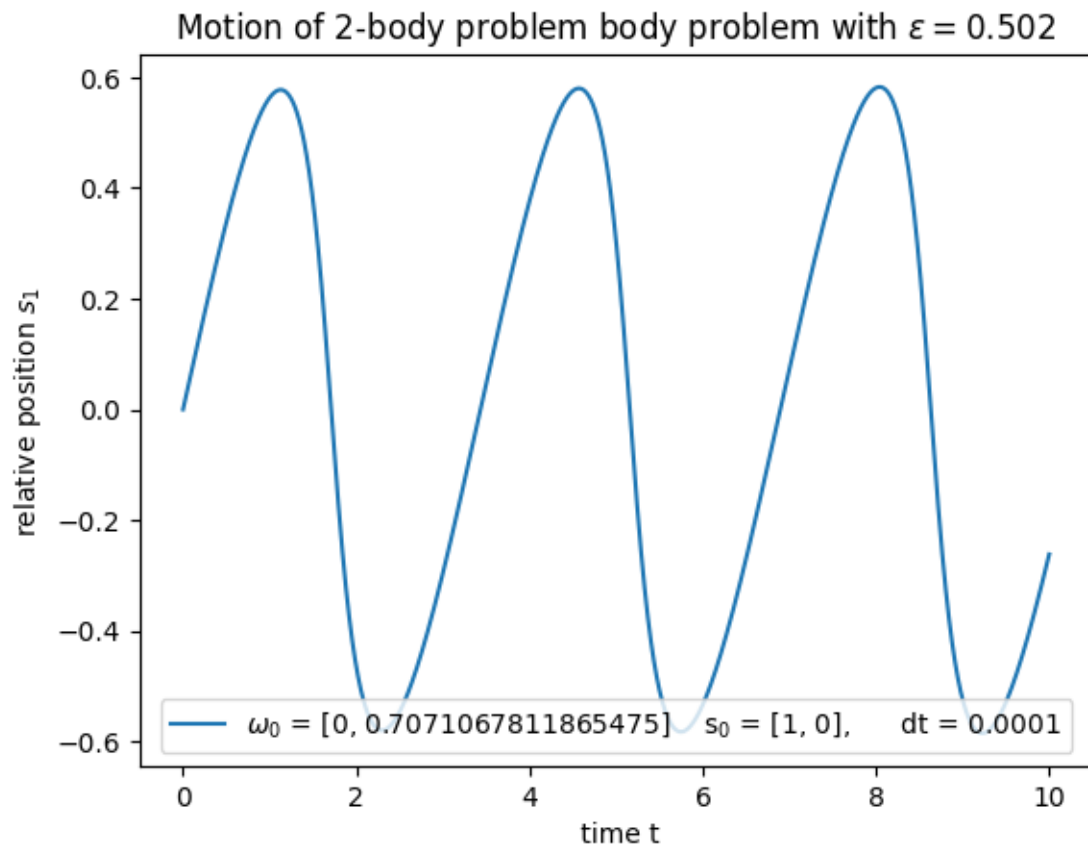
```
1.3480258576967694
```

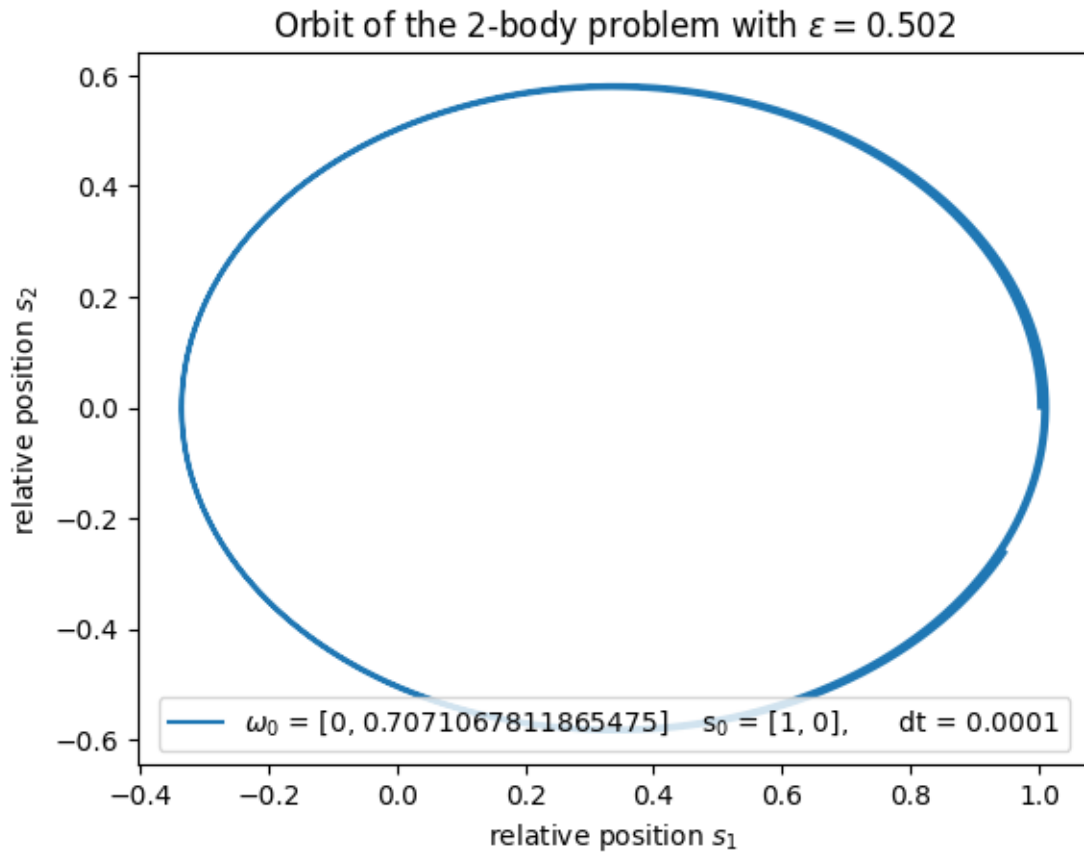
```
1.1675373818253267
```

```
epsilon calculated geometrically :0.577
```

```
[ ]: (array([[ 1.00000000e+00,  0.00000000e+00],
             [ 1.00000000e+00,  7.07106781e-05],
             [ 9.99999990e-01,  1.41421356e-04],
             ...,
             [ 9.47099038e-01, -2.61851117e-01],
             [ 9.47136525e-01, -2.61786428e-01],
             [ 9.47174002e-01, -2.61721736e-01]]),
      array([[ 0.00000000e+00,  7.07106781e-01],
             [-1.00000000e-04,  7.07106781e-01],
             [-1.99999999e-04,  7.07106774e-01],
             ...,
             [ 3.74868772e-01,  6.46891334e-01],
             [ 3.74768950e-01,  6.46918932e-01],
```

[3.74669130e-01, 6.46946522e-01]]))



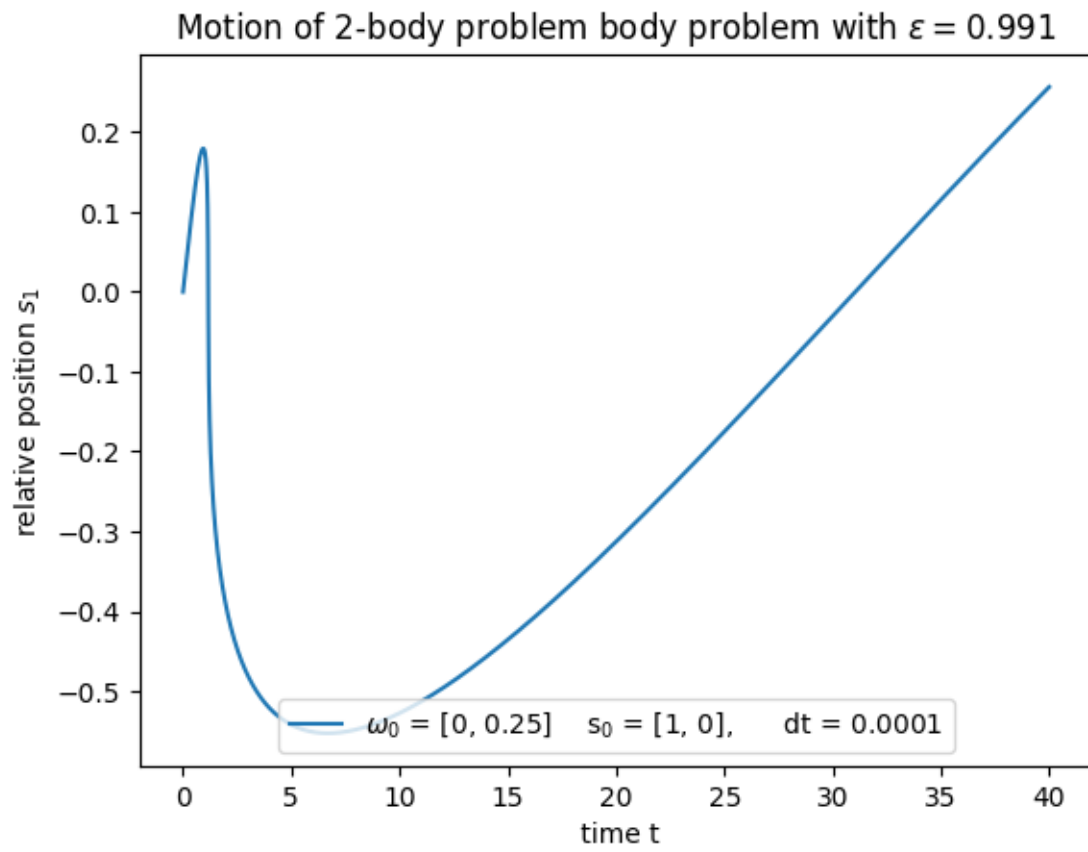


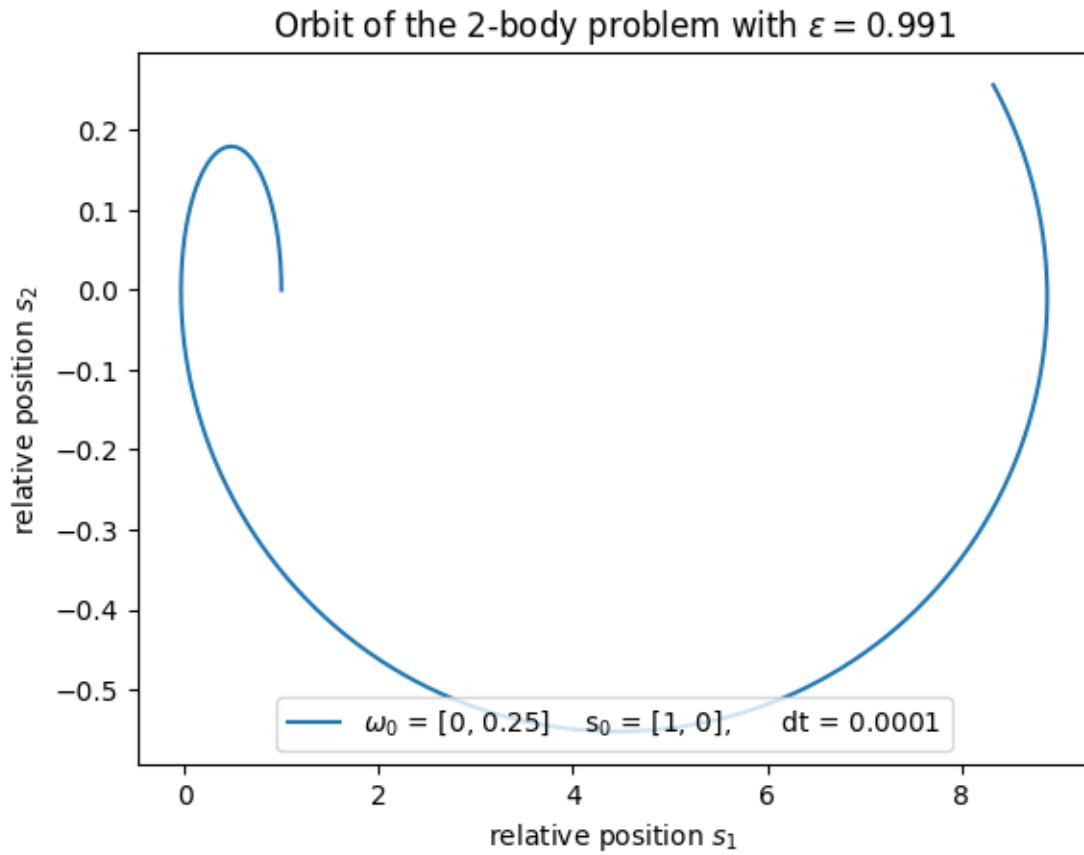
```
[ ]: euler([0,1/4],[1,0], 0.0001,40)
```

```
0.9374999990624999
8.916320592151196
0.8093261724427312
epsilon calculated geometrically :11.0
```

```
[ ]: (array([[1.00000000e+00, 0.00000000e+00],
            [1.00000000e+00, 2.50000000e-05],
            [9.99999990e-01, 5.00000000e-05],
            ...,
            [8.32886475e+00, 2.56626550e-01],
            [8.32885251e+00, 2.56629291e-01],
            [8.32884027e+00, 2.56632031e-01]]),
      array([[ 0.00000000e+00,  2.50000000e-01],
            [-1.00000000e-04,  2.50000000e-01],
            [-2.00000000e-04,  2.49999998e-01],
            ...,
            [-1.22373708e-01,  2.74071714e-02],
            [-1.22375148e-01,  2.74071271e-02],
```

$[-1.22376587\text{e-}01, \ 2.74070827\text{e-}02]]))$



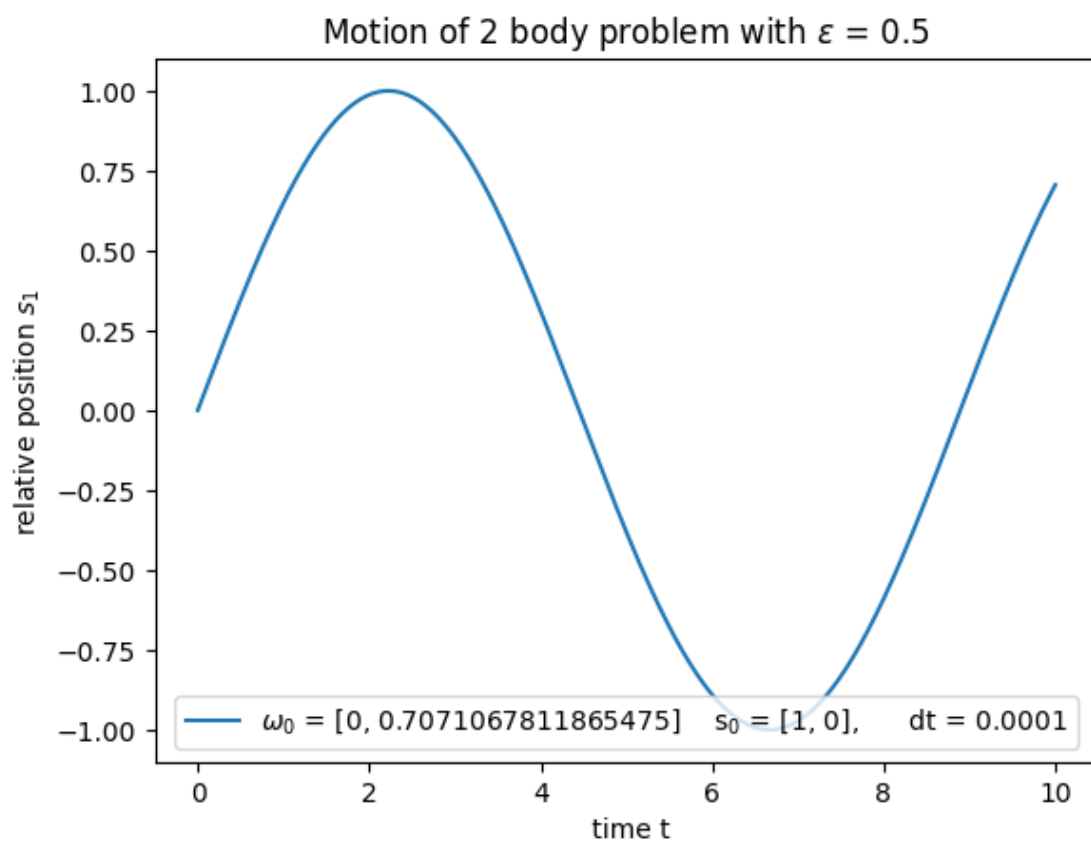


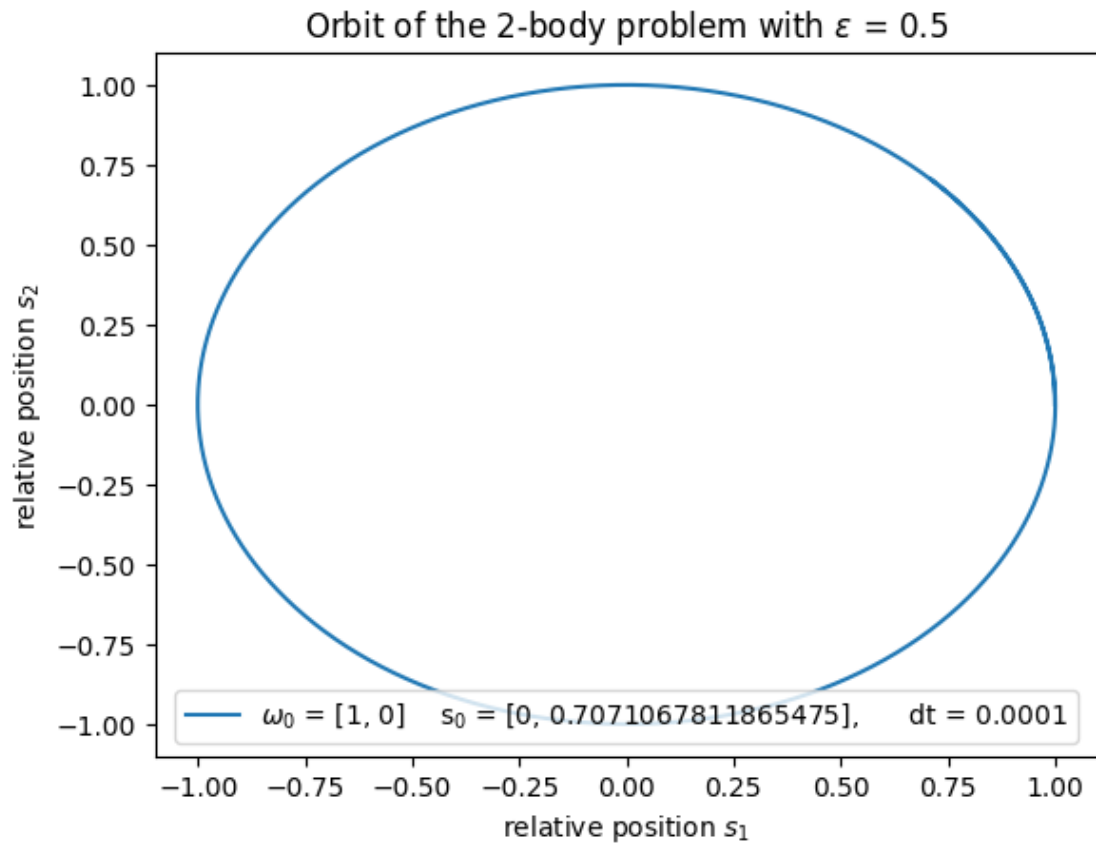
```
[ ]: leap_frog([0,1/np.sqrt(2)], [1,0], 0.0001, 10)
```

```
2.001332646540053
```

```
2.000888510189057
```

```
epsilon calculated geometrically :0.0211
```



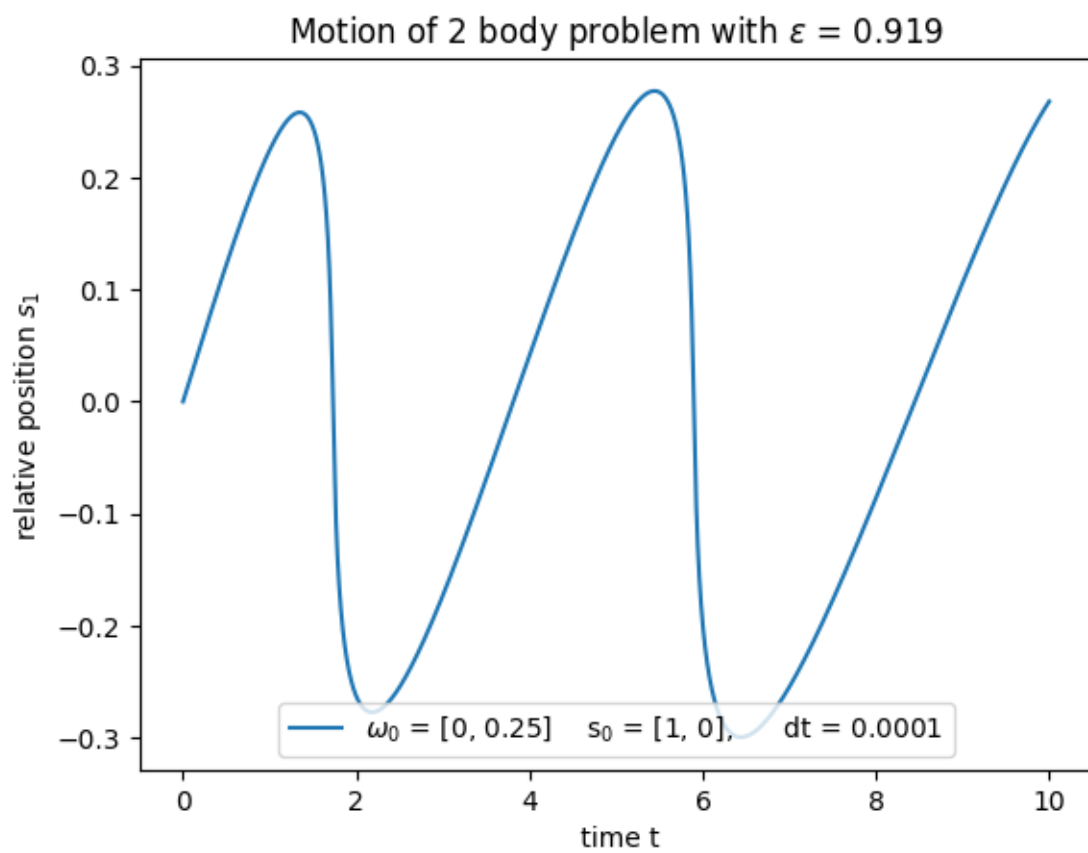


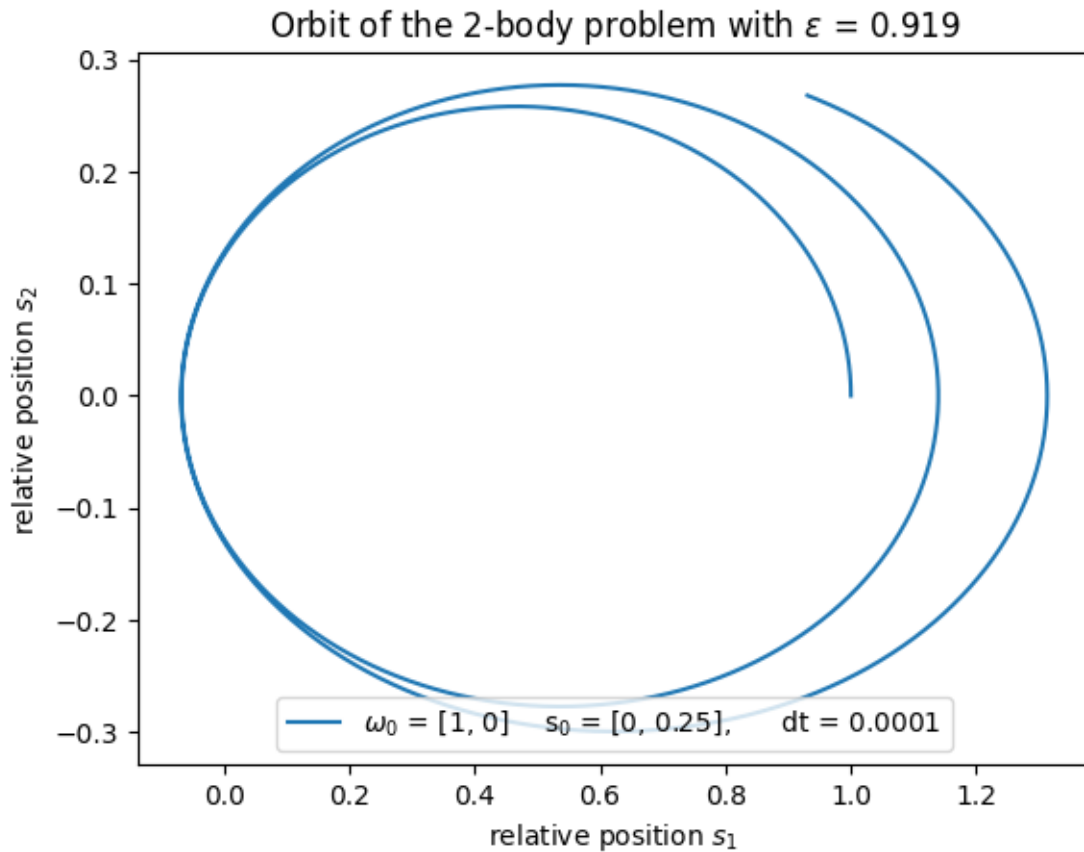
```
[ ]: leap_frog([0,1/4],[1,0], 0.0001,10)
```

```
1.3808830743106262
```

```
0.5770487824456738
```

```
epsilon calculated geometrically :0.909
```





3 Error Analysis

3.1 a) Euler methode

```
[ ]: euler([0.0,1.0],[1,0], 0.1,30)
```

```
0.015037437733209797
4.752794399746748
5.07943758093948
epsilon calculated geometrically :0.353
```

```
[ ]: (array([[ 1.00000000e+00,  0.00000000e+00],
 [ 1.00000000e+00,  1.00000000e-01],
 [ 9.90000000e-01,  2.00000000e-01],
 [ 9.70148147e-01,  2.99014815e-01],
 [ 9.40687451e-01,  3.96088449e-01],
 [ 9.01954090e-01,  4.90304103e-01],
 [ 8.54374015e-01,  5.80794736e-01],
 [ 7.98457639e-01,  6.66753737e-01],
```

[7.34792507e-01, 7.47445214e-01],
 [6.64033986e-01, 8.22213342e-01],
 [5.86894173e-01, 8.90490295e-01],
 [5.04129334e-01, 9.51802285e-01],
 [4.16526338e-01, 1.00577338e+00],
 [3.24888623e-01, 1.05212687e+00],
 [2.30022234e-01, 1.09068418e+00],
 [1.32722525e-01, 1.12136137e+00],
 [3.37619908e-02, 1.14416352e+00],
 [-6.61203620e-02, 1.15917729e+00],
 [-1.66227826e-01, 1.16656226e+00],
 [-2.65912845e-01, 1.16654122e+00],
 [-3.64581885e-01, 1.15939018e+00],
 [-4.61698396e-01, 1.14542831e+00],
 [-5.56784057e-01, 1.12500823e+00],
 [-6.49418509e-01, 1.09850693e+00],
 [-7.39237845e-01, 1.06631757e+00],
 [-8.25932122e-01, 1.02884209e+00],
 [-9.09242149e-01, 9.86484978e-01],
 [-9.88955794e-01, 9.39647949e-01],
 [-1.06490399e+00, 8.88725588e-01],
 [-1.13695665e+00, 8.34101917e-01],
 [-1.20501859e+00, 7.76147744e-01],
 [-1.26902553e+00, 7.15218725e-01],
 [-1.32894041e+00, 6.51654023e-01],
 [-1.38474982e+00, 5.85775488e-01],
 [-1.43646077e+00, 5.17887249e-01],
 [-1.48409775e+00, 4.48275646e-01],
 [-1.52770009e+00, 3.77209434e-01],
 [-1.56731955e+00, 3.04940182e-01],
 [-1.60301824e+00, 2.31702839e-01],
 [-1.63486676e+00, 1.57716401e-01],
 [-1.66294257e+00, 8.31846496e-02],
 [-1.68732860e+00, 8.29694296e-03],
 [-1.70811201e+00, -6.67709755e-02],
 [-1.72538318e+00, -1.41856164e-01],
 [-1.73923477e+00, -2.16807680e-01],
 [-1.74976099e+00, -2.91485793e-01],
 [-1.75705695e+00, -3.65761231e-01],
 [-1.76121809e+00, -4.39514454e-01],
 [-1.76233980e+00, -5.12634966e-01],
 [-1.76051696e+00, -5.85020663e-01],
 [-1.75584371e+00, -6.56577222e-01],
 [-1.74841314e+00, -7.27217523e-01],
 [-1.73831713e+00, -7.96861116e-01],
 [-1.72564619e+00, -8.65433716e-01],
 [-1.71048932e+00, -9.32866742e-01],

[-1.69293394e+00, -9.99096883e-01],
 [-1.67306581e+00, -1.06406570e+00],
 [-1.65096899e+00, -1.12771924e+00],
 [-1.62672586e+00, -1.19000773e+00],
 [-1.60041704e+00, -1.25088522e+00],
 [-1.57212145e+00, -1.31030931e+00],
 [-1.54191631e+00, -1.36824091e+00],
 [-1.50987716e+00, -1.42464391e+00],
 [-1.47607789e+00, -1.47948506e+00],
 [-1.44059081e+00, -1.53273367e+00],
 [-1.40348664e+00, -1.58436145e+00],
 [-1.36483461e+00, -1.63434237e+00],
 [-1.32470246e+00, -1.68265243e+00],
 [-1.28315655e+00, -1.72926954e+00],
 [-1.24026183e+00, -1.77417340e+00],
 [-1.19608200e+00, -1.81734535e+00],
 [-1.15067948e+00, -1.85876825e+00],
 [-1.10411549e+00, -1.89842640e+00],
 [-1.05645011e+00, -1.93630542e+00],
 [-1.00774236e+00, -1.97239216e+00],
 [-9.58050197e-01, -2.00667463e+00],
 [-9.07430618e-01, -2.03914193e+00],
 [-8.55939692e-01, -2.06978416e+00],
 [-8.03632618e-01, -2.09859237e+00],
 [-7.50563773e-01, -2.12555851e+00],
 [-6.96786768e-01, -2.15067538e+00],
 [-6.42354495e-01, -2.17393655e+00],
 [-5.87319173e-01, -2.19533639e+00],
 [-5.31732403e-01, -2.21486994e+00],
 [-4.75645209e-01, -2.23253296e+00],
 [-4.19108085e-01, -2.24832185e+00],
 [-3.62171042e-01, -2.26223364e+00],
 [-3.04883653e-01, -2.27426598e+00],
 [-2.47295091e-01, -2.28441711e+00],
 [-1.89454177e-01, -2.29268583e+00],
 [-1.31409419e-01, -2.29907151e+00],
 [-7.32090501e-02, -2.30357407e+00],
 [-1.49010736e-02, -2.30619397e+00],
 [4.34667029e-02, -2.30693223e+00],
 [1.01846627e-01, -2.30579039e+00],
 [1.60191167e-01, -2.30277054e+00],
 [2.18452870e-01, -2.29787529e+00],
 [2.76584334e-01, -2.29110784e+00],
 [3.34538168e-01, -2.28247193e+00],
 [3.92266959e-01, -2.27197185e+00],
 [4.49723240e-01, -2.25961250e+00],
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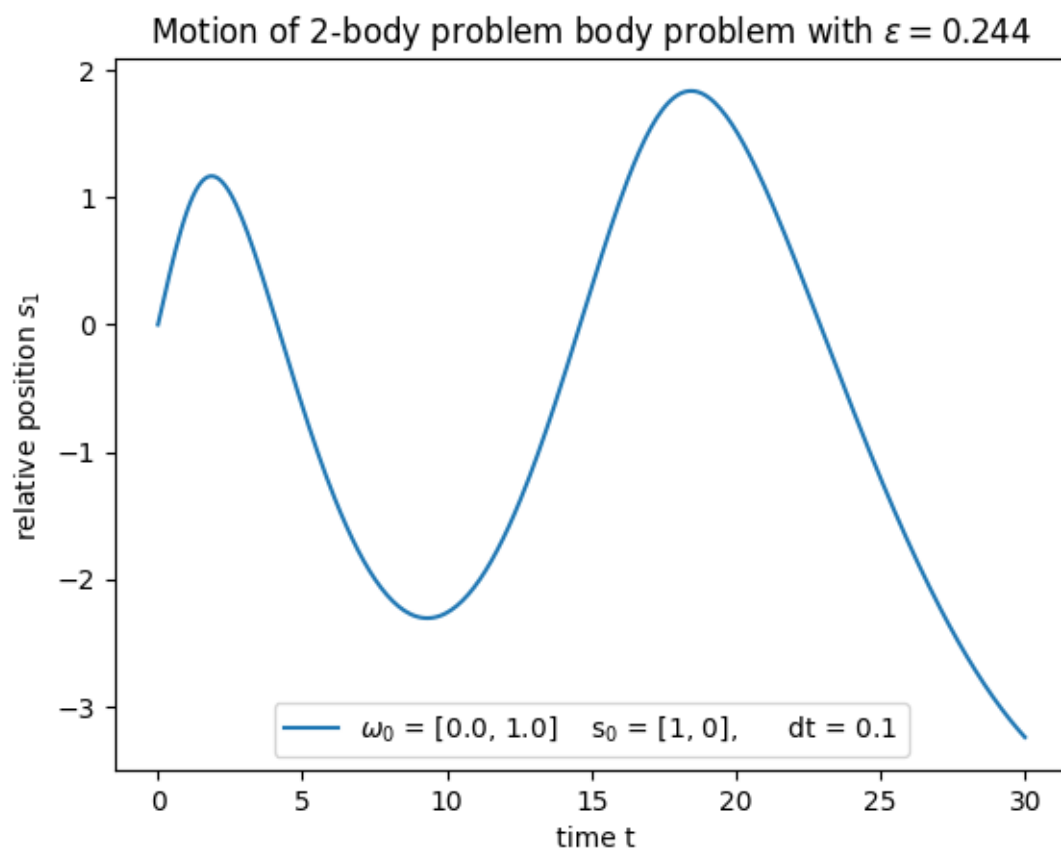
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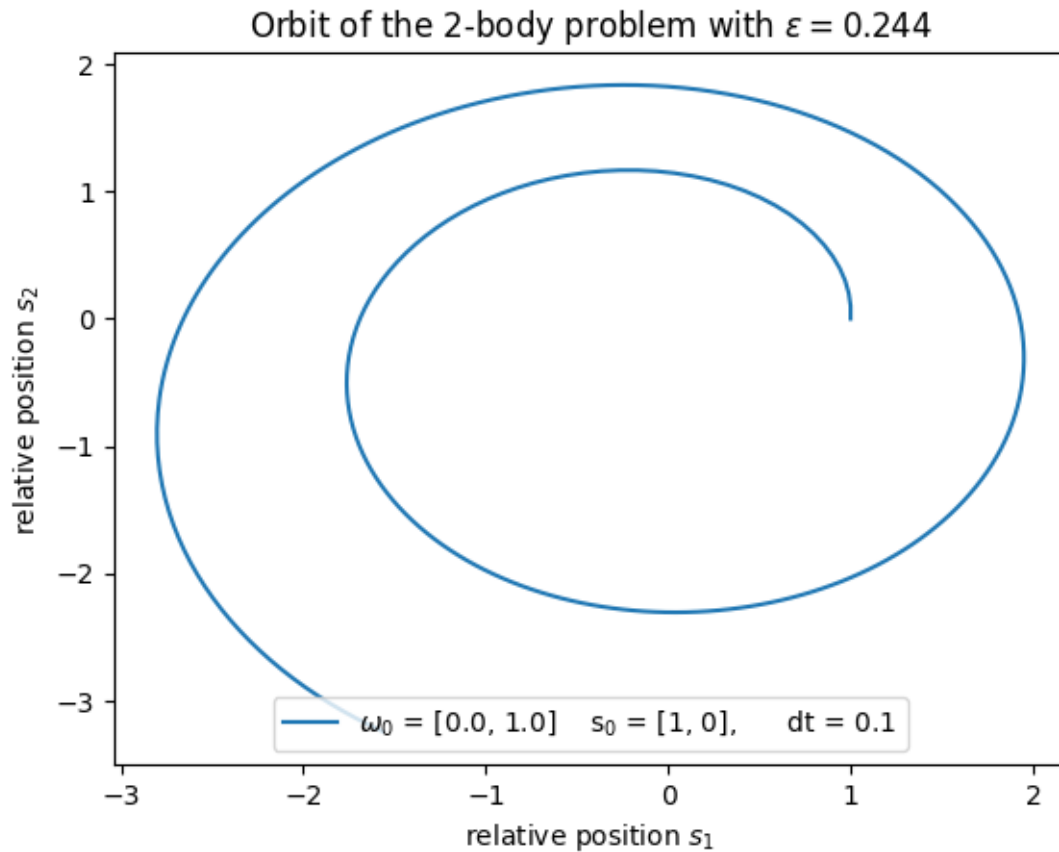
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```
[ ]: euler([0,1],[1,0], 0.01,30)
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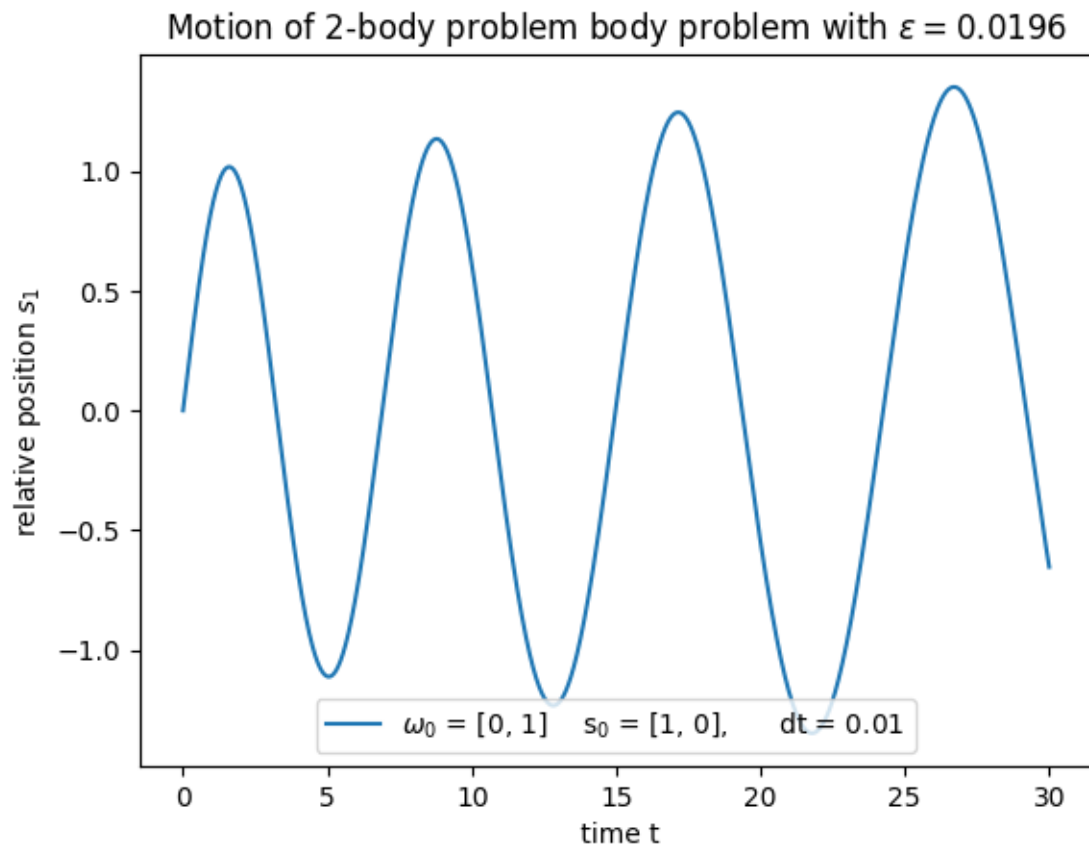
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2.7516943302449928
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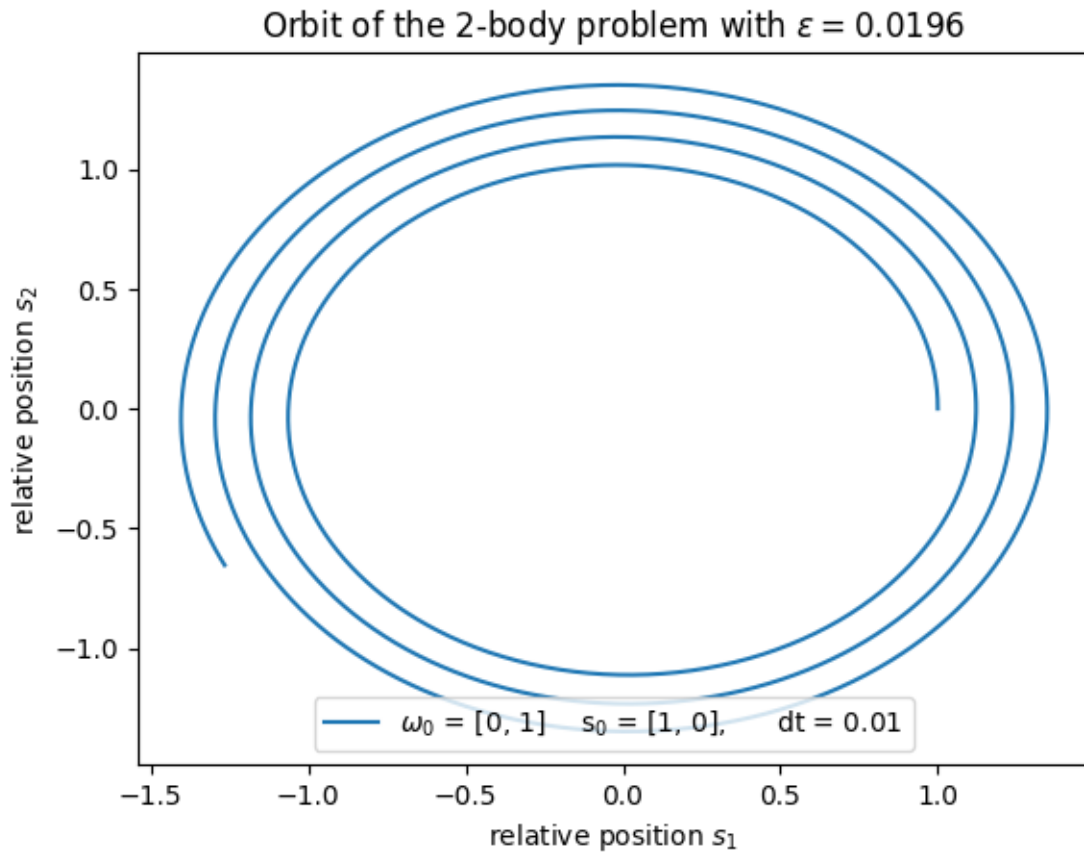
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epsilon calculated geometrically :0.203
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```

[0.36307115, -0.75120722]]))





```
[ ]: euler([0,1],[1,0], 0.001,30)
```

```
1.5000003749539096e-06
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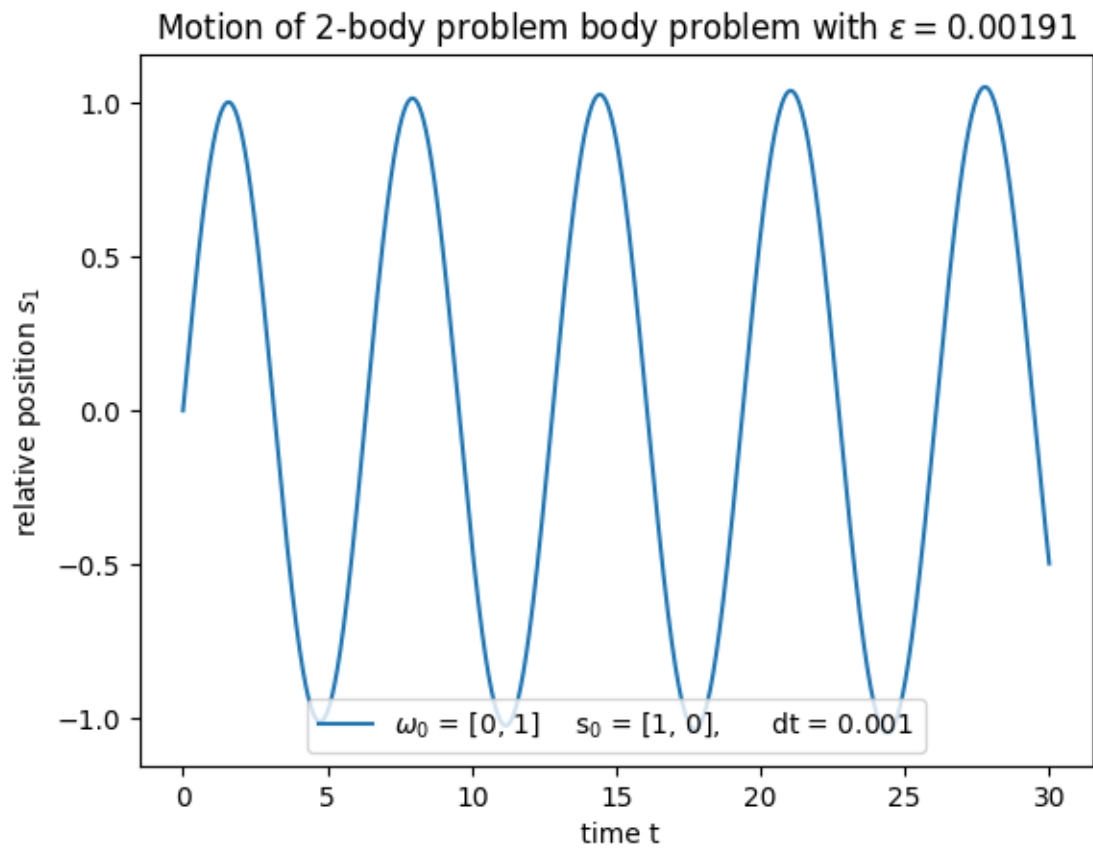
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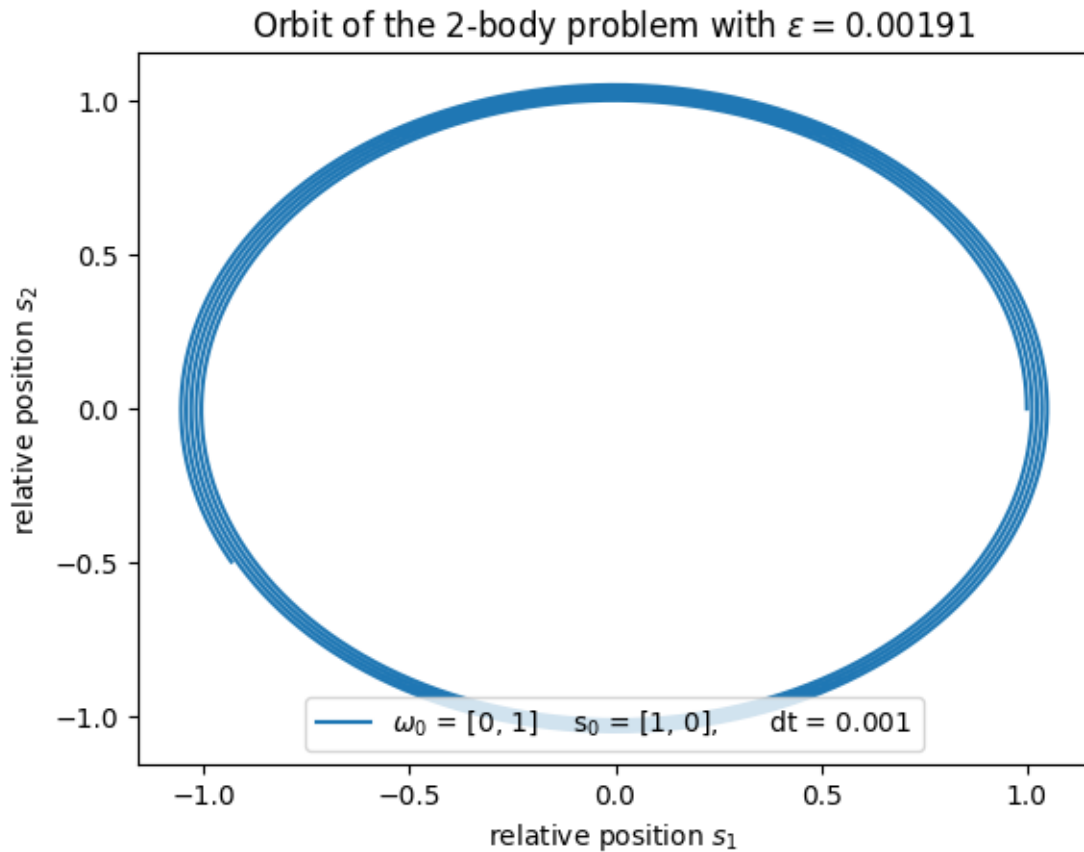
```
2.0993082373604404
```

```
epsilon calculated geometrically :0.0766
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            [-0.002        ,  0.999999    ],
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```


[0.45441717, -0.85931387]]))





```
[ ]: euler([0,1],[1,0], 0.0001,30)
```

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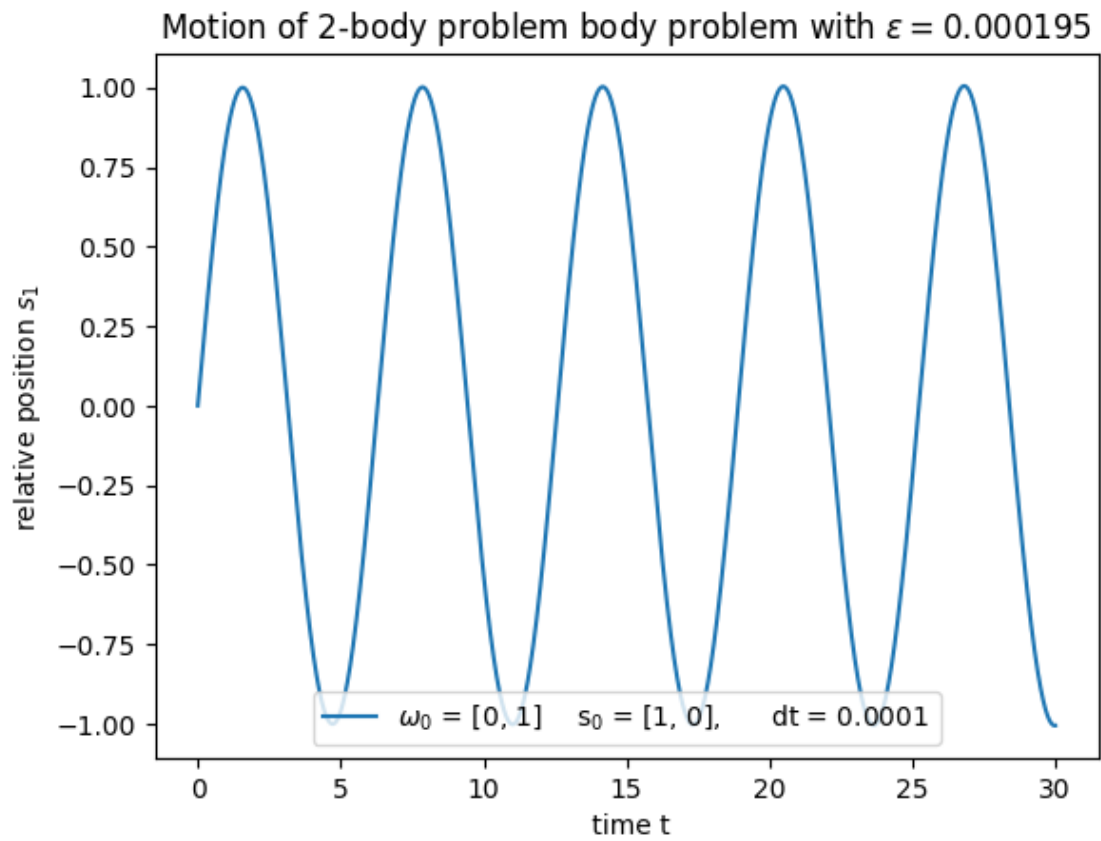
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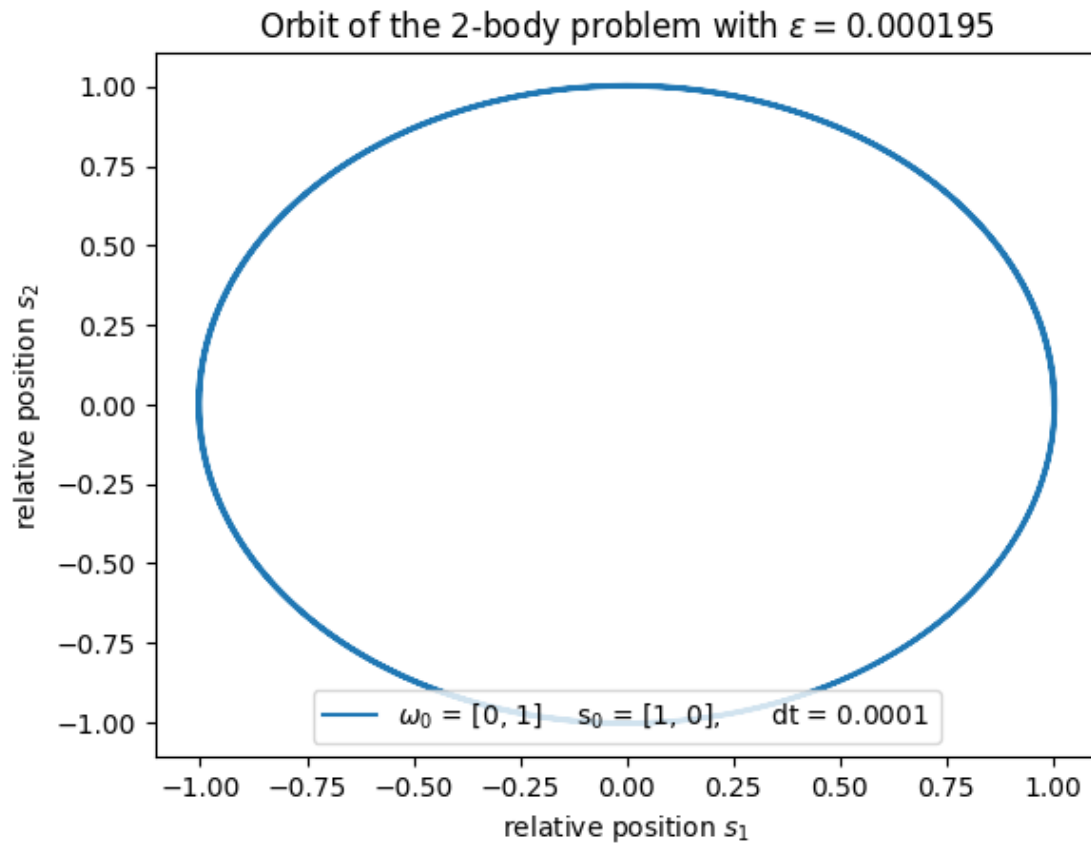
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epsilon calculated geometrically :0.025
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[9.96658677e-01, 2.10453633e-02]]))





```
[ ]: euler([0,1/np.sqrt(2)], [1,0], 0.1,10)
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```
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```
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```
epsilon calculated geometrically :0.898
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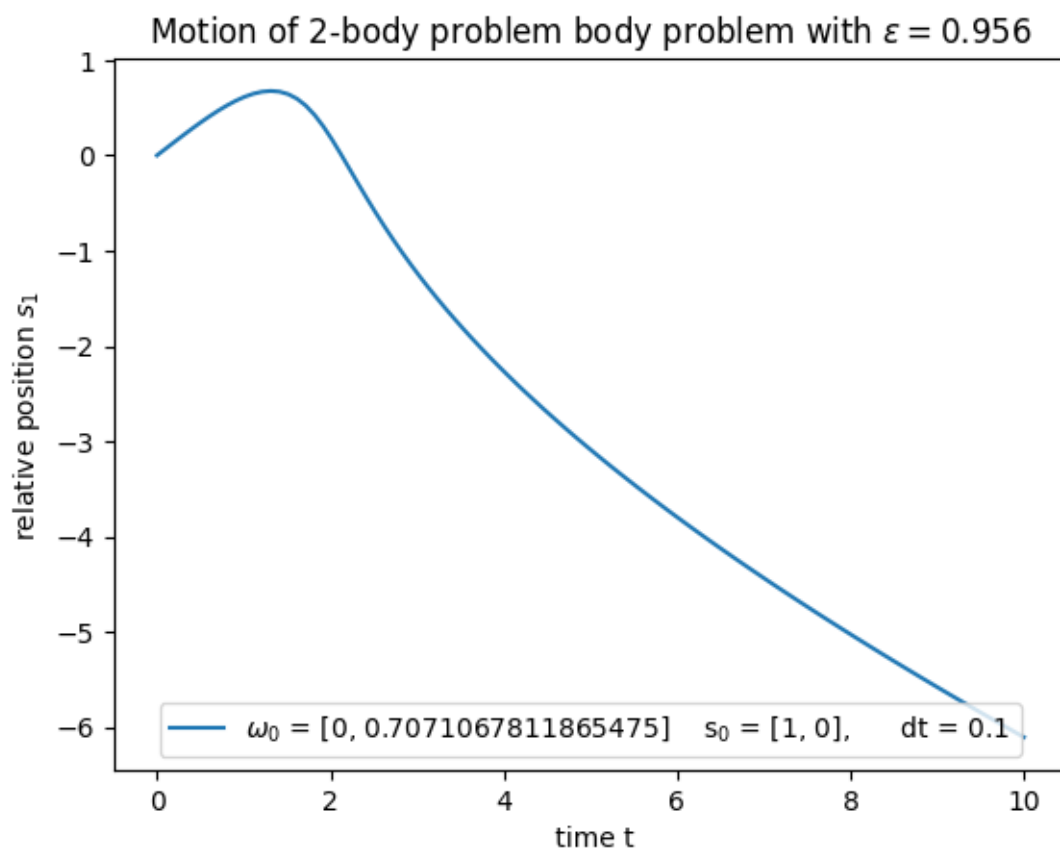
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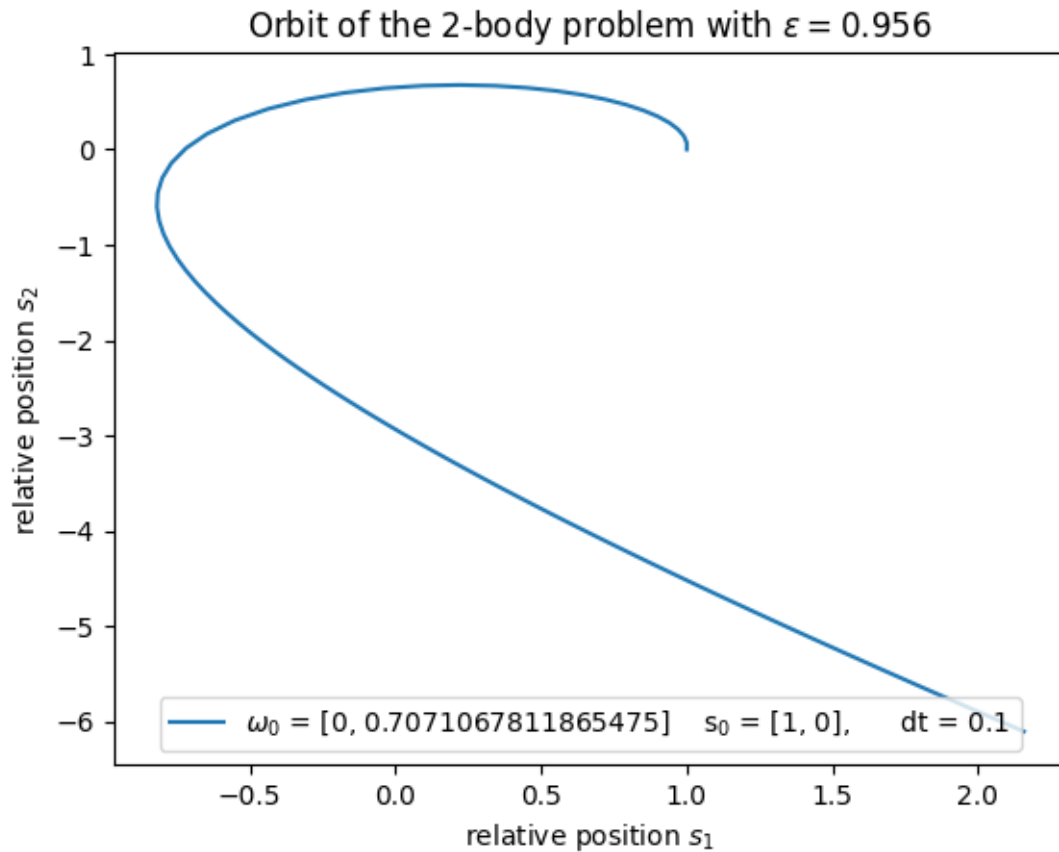
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[ 0.40973575, -0.56132895],
[ 0.40880389, -0.55810061],
[ 0.40788038, -0.5549522 ],
[ 0.40696534, -0.55188058],
[ 0.40605884, -0.5488828 ],
[ 0.40516096, -0.54595604],
[ 0.40427175, -0.54309764],
[ 0.40339122, -0.54030507],
[ 0.40251939, -0.53757593],
[ 0.40165626, -0.53490794],
[ 0.40080181, -0.53229893],
[ 0.39995602, -0.52974683],
[ 0.39911884, -0.52724967],
[ 0.39829023, -0.52480557],
[ 0.39747014, -0.52241274],
[ 0.39665851, -0.52006948],
[ 0.39585527, -0.51777416]]))

```



```
[ ]: euler([0,1/np.sqrt(2)], [1,0], 0.01,10)
```

```
0.4999250009382423
```

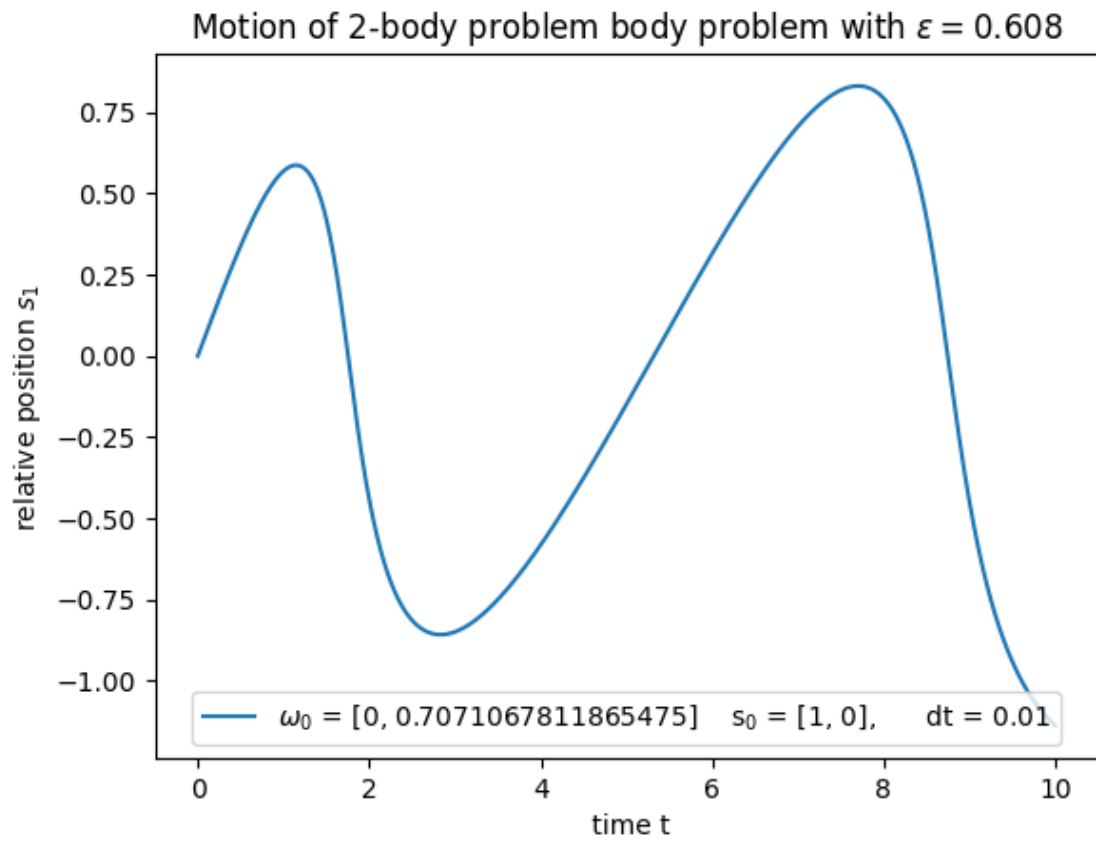
```
2.1810763825379444
```

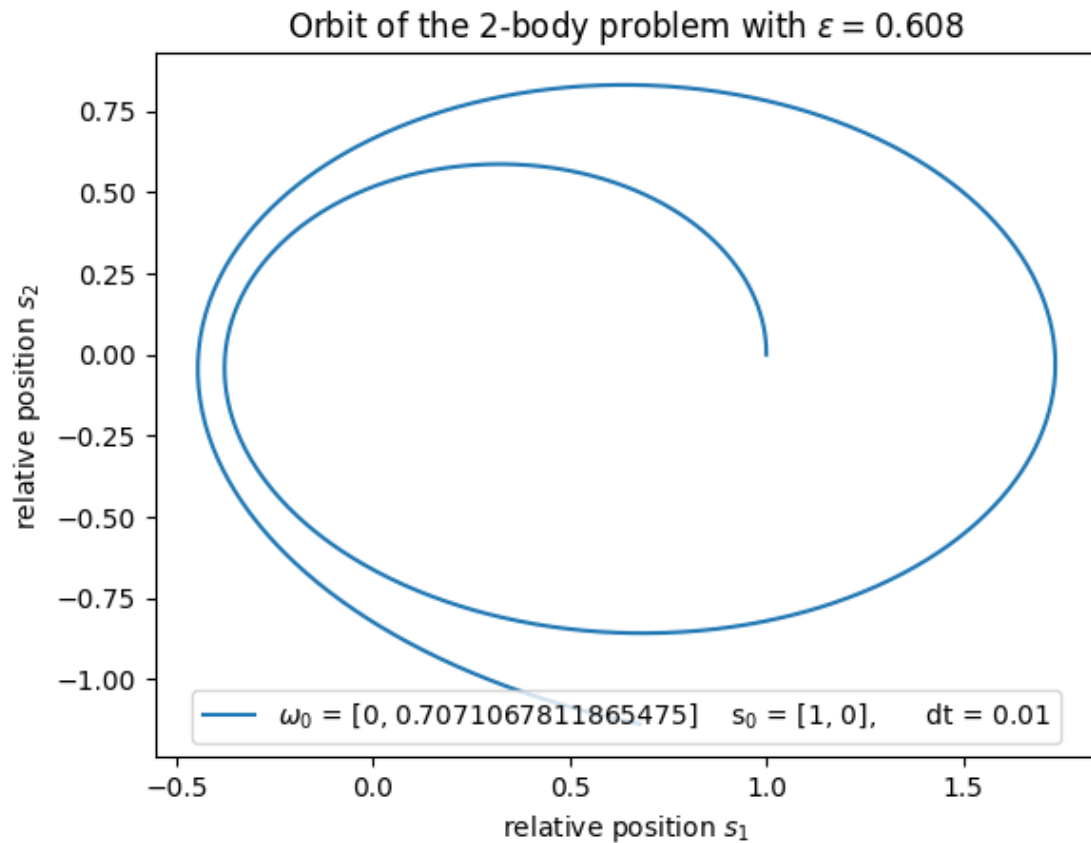
```
1.9697271277033543
```

```
epsilon calculated geometrically :0.476
```

```
[ ]: (array([[ 1.         ,  0.         ],
            [ 1.         ,  0.00707107],
            [ 0.9999      ,  0.01414214],
            ...,
            [ 0.66002372, -1.13378841],
            [ 0.66931768, -1.13618733],
            [ 0.67858241, -1.13853604]]),
      array([[ 0.         ,  0.70710678],
            [-0.01        ,  0.70710678],
            [-0.01999925,  0.70703608],
            ...,
            [ 0.92939642, -0.23989199],
            [ 0.92647329, -0.23487064],
```

[0.9235544 , -0.22991574]]))





```
[ ]: euler([0,1/np.sqrt(2)], [1,0], 3.8,10)
```

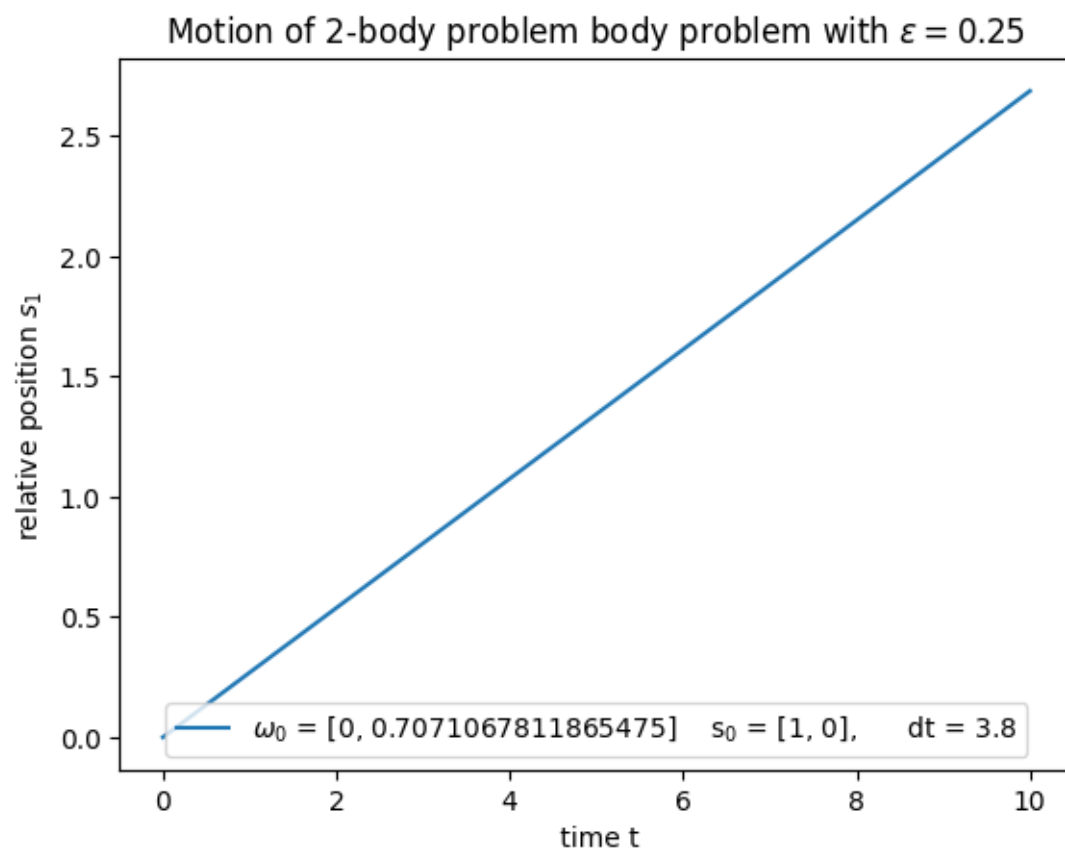
```
0.0
```

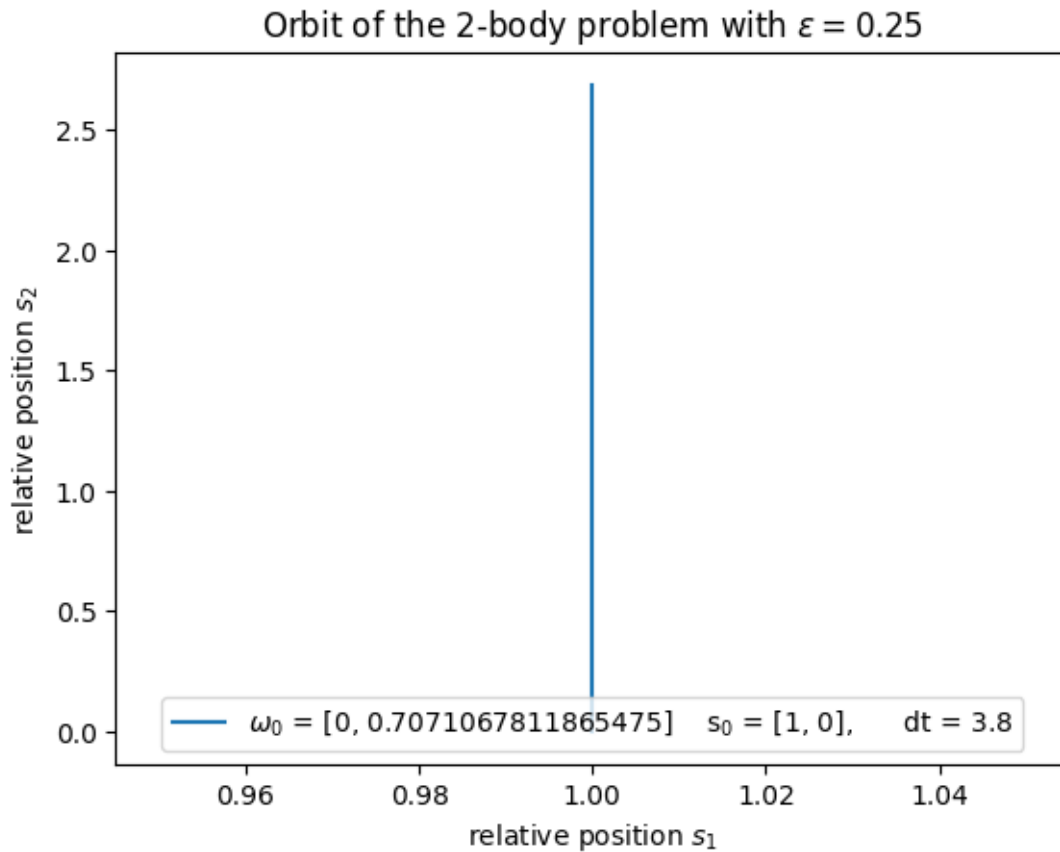
```
0.0
```

```
2.68700576850888
```

```
epsilon calculated geometrically :1.0
```

```
[ ]: (array([[1.      , 0.      ],
             [1.      , 2.68700577]]),
      array([[0.      , 0.70710678],
             [-3.8     , 0.70710678]]))
```





```
[ ]: euler([0,1/np.sqrt(2)], [1,0], 0.0001,10)
```

```
0.49999999250000005
```

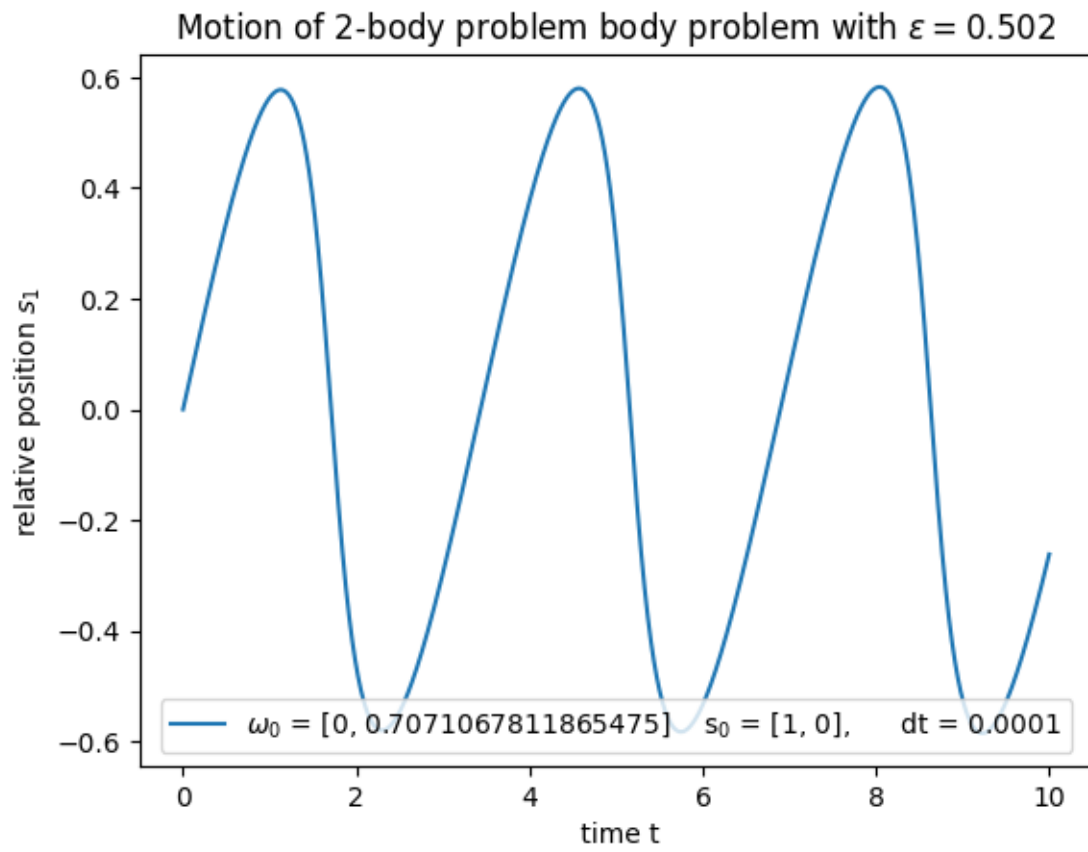
```
1.3480258576967694
```

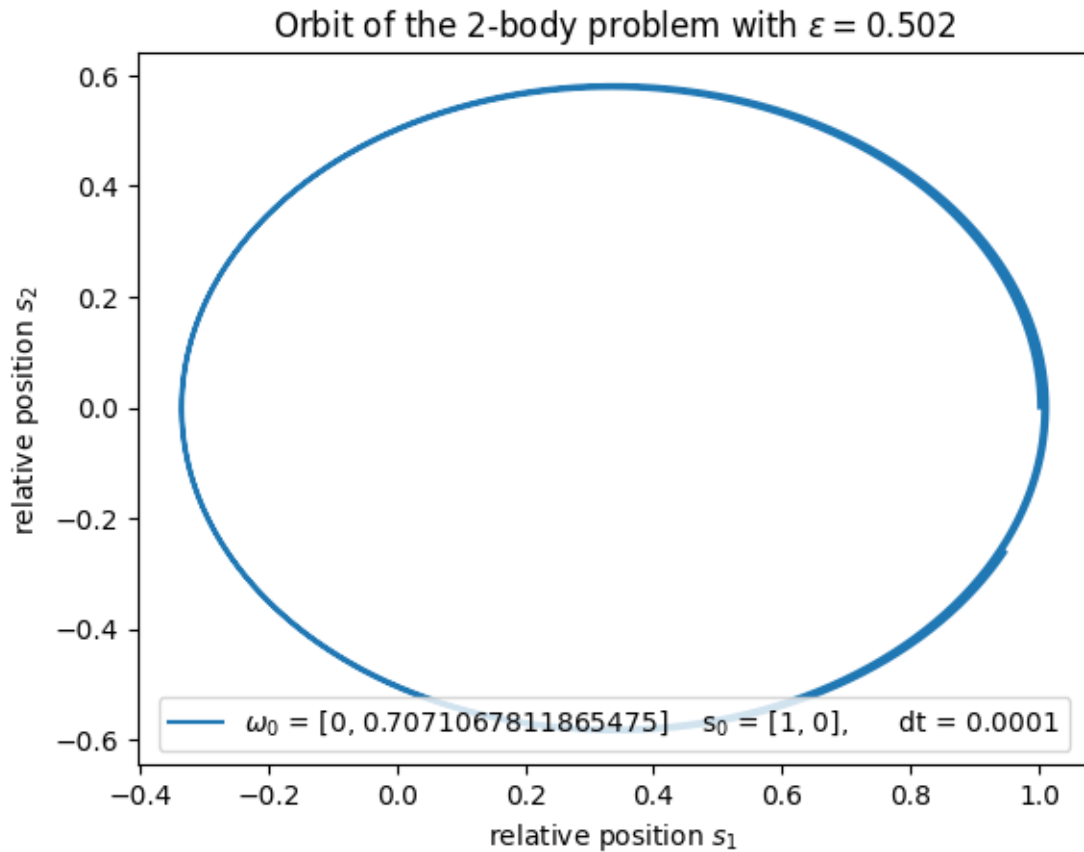
```
1.1675373818253267
```

```
epsilon calculated geometrically :0.577
```

```
[ ]: (array([[ 1.00000000e+00,  0.00000000e+00],
             [ 1.00000000e+00,  7.07106781e-05],
             [ 9.99999990e-01,  1.41421356e-04],
             ...,
             [ 9.47099038e-01, -2.61851117e-01],
             [ 9.47136525e-01, -2.61786428e-01],
             [ 9.47174002e-01, -2.61721736e-01]]),
      array([[ 0.00000000e+00,  7.07106781e-01],
             [-1.00000000e-04,  7.07106781e-01],
             [-1.99999999e-04,  7.07106774e-01],
             ...,
             [ 3.74868772e-01,  6.46891334e-01],
             [ 3.74768950e-01,  6.46918932e-01],
```

[3.74669130e-01, 6.46946522e-01]]))





```
[ ]: #computing the energie of the motion and plotting it
```

```
def euler_0(w0,s0,h,e):
    #the function will take the initial velocity
    # initial position, the whished step size
    # and the desired upper bondary for the integral

    G = 1 #grativational constant
    m = 1 #mass of both bodies
    N = int(e/h) #time steps

    s = np.zeros((N,2))
    w = np.zeros((N,2))
    LRL = np.zeros((N,2))
    s[0] = s0
    w[0] = w0
    #leap-frog algorithm
    for i in range(N-1):
        s[i+1] = s[i] + h*w[i]
```



```

        w[i+1] = w[i]-h*s[i]/np.linalg.norm(s[i])**3
    return w

def energy_em(w0,s0,dt,e):
    # dt should be a list with initila and final ime step for which the
    ↪ difference of energy has
    # to be computed
    # w0 is the initial velocity, s0, initila position
    # and e the time span in which the motion will take place

    dt = np.arange(dt[0], dt[-1], 0.1) #list of dt to be consider
    E = np.array([])

    #computiong the error in the energy
    for i in dt:
        w = euler_0(w0,s0,i,e)
        e_i = 1/2*np.dot(w[0],w[0]) #energy at beginning of motion
        e_f = 1/2*np.dot(w[-1],w[-1]) #energy at end of motion
        E = np.append(E, abs(e_f-e_i))

    #plotting error as funktion of dt
    plt.loglog(dt,E)
    plt.title(r'Error of energy as fucntion of time step, Euler Methode')
    plt.xlabel(r'time step $dt$')
    plt.ylabel(r'Error of energy $\Delta E$')
    print(dt)
    print()
    print(E)

```

```
[ ]: energy_em([0,1/np.sqrt(2)], [1,0], [0.0001,6],10)
```

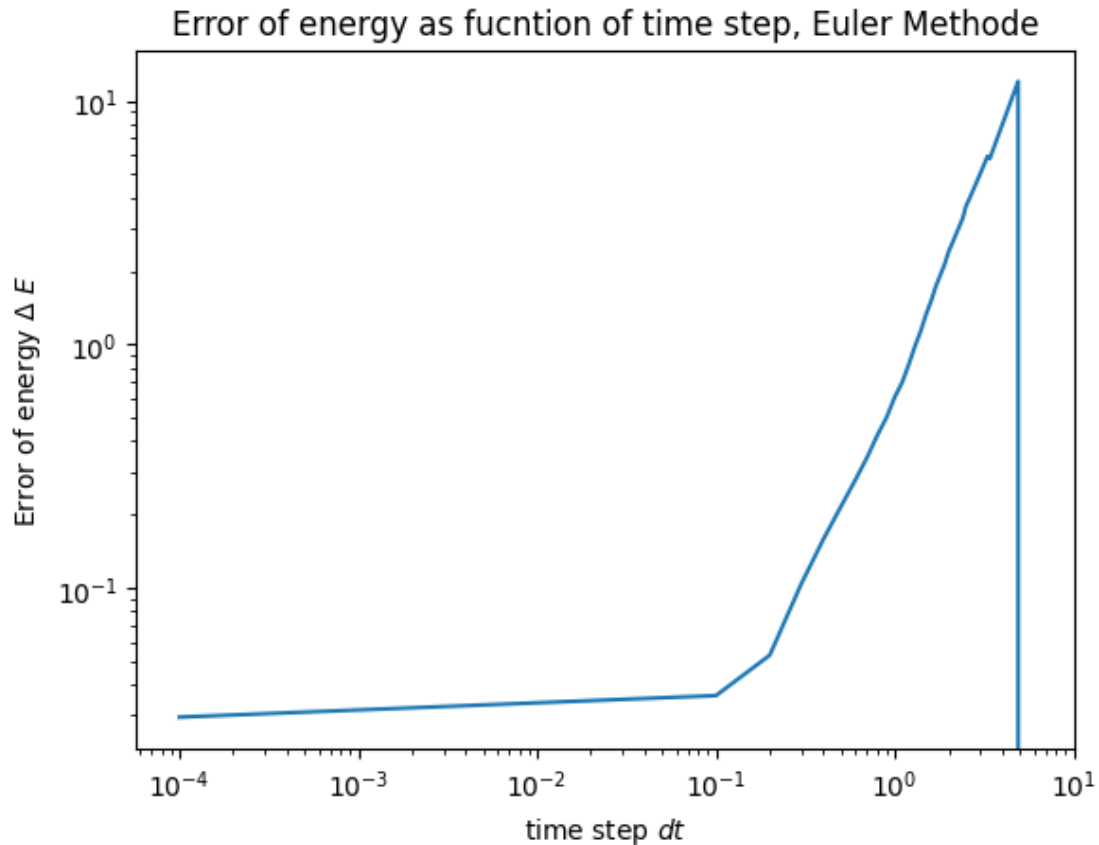
```

1.0000e-04 1.0010e-01 2.0010e-01 3.0010e-01 4.0010e-01 5.0010e-01
6.0010e-01 7.0010e-01 8.0010e-01 9.0010e-01 1.0001e+00 1.1001e+00
1.2001e+00 1.3001e+00 1.4001e+00 1.5001e+00 1.6001e+00 1.7001e+00
1.8001e+00 1.9001e+00 2.0001e+00 2.1001e+00 2.2001e+00 2.3001e+00
2.4001e+00 2.5001e+00 2.6001e+00 2.7001e+00 2.8001e+00 2.9001e+00
3.0001e+00 3.1001e+00 3.2001e+00 3.3001e+00 3.4001e+00 3.5001e+00
3.6001e+00 3.7001e+00 3.8001e+00 3.9001e+00 4.0001e+00 4.1001e+00
4.2001e+00 4.3001e+00 4.4001e+00 4.5001e+00 4.6001e+00 4.7001e+00
4.8001e+00 4.9001e+00 5.0001e+00 5.1001e+00 5.2001e+00 5.3001e+00
5.4001e+00 5.5001e+00 5.6001e+00 5.7001e+00 5.8001e+00 5.9001e+00]

[ 0.02945838  0.03610472  0.0529069   0.10374135  0.15885144  0.2156045
 0.27591036  0.34284985  0.42363654  0.50068126  0.60675802  0.69698256
 0.82866947  0.98541209  1.13792351  1.33209797  1.50990621  1.73431401
 1.92695346  2.12763023  2.39696296  2.60742929  2.82609298  3.05307877
 3.28856534  3.68189761  3.92516049  4.17829417  4.4414105   4.71459728
 4.99792264  5.29143866  5.59518446  5.90918866  5.78034001  6.12535001]

```

6.48036001	6.84537001	7.22038001	7.60539001	8.0004	8.40541001
8.82042	9.24543	9.68044001	10.12545	10.58046001	11.04547001
11.52048001	12.00549001	0.	0.	0.	0.
0.	0.	0.	0.	0.	0.



The results are as expected till $dt > 4$, at this point the simulation is so erroneous, that the result of the Euler Method is a linear trajectory, for which the Energy is constant. A constant energy is the ideal case, as it is visible for $dt \rightarrow 0$, however if the trajectory is erroneously integrated, this result can't be taken as a correct one.

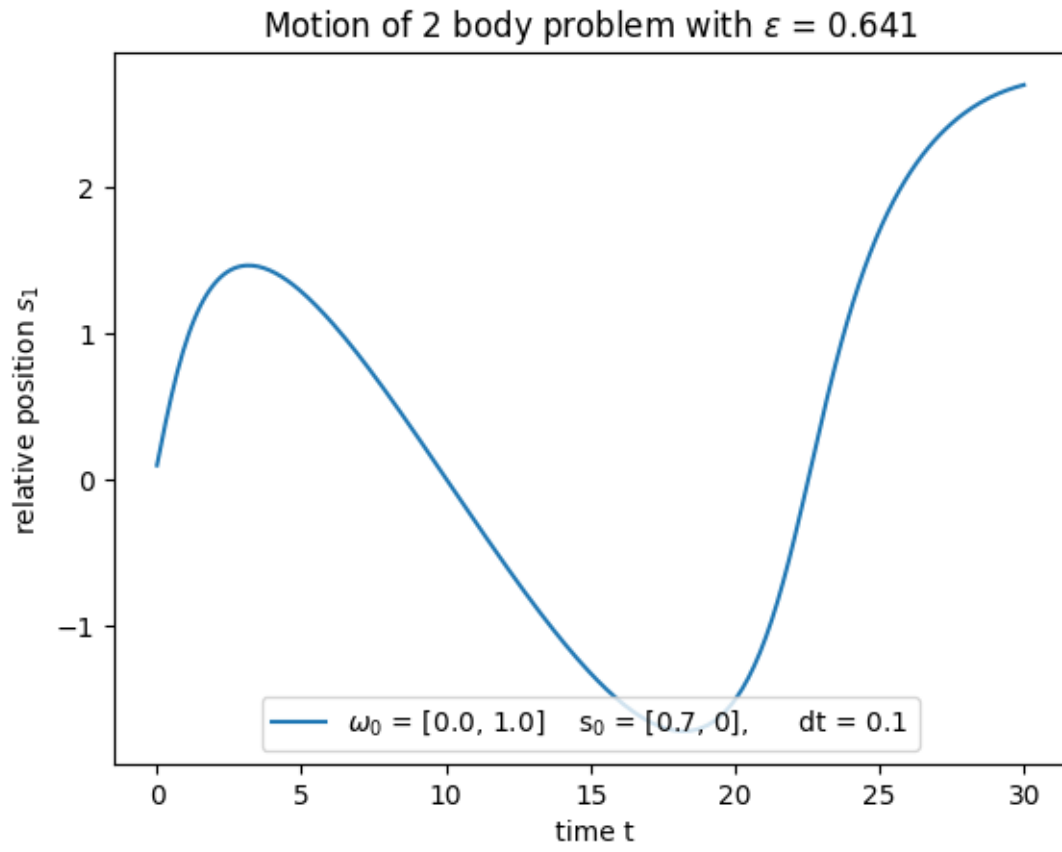
3.2 b) leap-frog algorithm

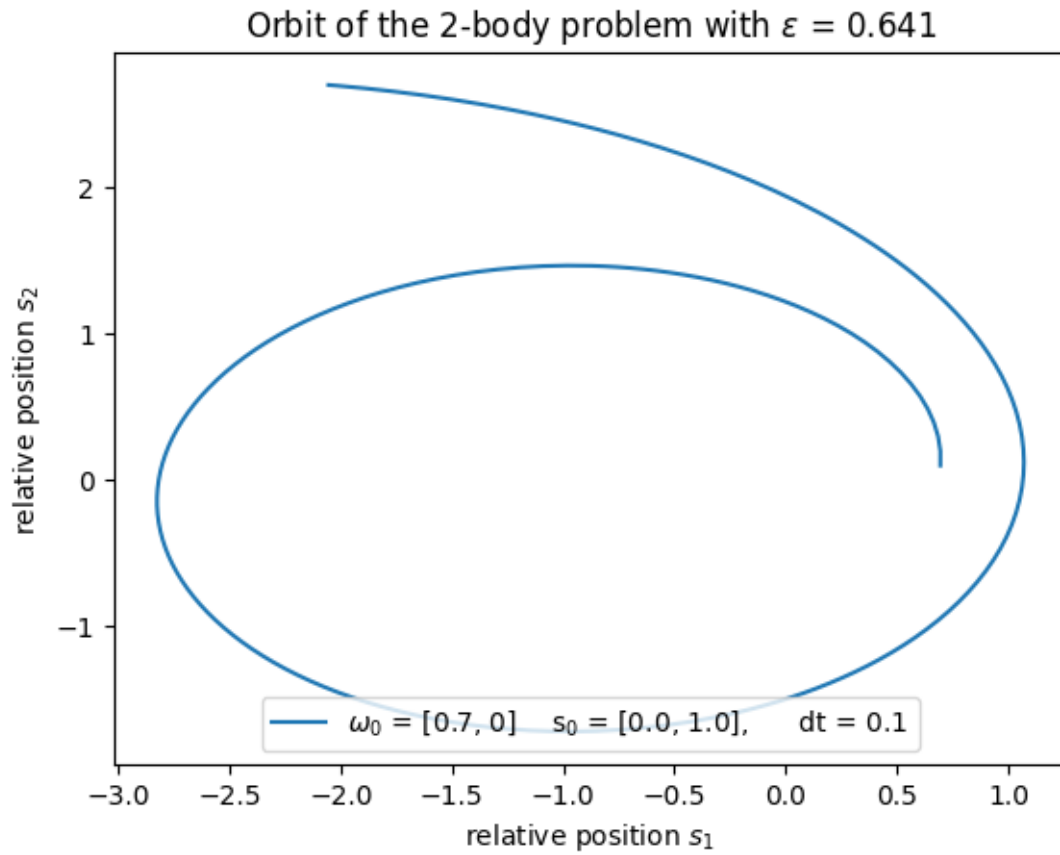
```
[ ]: leap_frog([0.0,1.0],[0.7,0], 0.1,30)
```

```
3.895588310504747
```

```
4.416367359310763
```

```
epsilon calculated geometrically :0.534
```



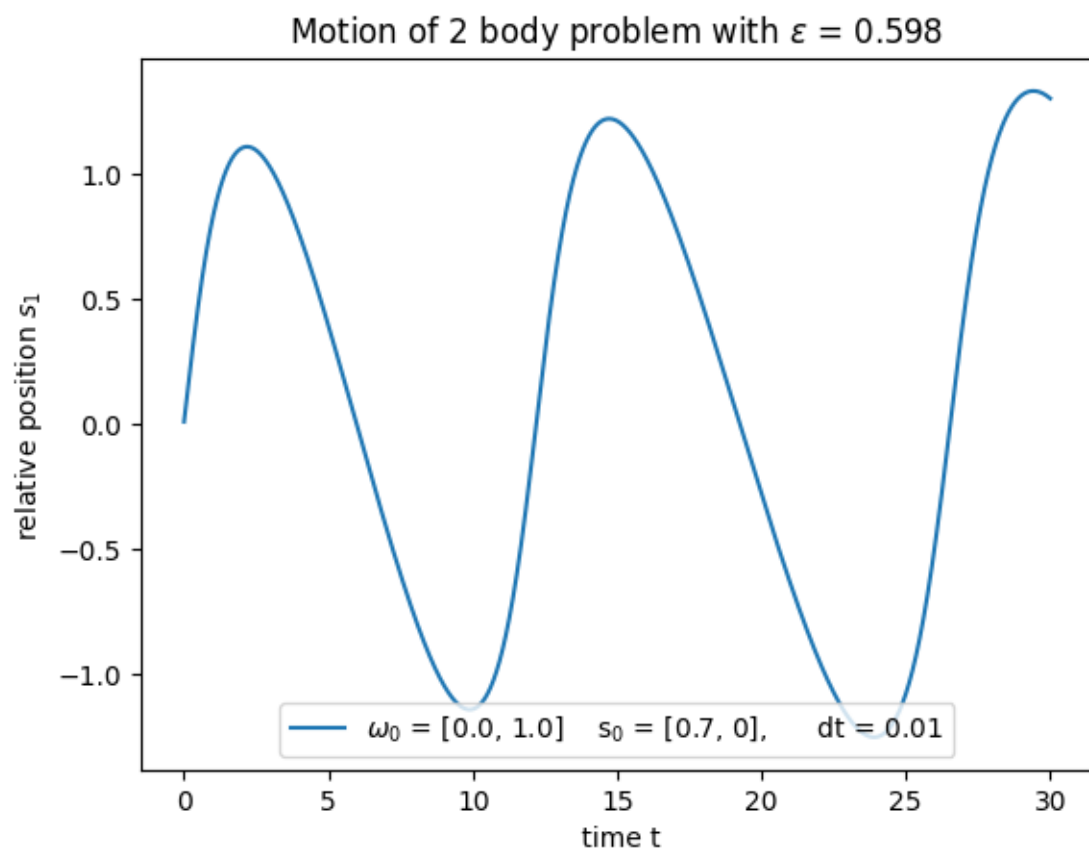


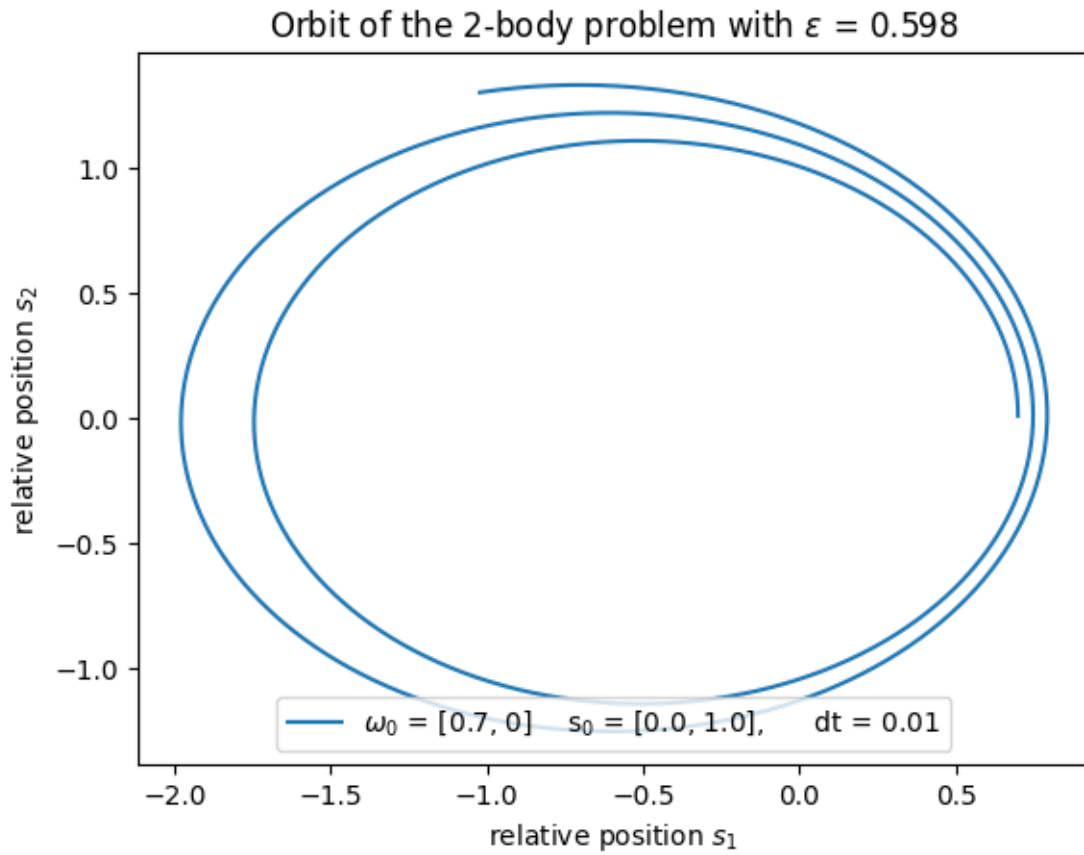
```
[ ]: leap_frog([0.0,1.0],[0.7,0], 0.01,30)
```

2.773665244342849

2.5827428157153465

epsilon calculated geometrically :0.365



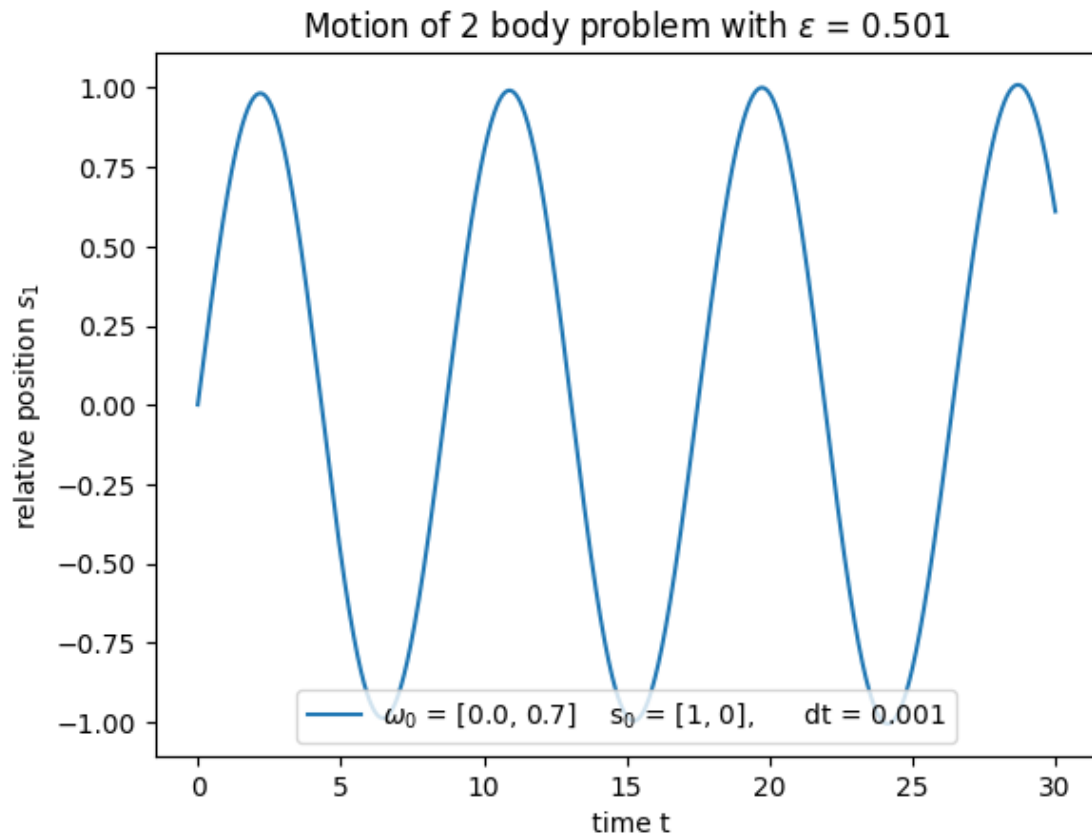


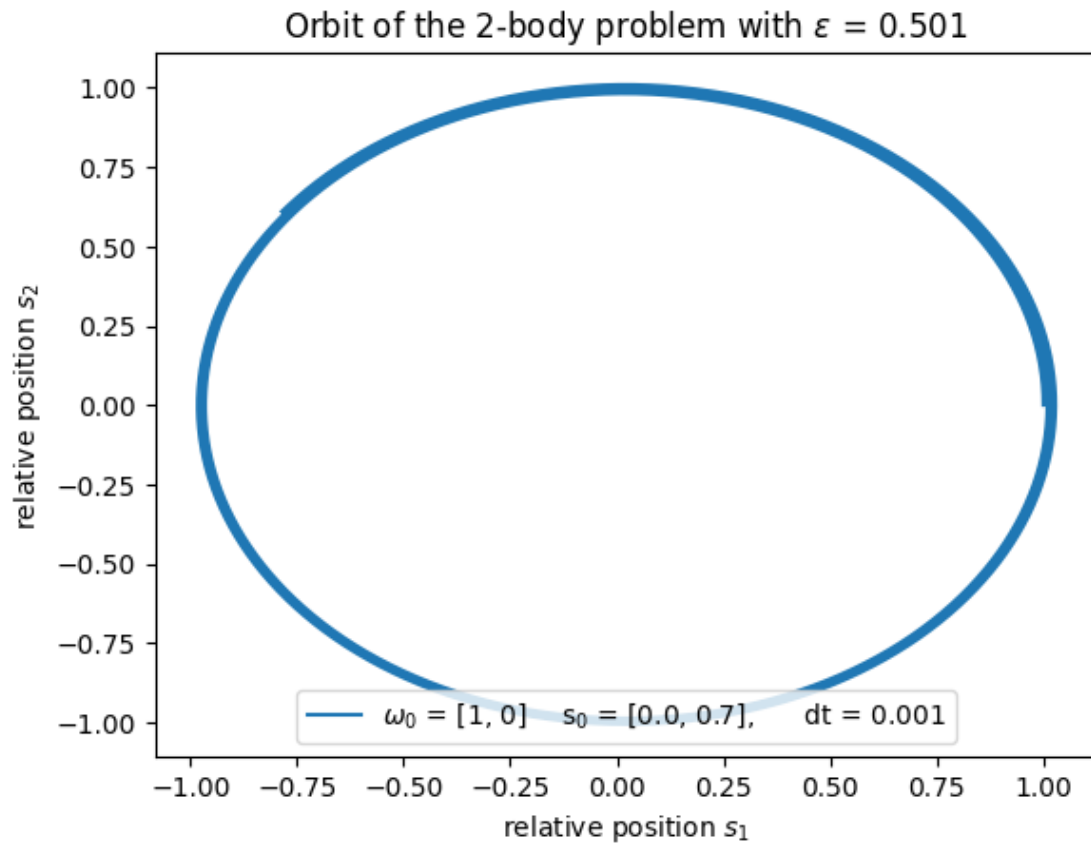
```
[ ]: leap_frog([0.0,0.7],[1,0], 0.001,30)
```

```
2.010055015594763
```

```
2.0139096011602406
```

```
episilon calculated geometrically :0.062
```



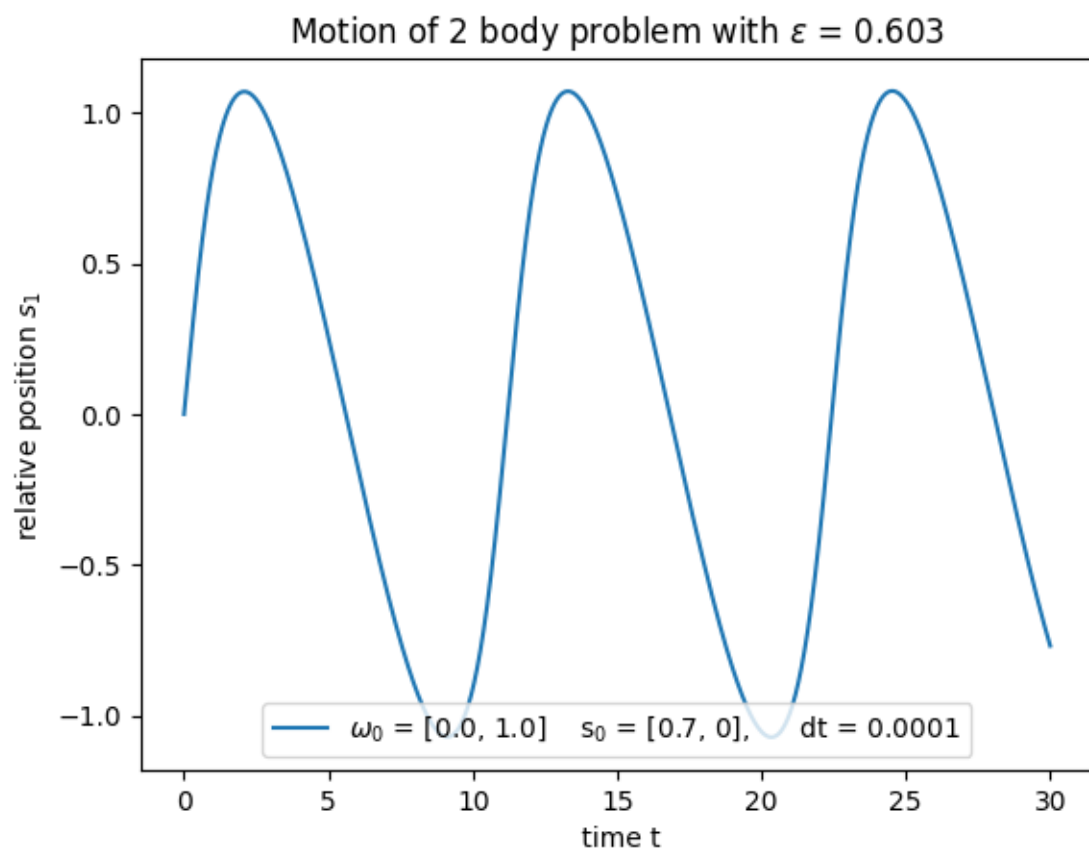


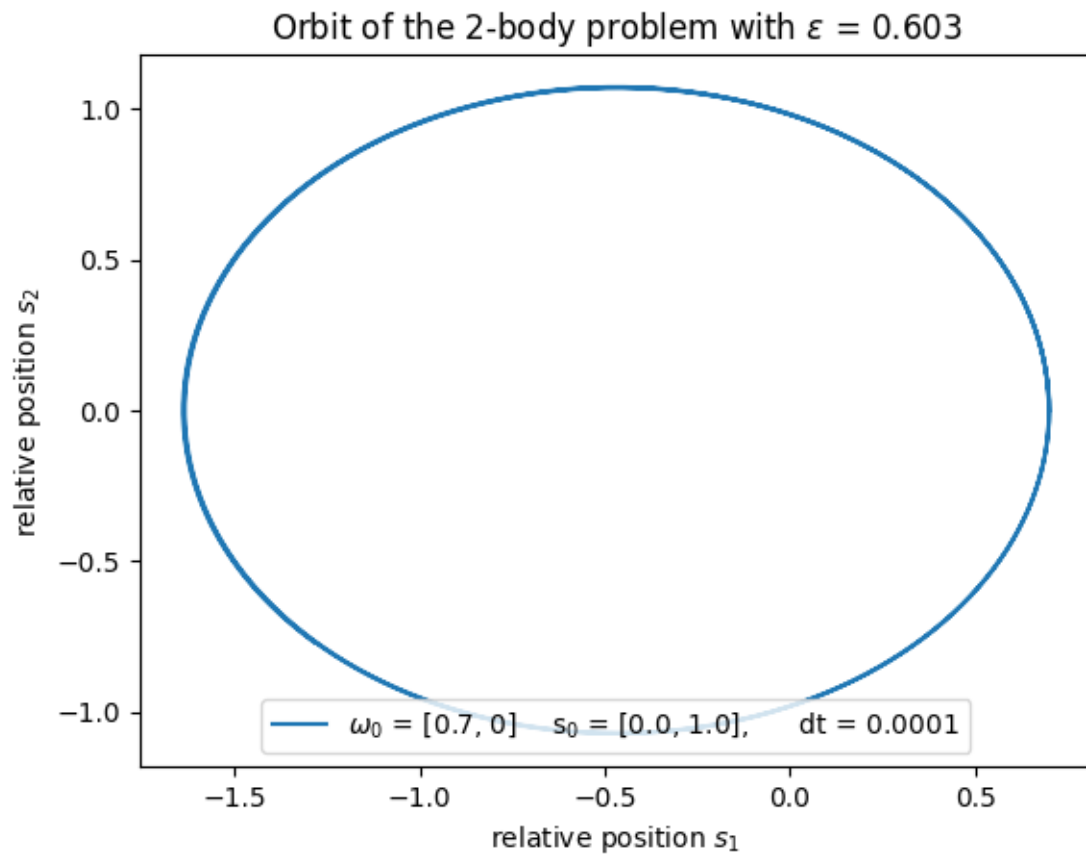
```
[ ]: leap_frog([0.0,1.0],[0.7,0], 0.0001,30)
```

```
2.3399526917437283
```

```
2.1430118545307777
```

```
epsilon calculated geometrically :0.402
```

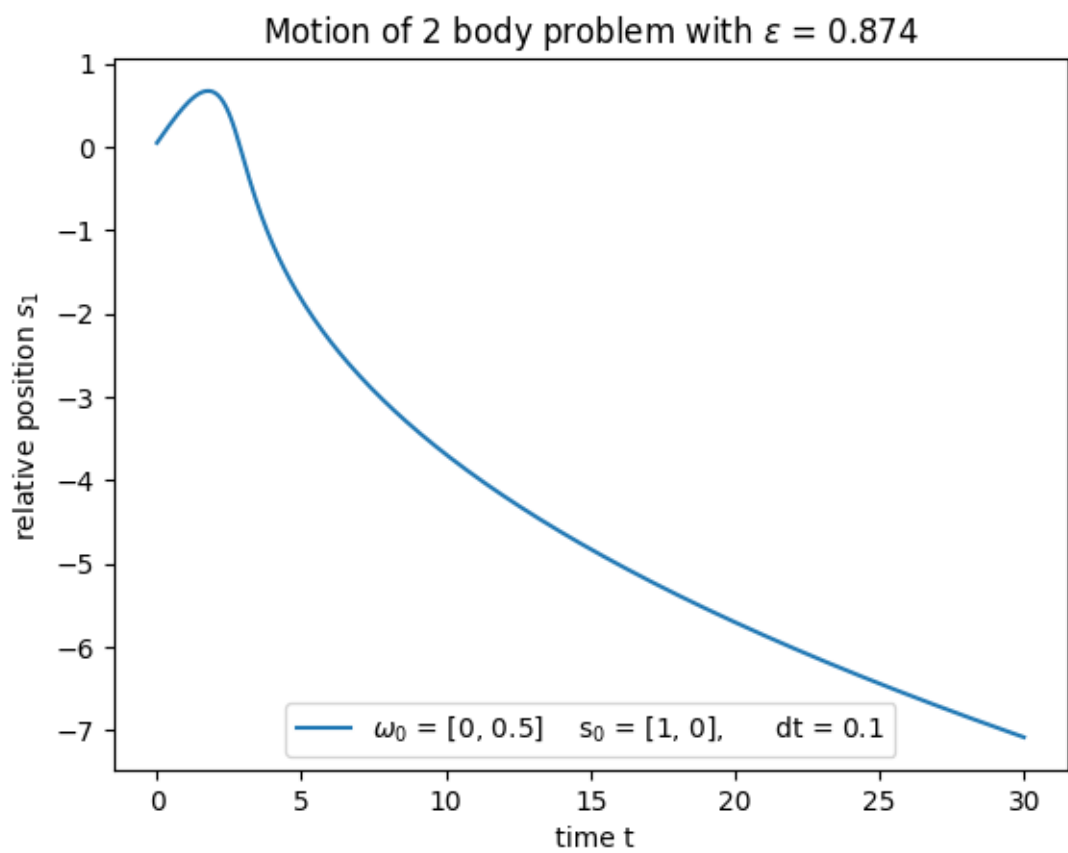


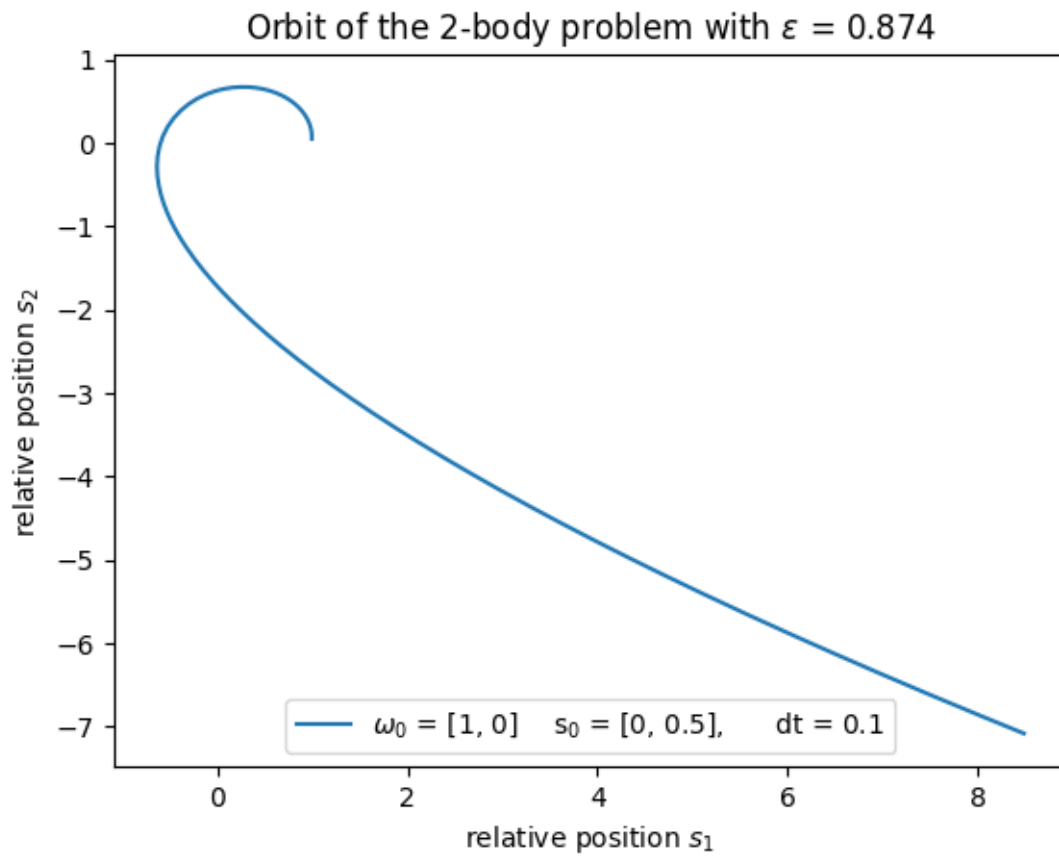
```
[ ]: leap_frog([0,1/2],[1,0], 0.1,30)
```

```
9.115157067317028
```

```
7.758732521682457
```

```
epsilon calculated geometrically :0.525
```



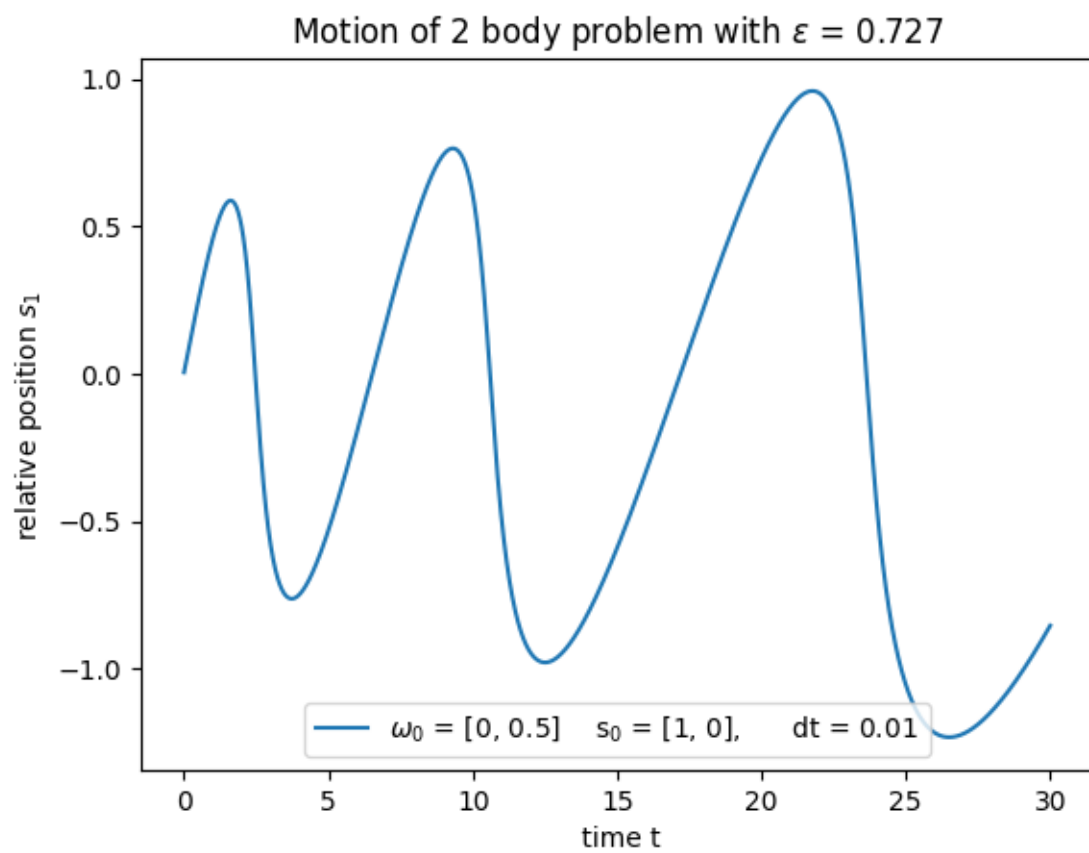


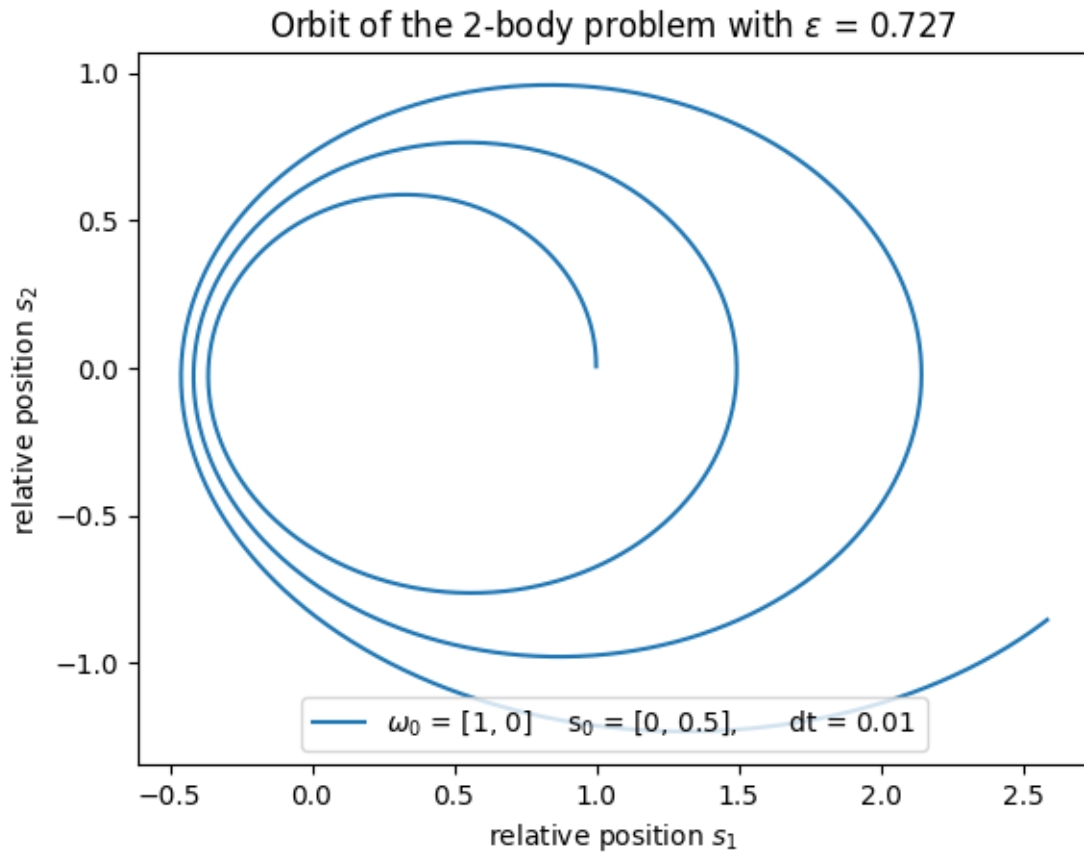
```
[ ]: leap_frog([0,1/2],[1,0], 0.01,30)
```

```
3.0464735953711406
```

```
2.191182322829405
```

```
epsilon calculated geometrically :0.695
```



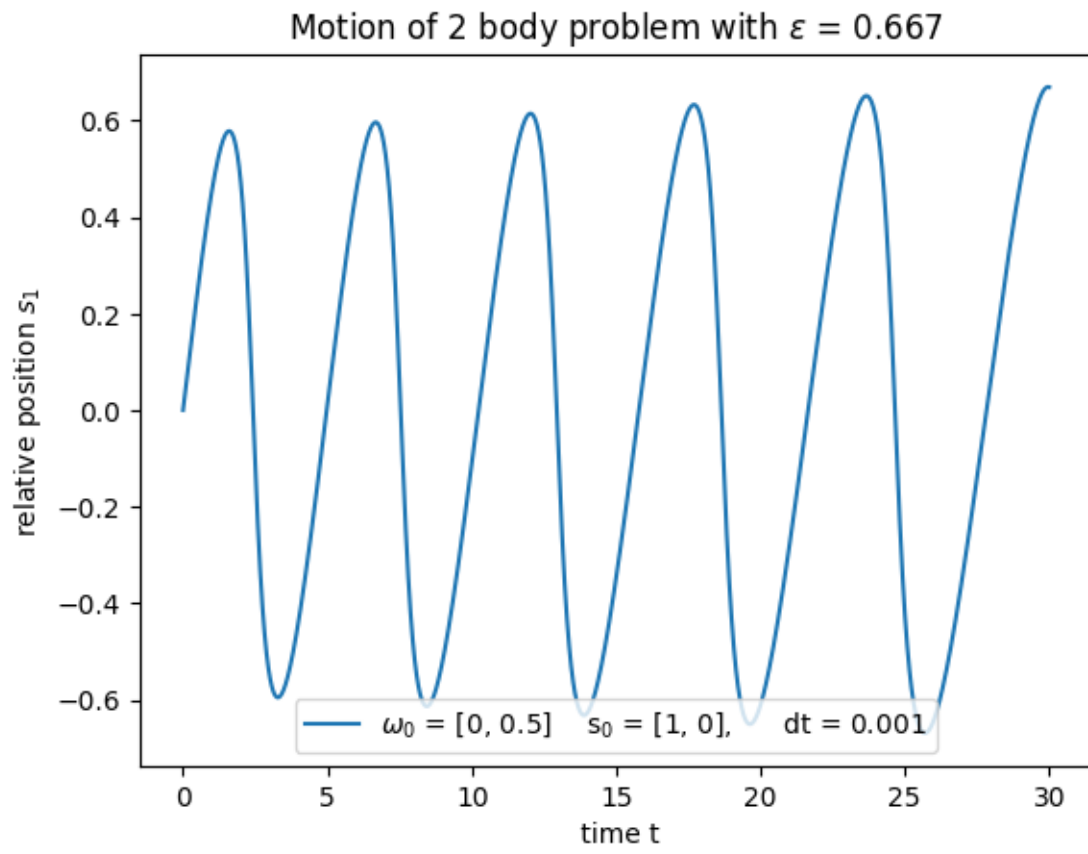


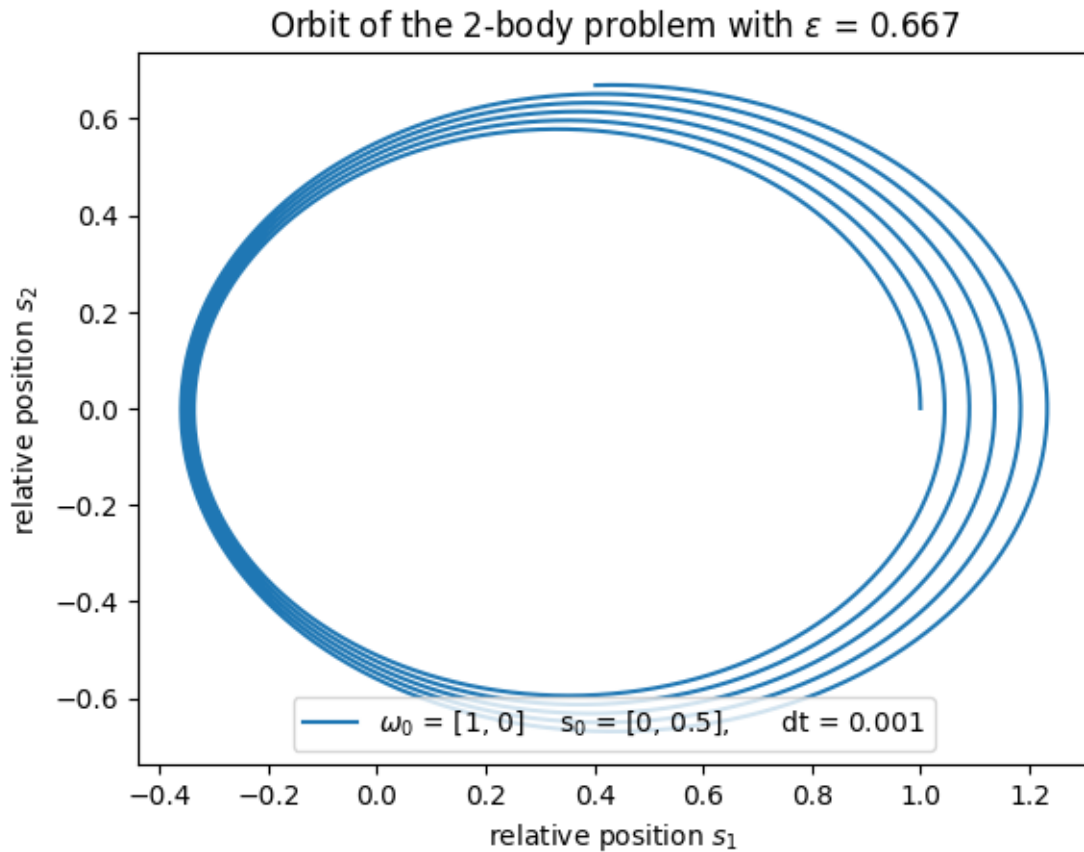
```
[ ]: leap_frog([0,1/2],[1,0], 0.001,30)
```

```
1.5934798580931757
```

```
1.3383089210857317
```

```
epsilon calculated geometrically :0.543
```





```
[ ]: leap_frog([0,1/2],[1,0], 0.0001,30)
```

```
[ ]: def leap_frog_0(w0,s0,h,e):
    #the function will take the initial velocity
    # initial position, the whished step size
    # and the desired upper bondary for the integral
    G = 1 #grativational constant
    m = 1 #mass of both bodies
    N = int(e/h) #time steps

    s = np.zeros((N,2))
    w = np.zeros((N,2))
    LRL = np.zeros((N,2))

    w[0] = w0
    s[0] = s0

    #First the programm will compute the s_1/2 step using the
    #euler forward methode
```



```

a = 0.5*h/np.linalg.norm(s0)**3
wn = w[0] - h/2*s[0]/np.linalg.norm(s[0])
s[0] = s0 + h*wn

#leap-frog algorithm
for i in range(N-1):
    w[i + 1] = w[i]-h/2*s[i]/np.linalg.norm(s[i])**3
    s[i + 1] = s[i] + w[i]*h
return w

def energy_lf(w0,s0,dt,e):
    # dt should be a list with initila and final ime step for which the
    ↪ difference of energy has
    # to be computed
    # w0 is the initial velocity, s0, initila position
    # and e the time span in which the motion will take place

    dt = np.arange(dt[0], dt[-1], 0.1) #list of dt to be consider
    E = np.array([])
    #computiong the error in the energy
    for i in dt:
        w = leap_frog_0(w0,s0,i,e)
        print(w)
        print()
        e_i = 1/2*np.dot(w[0],w[0]) #energy at beginning of motion
        e_f = 1/2*np.dot(w[-1],w[-1]) #energy at end of motion
        E = np.append(E, abs(e_f-e_i))

    #plotting error as funktion of dt
    plt.loglog(dt,E)
    print(dt)
    print()
    print(E)

```

```
[ ]: energy_lf([0,1/np.sqrt(2)], [1,0], [0.0001,6],10)
```

```

[[ 0.00000000e+00  7.07106781e-01]
 [-5.00000001e-05  7.07106778e-01]
 [-9.99999991e-05  7.07106771e-01]
 ...
 [-4.98254992e-01  5.01266286e-01]
 [-4.98290382e-01  5.01231103e-01]
 [-4.98325770e-01  5.01195918e-01]]

[[ 0.          0.70710678]
 [-0.05017395  0.70353752]
 [-0.09923215  0.69655773]
 [-0.1469802   0.6863332 ]

```

[-0.19323985 0.67303579]
[-0.23784881 0.65684262]
[-0.28066058 0.63793509]
[-0.32154416 0.61649799]
[-0.36038387 0.59271861]
[-0.39707909 0.56678581]
[-0.43154387 0.53888916]
[-0.46370666 0.50921802]
[-0.49350989 0.47796074]
[-0.52090949 0.44530377]
[-0.54587448 0.41143092]
[-0.56838632 0.37652259]
[-0.58843847 0.34075505]
[-0.60603563 0.3042998]
[-0.62119323 0.26732298]
[-0.63393662 0.2299848]
[-0.64430047 0.19243911]
[-0.65232797 0.15483293]
[-0.65807016 0.11730615]
[-0.66158514 0.07999122]
[-0.66293737 0.04301293]
[-0.66219692 0.00648828]
[-0.65943876 -0.02947366]
[-0.65474205 -0.06477174]
[-0.64818948 -0.09931277]
[-0.63986662 -0.13301139]
[-0.62986126 -0.16579001]
[-0.61826288 -0.19757866]
[-0.60516208 -0.22831474]
[-0.59065004 -0.25794286]
[-0.57481807 -0.28641457]
[-0.55775716 -0.31368808]
[-0.5395576 -0.33972798]
[-0.5203086 -0.36450492]
[-0.500098 -0.3879953]
[-0.47901194 -0.41018094]
[-0.45713465 -0.43104877]
[-0.43454823 -0.45059048]
[-0.41133245 -0.46880221]
[-0.38756464 -0.48568421]
[-0.36331951 -0.50124053]
[-0.33866911 -0.51547872]
[-0.31368272 -0.52840954]
[-0.28842684 -0.54004663]
[-0.26296512 -0.55040631]
[-0.23735835 -0.55950723]
[-0.21166453 -0.56737018]
[-0.18593879 -0.57401784]

[-0.16023353 -0.57947456]
[-0.13459837 -0.58376615]
[-0.10908025 -0.58691965]
[-0.0837235 -0.58896323]
[-0.05856987 -0.58992593]
[-0.03365865 -0.58983759]
[-0.00902669 -0.58872862]
[0.01529145 -0.58662996]
[0.0392635 -0.58357287]
[0.0628593 -0.5795889]
[0.08605081 -0.57470973]
[0.10881195 -0.56896712]
[0.13111855 -0.56239281]
[0.15294828 -0.55501844]
[0.17428053 -0.54687553]
[0.19509634 -0.53799536]
[0.21537833 -0.52840897]
[0.23511063 -0.51814709]
[0.25427879 -0.50724011]
[0.27286969 -0.49571806]
[0.29087152 -0.48361054]
[0.30827367 -0.47094676]
[0.32506668 -0.45775546]
[0.34124217 -0.44406495]
[0.35679279 -0.42990304]
[0.37171217 -0.41529709]
[0.38599486 -0.40027393]
[0.39963627 -0.38485996]
[0.41263261 -0.36908103]
[0.42498091 -0.35296254]
[0.43667889 -0.33652937]
[0.44772497 -0.31980595]
[0.45811824 -0.30281618]
[0.46785839 -0.28558353]
[0.47694569 -0.26813096]
[0.48538099 -0.250481]
[0.49316562 -0.2326557]
[0.50030144 -0.21467668]
[0.50679077 -0.19656512]
[0.51263638 -0.17834177]
[0.51784145 -0.16002694]
[0.52240959 -0.14164057]
[0.52634478 -0.12320217]
[0.52965138 -0.10473087]
[0.53233409 -0.08624542]
[0.53439797 -0.06776419]
[0.5358484 -0.0493052]]

```

[[ 0.          0.70710678]
 [-0.10100482  0.69252344]
 [-0.19339126  0.66584545]
 [-0.2767528   0.62923067]
 [-0.35088018  0.58458113]
 [-0.41572982  0.53358924]
 [-0.47139882  0.47776659]
 [-0.51810354  0.41846244]
 [-0.55616     0.35687609]
 [-0.58596566  0.29406605]
 [-0.60798199  0.23095762]
 [-0.6227183   0.16835019]
 [-0.63071658  0.10692456]
 [-0.63253789  0.04725088]
 [-0.62875023 -0.01020301]
 [-0.61991809 -0.0650627 ]
 [-0.60659374 -0.11703816]
 [-0.58931018 -0.16591456]
 [-0.56857571 -0.21154316]
 [-0.54486989 -0.25383268]
 [-0.51864094 -0.29274091]
 [-0.49030412 -0.32826708]
 [-0.46024121 -0.36044478]
 [-0.42880062 -0.38933557]
 [-0.39629825 -0.41502336]
 [-0.36301868 -0.43760941]
 [-0.32921681 -0.45720805]
 [-0.29511965 -0.47394295]
 [-0.26092828 -0.48794407]
 [-0.22681983 -0.49934499]
 [-0.19294955 -0.50828085]
 [-0.15945269 -0.51488653]
 [-0.12644648 -0.51929531]
 [-0.09403188 -0.52163778]
 [-0.06229528 -0.52204099]
 [-0.03131005 -0.52062785]
 [-0.00113802 -0.5175167 ]
 [ 0.02816922 -0.512821  ]
 [ 0.05656913 -0.50664919]
 [ 0.08402708 -0.49910458]
 [ 0.11051545 -0.49028541]
 [ 0.13601265 -0.48028491]
 [ 0.16050243 -0.46919137]
 [ 0.18397309 -0.45708837]
 [ 0.20641694 -0.44405489]
 [ 0.22782967 -0.43016555]
 [ 0.24820994 -0.41549082]
 [ 0.26755887 -0.4000972 ]

```

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