## hw3 ex12

## February 20, 2020

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[1]: #!/usr/bin/env python3
     # -*- coding: utf-8 -*-
     Created on Thu Feb 6 13:33:08 2020
     @author: aswart
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     # python 3 version 2/15
     import numpy as np
     import matplotlib.pyplot as plt
     def rk4(x,t,tau,derivsRK):
     #% Runge-Kutta integrator (4th order)
     #% Input arguments -
     \#\% x = current value of dependent variable
     #% t = independent variable (usually time)
        tau = step size (usually timestep)
     #%
         derivsRK = right hand side of the ODE; derivsRK is the
     #%
     #%
                    name of the function which returns dx/dt
                    Calling format derivsRK(x,t).
     #%
     #% Output arguments -
     #% xout = new value of x after a step of size tau
         half_tau = 0.5*tau
         F1 = derivsRK(x,t)
         t_half = t + half_tau
         xtemp = x + half tau*F1
         F2 = derivsRK(xtemp,t_half)
         xtemp = x + half_tau*F2
         F3 = derivsRK(xtemp,t_half)
         t_full = t + tau
         xtemp = x + tau*F3
         F4 = derivsRK(xtemp,t_full)
         xout = x + tau/6.*(F1 + F4 + 2.*(F2+F3))
         return xout
     def rka(x,t,tau,err,derivsRK):
```

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#% Adaptive Runge-Kutta routine
#% Inputs
#%
    \boldsymbol{x}
              Current value of the dependent variable
#%
              Independent variable (usually time)
#%
   tau
              Step size (usually time step)
              Desired fractional local truncation error
#%
   err
#% derivsRK Right hand side of the ODE; derivsRK is the
#%
              name of the function which returns dx/dt
#%
               Calling format derivsRK(x,t).
#% Outputs
            New value of the dependent variable
    xSmall
#%
#%
              New value of the independent variable
#%
   tau
               Suggested step size for next call to rka
#%* Set initial variables
   tSave = t; xSave = x # Save initial values
    safe1 = .9; safe2 = 4. # Safety factors
    eps = np.spacing(1) # smallest value
#%* Loop over maximum number of attempts to satisfy error bound
   maxTry = 100
   for iTry in range(1,maxTry):
#%* Take the two small time steps
       half tau = 0.5 * tau
       xTemp = rk4(xSave,tSave,half_tau,derivsRK)
       t = tSave + half tau
       xSmall = rk4(xTemp,t,half_tau,derivsRK)
  #%* Take the single big time step
       t = tSave + tau
        xBig = rk4(xSave,tSave,tau,derivsRK)
  #%* Compute the estimated truncation error
        scale = err * (np.abs(xSmall) + np.abs(xBig))/2.
        xDiff = xSmall - xBig
        errorRatio = np.max( [np.abs(xDiff)/(scale + eps)] )
        #print safe1, tau, errorRatio
  #%* Estimate news tau value (including safety factors)
       tau old = tau
       tau = safe1*tau_old*errorRatio**(-0.20)
        tau = np.max([tau,tau_old/safe2])
        tau = np.min([tau,safe2*tau_old])
```

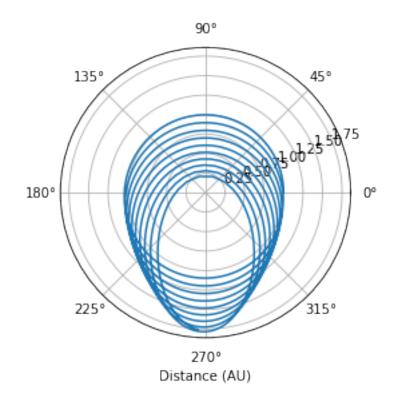
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#%* If error is acceptable, return computed values
        if errorRatio < 1 :</pre>
          \# xSmall = xSmall \#\% + (xDiff)/15
          \# xSmall = (16.*xSmall - xBig)/15. <math>\# correction
            return xSmall, t, tau
#%* Issue error message if error bound never satisfied
    print ('ERROR: Adaptive Runge-Kutta routine failed')
    return
def gravrk(s,t):
#% Returns right-hand side of Kepler ODE; used by Runge-Kutta routines
#% Inputs
#% s State vector [r(1) \ r(2) \ v(1) \ v(2)]
#%
    t
            Time (not used)
#% Output
#% deriv Derivatives [dr(1)/dt dr(2)/dt dv(1)/dt dv(2)/dt]
    GM = 4*np.pi**2
#%* Compute acceleration
    r = np.array([s[0], s[1]]) # Unravel the vector s into position and
\rightarrow velocity
    v = np.array([s[2],s[3]])
    accel = -GM*r/np.linalg.norm(r)**3 # Gravitational acceleration
#%* Return derivatives [dr(1)/dt \ dr(2)/dt \ dv(1)/dt \ dv(2)/dt]
    derivs = np.array([v[0], v[1], accel[0], accel[1]])
    return derivs
def gravrk_ex12(s,t):
#% Returns right-hand side of Kepler ODE; used by Runge-Kutta routines
#% Inputs
#%
    S
            State vector [r(1) \ r(2) \ v(1) \ v(2)]
#% t
            Time (not used)
#% Output
    deriv Derivatives [dr(1)/dt dr(2)/dt dv(1)/dt dv(2)/dt]
#%
    GM = 4*np.pi**2
    # Hardcoded in vars since
    # Im not sure on an elegant way to add them to function
    r0 = np.array([1.0,0])
    init_Grav_accel = .01*-GM*r0/np.linalg.norm(r0)**3
#%* Compute acceleration
    r = np.array([s[0], s[1]]) # Unravel the vector s into position and
\rightarrow velocity
```

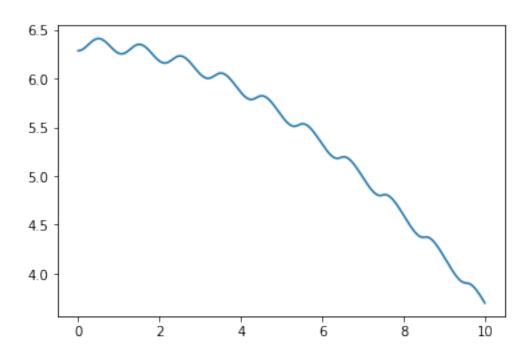
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v = np.array([s[2],s[3]])
    # Gravitational acceleration
    accel = -GM*r/np.linalg.norm(r)**3 + init_Grav_accel
#%* Return derivatives [dr(1)/dt \ dr(2)/dt \ dv(1)/dt \ dv(2)/dt]
    derivs = np.array([v[0], v[1], accel[0], accel[1]])
    return derivs
# orbit - Program to compute the orbit of a comet.
#clear all; help orbit; % Clear memory and print header
def orbit(input_dict = {}, calc_info = False, plot_momentum = False,
          inter_input = False, plot_traj = True, plot_energy = False):
    if inter input:
        # Set initial position and velocity of the comet.
        r0 = float(input("Enter initial radial distance (AU): "))
        # v0 = float(input("Enter initial tangential velocity (AU/yr): "))
        vstr = str(input("Enter initial tangential velocity (AU/yr) as a number ⊔
→or multiple of Pi (e.g.,2*pi): "))
        # modify input to allow 'pi'
        vinp = vstr.split('*')
        if (vinp[-1].lower() == 'pi'):
          v0 = float(vinp[0])*np.pi
        else:
          v0 = float(vinp[0])
        nStep = int(input("Enter number of steps: "))
        tau = float(input("Enter time step (yr): "))
        NumericalMethod=0
        while(NumericalMethod not in np.array([1,2,3,4,5,6,7])):
            NumericalMethod = int(input("Choose a number for a numerical method:
\hookrightarrow \n
         1-Euler, 2-Euler-Cromer, 3-Runge-Kutta 4-Adaptive R-K: "))
    elif input dict:
        r0 = input_dict['r0']
        v0 = input_dict['v0']
        nStep = input_dict['nStep']
        tau = input_dict['tau']
        NumericalMethod = input_dict['NumericalMethod']
    else:
        r0 = 1
        v0 = 2*np.pi
        nStep = 1000
        tau = .01
        NumericalMethod = 3
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r = np.array([r0, 0.])
   v = np.array([0., v0])
   state = np.array([ r[0], r[1], v[0], v[1] ]) # Used by R-K routines
   #Set physical parameters (mass, G*M)
   GM = 4*np.pi**2  # Grav. const. * Mass of Sun (au^3/yr^2)
   mass = 1. # Mass of comet
   adaptErr = 1.e-4 # Error parameter used by adaptive Runge-Kutta
   time = 0.0
   #%* Loop over desired number of steps using specified
   #% numerical method.
   for istep in range(0,nStep):
     #%* Record position and energy for plotting.
     # Initially set the arrays for the first step
     if istep == 0:
         rplot = np.linalg.norm(r)
         thplot = np.arctan2(r[1],r[0])
         tplot = time
         kinetic = .5*mass*np.linalg.norm(v)**2
         potential= - GM*mass/np.linalg.norm(r)
         momentum = [np.linalg.norm(np.cross(r,mass*v))]
     else:
         rplot = np.append(rplot,np.linalg.norm(r))
                                                      #Record position
→ for polar plot
         thplot = np.append(thplot,np.arctan2(r[1],r[0]))
         tplot = np.append(tplot,time)
         kinetic = np.append(kinetic,0.5*mass*np.linalg.norm(v)**2) # Record
\rightarrow energies
         potential= np.append(potential, - GM*mass/np.linalg.norm(r))
         momentum.append(np.linalg.norm(np.cross(r, mass*v)))
     #%* Calculate new position and velocity using desired method.
     if NumericalMethod == 1 :
       accel = -GM*r/np.linalg.norm(r)**3
      r = r + tau*v
                                # Euler step
      v = v + tau*accel
      time = time + tau
     elif NumericalMethod == 2:
      accel = -GM*r/np.linalg.norm(r)**3
      v = v + tau*accel
      r = r + tau*v
                                 # Euler-Cromer step
      time = time + tan
     elif NumericalMethod == 3:
       state = rk4(state,time,tau,gravrk_ex12)
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r = np.array([state[0], state[1]])
                                          # 4th order Runge-Kutta
      v = np.array([state[2], state[3]])
      time = time + tau
       [state, time, tau] = rka(state,time,tau,adaptErr,gravrk)
      r = np.array([state[0], state[1]])
                                          # Adaptive Runge-Kutta
      v = np.array([state[2], state[3]])
     # If sometime after first step and theta goes from neg to pos, then we_
→ know we've completed an orbit
      if istep != 0 and (thplot[-2:]*[-1,1] > 0).all():
#
          break
   #print(thplot[-5:])
  if plot_traj:
       #%* Graph the trajectory of the comet.
      plt.figure(1); plt.clf() #Clear figure 1 window and bring forward
      plt.polar(thplot,rplot,'-') # Use polar plot for graphing orbit
      plt.xlabel('Distance (AU)')
      plt.grid(True)
   if plot_energy:
       #%* Graph the energy of the comet versus time.
      plt.figure(2); plt.clf() # Clear figure 2 window and bring forward
      totalE = kinetic + potential
                                    # Total energy
      plt.plot(tplot,kinetic,'-.',tplot,potential,'--',tplot,totalE,'-')
      #plt.legend('Kinetic', 'Potential', 'Total')
      plt.xlabel('Time (yr)'); plt.ylabel('Energy (M AU^2/yr^2)')
      plt.grid(True)
      plt.show()
   if plot_momentum:
       # Plots angular momentum as a function of time
      plt.figure(3)
      plt.plot(tplot,momentum)
  return rplot, thplot
```

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[2]: # non-elliptical
input_dict = {
    'r0': 1,
    'v0': 2*np.pi,
    'nStep': 1000,
    'tau': .01,
    'NumericalMethod': 3
    }
    rplot, thplot = orbit(input_dict,plot_momentum = True)
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