Chapter 7. Secular Perturbations

Action by Daniel Niño-Villegas, University of Antioquia

In this notebook we will present part of the theoretical background of the chapter devoted to *Secular Perturbations* and some mathematical and numerical results which are interested for the theory.

For details on the theory please refer directly to the book:

Murray, C. D., & Dermott, S. F. (1999). Solar system dynamics. Cambridge university press.

Preliminaries

Prerrequisites

```
In [1]: 1 #!pip install -q rebound
2 #!pip install -q Fraction
3 #!pip install -q ipywidgets
```

Other libraries

```
In [2]:

1 #Global packages
2 import numpy as np
3 import matplotlib.pyplot as plt
4 import matplotlib.animation as animation
5 import rebound as rb
6 import celluloid as cell
7 from secular import *

8

9 #Specific modules and routines
10 from tqdm import tqdm
11 from ipywidgets import interact, widgets, fixed
12 from fractions import Fraction
13 from IPython.core.display import HTML
```

Useful constants

```
In [3]: 1 deg = np.pi/180
2 rad = 1/deg
```

Plots aesthetics

Section 7.2 Secular Perturbations for Two Planets

Experiment: Laplace coefficients study (7.1)

In §6.9 we introduced the computation of the Laplace coefficients and their derivatives (up to second order), who are completely essential in the development of perturbation theory. This functions are until now completely "dark" in a sense that we don't know how they behave or even how their derivatives behave. So, we'll now show an interactive plot of $b_s^{(j)}(\alpha)$, D $b_s^{(j)}(\alpha)$ and D^2 $b_s^{(j)}(\alpha)$ where you can modify the parameters s and j to see how the coefficents change:

```
In [5]:
         1 def blap_plot(s,j):
                a = np.linspace(0, 0.9, 1000)
         2
                b = np.array([blap(x, s, j) for x in a])
         3
         4
                db = np.array([blap_dot(x, s, j) for x in a])
                ddb = np.array([blap_ddot(x, s, j) for x in a])
                i = int(i)
         8
                s = str(Fraction(s).numerator) + '/' + str(Fraction(s).denominator)
         9
                ss = fr'$s = {s}$' + r'$\hspace{20 pt}$' fr'$j = {j}$'
        10
        11
                plt.close()
                fig, axs = plt.subplots(3, 1, figsize=(4,6), sharex=True, dpi=110)
        12
        13
                axs[0].set_title(ss, fontsize=15)
        14
                axs[0].set_ylabel(r'$b_{s}^{(j)}(\lambda)$', fontsize=17)
        15
                axs[0].plot(a, b, 'k-')
        16
                axs[1].set\_xlabel(r'\$\alpha\$', fontsize=15)
                axs[1].set\_ylabel(r'$D \ \ b_{s}^{(j)}(\alpha)$', fontsize=15)
        17
        18
                axs[1].plot(a, db, 'k-')
        19
                axs[2].set_xlabel(r'$\alpha$', fontsize=15)
        20
                axs[2].set_ylabel(r'$D^2 \, b_{s}^{(j)}(\alpha)$', fontsize=15)
        21
                axs[2].plot(a, ddb, 'k-')
        22
                fig.tight_layout()
        23
                plt.subplots_adjust(hspace=0)
        24
                plt.show()
In [6]:
         1 opciones = dict(continuous_update=False)
         2 interact(blap_plot,
         3
                     s = widgets.FloatSlider(min=0.5, max=5, step=0.5, value=0.5, **opciones
                     j = widgets.FloatSlider(min=0, max=5, step=1, value=0, **opciones))
```

A Jupyter widget could not be displayed because the widget state could not be found. This could happen if the kernel storing the widget is no longer available, or if the widget state was not saved in the notebook. You may be able to create the widget by running the appropriate cells.

```
<function __main__.blap_plot(s, j)>
```

An important result of the previous plots is that the Laplace coefficients diverge rapidly near $\alpha=1$, and so their derivatives. You can also check that if you increase the value of the s parameter, the coefficient is virtually cero for a large number of α values, and diverge much quickly near $\alpha=1$. Be free to combine different parameter values to check all of this for yourself.

Experiment: excentricity and inclination vectors (7.2)

For the following experiment we'll suppose a system made up with a central body of mass m_c and two orbiting bodies with masses m_1 and m_2 , where $m_1 \ll m_c$ and $m_2 \ll m_c$; for convention we'll take the inner body to be m_1 and the outter body to be m_2 . In the previous notebook we developed the theory for a similar system where $m_1 \ll m_2$, but now we'll describe the motion for two particles ignoring this condition. Let's write the disturbing function for both bodies like we did before:

Initial conditions:

$$\mathcal{R}_1 = \frac{Gm_2}{a_1} \alpha_{12} \mathcal{R}_D^{(\text{sec})}, \qquad \mathcal{R}_2 = \frac{Gm_1}{a_2} \mathcal{R}_D^{(\text{sec})}$$
 (1)

Where:

$$\mathcal{R}_{D}^{(\text{sec})} = \frac{1}{8} \left[2\alpha_{12}D + \alpha_{12}^{2}D^{2} \right] b_{\frac{1}{2}}^{(0)} \left(e_{1}^{2} + e_{2}^{2} \right) - \frac{1}{2}\alpha_{12}b_{3/2}^{(1)} \left(s_{1}^{2} + s_{2}^{2} \right)$$

$$+ \frac{1}{4} \left[2 - 2\alpha_{12}D - \alpha_{12}^{2}D^{2} \right] b_{\frac{1}{2}}^{(1)} e_{1}e_{2} \cos(\varpi_{1} - \varpi_{2})$$

$$+ \alpha_{12}b_{3/2}^{(1)} s_{1} s_{2} \cos(\Omega_{1} - \Omega_{2})$$

$$(2)$$

Note that we've dropped the $\frac{1}{2}b_{\frac{1}{2}}^{(0)}(\alpha)$ term in $\mathcal{R}_{\mathrm{D}}^{(\mathrm{sec})}$ due to the fact that Lagrange's

equations for a_1 and a_2 are proportional to $\frac{\partial \mathcal{R}}{\partial \lambda_j}$, and in a secular regime we won´t consider terms associated with the mean anomalies. Using the recursion relations for the Laplace coefficients it can be found that:

$$2\alpha \frac{\mathrm{d}b_{1/2}^{(0)}}{\mathrm{d}\alpha} + \alpha^2 \frac{\mathrm{d}^2 b_{1/2}^{(0)}}{\mathrm{d}\alpha^2} = \alpha b_{3/2}^{(1)} \tag{3}$$

$$2b_{1/2}^{(1)} - 2\alpha \frac{\mathrm{d}b_{1/2}^{(1)}}{\mathrm{d}\alpha} - \alpha^2 \frac{\mathrm{d}^2b_{1/2}^{(1)}}{\mathrm{d}\alpha^2} = -\alpha b_{3/2}^{(2)} \tag{4}$$

Plugging (3) and (4) in (2) we find that:

$$\mathcal{R}_{D}^{(\text{sec})} = \frac{1}{8} \alpha b_{3/2}^{(1)} \left(e_{1}^{2} + e_{2}^{2} \right) - \frac{1}{2} \alpha_{12} b_{3/2}^{(1)} \left(s_{1}^{2} + s_{2}^{2} \right) - \frac{1}{4} \alpha b_{3/2}^{(2)} e_{1} e_{2} \cos(\varpi_{1} - \varpi_{2}) + \alpha_{12} b_{3/2}^{(1)} s_{1} s_{2} \cos(\Omega_{1} - \Omega_{2})$$
(5)

Now we can replace (5) in eqs. (1):

$$\mathcal{R}_{1} = \frac{Gm_{2}}{a_{1}} \left[\frac{1}{8} \alpha_{12}^{2} b_{3/2}^{(1)} (e_{1}^{2} + e_{2}^{2}) - \frac{1}{8} \alpha_{12}^{2} b_{3/2}^{(1)} (s_{1}^{2} + s_{2}^{2}) - \frac{1}{4} \alpha_{12}^{2} b_{3/2}^{(2)} e_{1} e_{2} \cos(\varpi_{1} - \varpi_{2}) + \frac{1}{4} \alpha_{12}^{2} b_{3/2}^{(1)} I_{1} I_{2} \cos(\Omega_{1} - \Omega_{2}) \right]$$

$$(6)$$

$$\mathcal{R}_{2} = \frac{Gm_{1}}{a_{2}} \left[\frac{1}{8} \alpha_{12} b_{3/2}^{(1)} (e_{1}^{2} + e_{2}^{2}) - \frac{1}{8} \alpha_{12} b_{3/2}^{(1)} (s_{1}^{2} + s_{2}^{2}) - \frac{1}{4} \alpha_{12} b_{3/2}^{(2)} e_{1} e_{2} \cos(\varpi_{1} - \varpi_{2}) + \frac{1}{4} \alpha_{12} b_{3/2}^{(1)} I_{1} I_{2} \cos(\Omega_{1} - \Omega_{2}) \right]$$

$$(7)$$

With this equations, we can now do a little handling:

$$\mathcal{R}_{1} = \frac{Gm_{2}}{a_{1}} \left[\frac{1}{8} \alpha_{12}^{2} b_{3/2}^{(1)} e_{1}^{2} - \frac{1}{8} \alpha_{12}^{2} b_{3/2}^{(1)} s_{1}^{2} \right.$$

$$- \frac{1}{4} \alpha_{12}^{2} b_{3/2}^{(2)} e_{1} e_{2} \cos(\varpi_{1} - \varpi_{2})$$

$$+ \frac{1}{4} \alpha_{12}^{2} b_{3/2}^{(1)} I_{1} I_{2} \cos(\Omega_{1} - \Omega_{2}) \right]$$

$$+ \underbrace{\frac{Gm_{2}}{a_{1}} \left[\frac{1}{8} \alpha_{12}^{2} b_{3/2}^{(1)} e_{2}^{2} - \frac{1}{8} \alpha_{12}^{2} b_{3/2}^{(1)} s_{2}^{2} \right]}_{(8)}$$

$$\mathcal{R}_{2} = \frac{Gm_{1}}{a_{2}} \left[\frac{1}{8} \alpha_{12} b_{3/2}^{(1)} e_{2}^{2} - \frac{1}{8} \alpha_{12} b_{3/2}^{(1)} s_{2}^{2} - \frac{1}{4} \alpha_{12} b_{3/2}^{(2)} e_{1} e_{2} \cos(\varpi_{1} - \varpi_{2}) + \frac{1}{4} \alpha_{12} b_{3/2}^{(1)} I_{1} I_{2} \cos(\Omega_{1} - \Omega_{2}) \right] + \underbrace{\frac{Gm_{1}}{a_{2}} \left[\frac{1}{8} \alpha_{12} b_{3/2}^{(1)} e_{1}^{2} - \frac{1}{8} \alpha_{12} b_{3/2}^{(1)} s_{1}^{2} \right]}_{(9)}$$

You can surely note that the underlined terms will have no effect in the Lagrange's equations because they point to the other particle, so the derivatives of these will be zero. Now we'll use Kepler's third law approximation:

$$G(m_c + m_1) \approx n_1^2 a_1^3, \quad G(m_c + m_2) \approx n_2^2 a_2^3$$

This yields:

$$\frac{G}{a_1} = \frac{n_1^2 a_1^2}{m_c + m_1}, \quad \frac{G}{a_2} = \frac{n_2^2 a_2^2}{m_c + m_2}$$
 (10)

Replacing $(\underline{10})$ in $(\underline{8})$ and $(\underline{9})$, and assuming that the approximations $s_1 \approx \frac{1}{2}I_1$,

 $s_2 \approx \frac{1}{2}I_2$ are valid we get:

$$\mathcal{R}_{1} = n_{1}^{2} a_{1}^{2} \frac{m_{2}}{m_{c} + m_{1}} \left[\frac{1}{8} \alpha_{12}^{2} b_{3/2}^{(1)} e_{1}^{2} - \frac{1}{8} \alpha_{12}^{2} b_{3/2}^{(1)} I_{1}^{2} - \frac{1}{4} \alpha_{12}^{2} b_{3/2}^{(2)} e_{1} e_{2} \cos(\varpi_{1} - \varpi_{2}) + \frac{1}{4} \alpha_{12}^{2} b_{3/2}^{(1)} I_{1} I_{2} \cos(\Omega_{1} - \Omega_{2}) \right]$$

$$(11)$$

$$\mathcal{R}_{2} = n_{2}^{2} a_{2}^{2} \frac{m_{1}}{m_{c} + m_{2}} \left[\frac{1}{8} \alpha_{12} b_{3/2}^{(1)} e_{2}^{2} - \frac{1}{8} \alpha_{12} b_{3/2}^{(1)} I_{2}^{2} - \frac{1}{4} \alpha_{12} b_{3/2}^{(2)} e_{1} e_{2} \cos(\varpi_{1} - \varpi_{2}) + \frac{1}{4} \alpha_{12} b_{3/2}^{(1)} I_{1} I_{2} \cos(\Omega_{1} - \Omega_{2}) \right]$$

$$(12)$$

Eqs. (11) and (12) are pretty similar, so it's not crazy to find a way to merge them in an unique expression. Let's proppose the next one:

$$\mathcal{R}_{j} = n_{j} a_{j}^{2} \left[\frac{1}{2} A_{jj} e_{j}^{2} + A_{jk} e_{j} e_{k} \cos(\varpi_{j} - \varpi_{k}) \right]$$

$$\frac{1}{2} B_{jj} I_{j}^{2} + B_{jk} I_{j} I_{k} \cos(\Omega_{j} - \Omega_{k})$$
(13)

Where:

$$A_{jj} = +n_j \frac{1}{4} \frac{m_k}{m_c + m_j} \alpha_{12} \bar{\alpha}_{12} b_{3/2}^{(1)} (\alpha_{12})$$
 (14)

$$A_{jk} = -n_j \frac{1}{4} \frac{m_k}{m_c + m_j} \alpha_{12} \bar{\alpha}_{12} b_{3/2}^{(2)} (\alpha_{12})$$
 (15)

$$B_{jj} = -n_j \frac{1}{4} \frac{m_k}{m_c + m_j} \alpha_{12} \bar{\alpha}_{12} b_{3/2}^{(1)} (\alpha_{12})$$
 (16)

$$B_{jk} = +n_j \frac{1}{4} \frac{m_k}{m_c + m_j} \alpha_{12} \bar{\alpha}_{12} b_{3/2}^{(1)} (\alpha_{12})$$
 (17)

In this equations you should be care that j=1,2, k=2,1 and $j\neq k$. Also $\bar{\alpha}_{12}=\alpha_{12}$ if j=1 and $\bar{\alpha}_{12}=1$ if j=2. These numbers (who have frequency units) constitute two matrices ${\bf A}$ and ${\bf B}$:

$$\mathbf{A} = \begin{pmatrix} A_{11} & A_{12} \\ A_{21} & A_{22} \end{pmatrix} \tag{18}$$

$$\mathbf{B} = \begin{pmatrix} B_{11} & B_{12} \\ B_{21} & B_{22} \end{pmatrix} \tag{19}$$

For upcoming developments it's convenient to define the vertical and horizontal components of the so called *eccentricity vectors*:

$$h_j = e_j \sin \varpi_j, \quad k_j = e_j \cos \varpi_j$$
 (20)

$$p_i = I_i \sin \Omega_i, \quad q_i = I_i \cos \Omega_i$$
 (21)

It can be shown immediately that:

$$h_i^2 + k_i^2 = e_i^2 (22)$$

$$p_i^2 + q_i^2 = I_i^2 (23)$$

$$h_j h_k + k_j k_k = e_j e_k \cos(\varpi_j - \varpi_k)$$
 (24)

$$p_i p_k + q_i q_k = I_i I_k \cos(\Omega_i - \Omega_k) \tag{25}$$

Replacing Eqs. (22), (23), (24) and (25) in (13) yields:

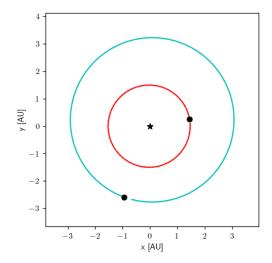
$$\mathcal{R}_{j} = n_{j} a_{j}^{2} \left[\frac{1}{2} A_{jj} \left(h_{j}^{2} + k_{j}^{2} \right) + A_{jk} \left(h_{j} h_{k} + k_{j} k_{k} \right) + \frac{1}{2} B_{jj} \left(p_{j}^{2} + q_{j}^{2} \right) + B_{jk} \left(p_{j} p_{k} + q_{j} q_{k} \right) \right]$$
(26)

These four vectors are completely dark in a sense that they are just mathematical tools for the theory purpose, so our goal is to visualize their behaviour and their evolution in a real physical system. Let's now take a system made up with two similar mass bodies orbiting a central point mass m_c , the following code cells will show the procedure to simulate the time evolution of this system (just like we did in §6.9) and the eccentricity vectors h_j and p_j using rebound

Let's first define the initial conditions for the a, e, and ϖ of the two bodies and create the simulation:

```
In [7]:
        1 #Partticle #1
        2 m1 = 1e-4
        3 a1 = 1.5
        4 e1 = 0.02
           pomega1 = 10*deg
        6
        7
           #Particle #2
           m2 = 2e-4
        9 a2 = 3
        10 e2 = 0.08
        11 pomega2 = 250*deg
        12
        13 #Smulation t=0
        14 sim = rb.Simulation()
        15 sim.units = ('au', 'msun', 'yr')
        16 sim.add(m=1, hash='Sun')
        17 sim.add(m=m1, a=a1, e=e1, pomega=pomega1, hash='Particle #1')
        18 sim.add(m=m2, a=a2, e=e2, pomega=pomega2, hash='Particle #2')
        19 #sim.save('tmp/system.bin')
```

The initial state of the system can be visualized in the next way:



Now we'll calculate the mean orbit periods who are very useful as a time step to simulate:

```
In [9]: 1 #Orbital periods
2 P1 = sim.particles['Particle #1'].P
3 n1 = 2*np.pi/P1
4 P2 = sim.particles['Particle #2'].P
5 n2 = 2*np.pi/P2
```

Now let's define the integration parameters in terms of the orbit period of the inner body:

```
In [10]: 1 #Integration parameters
2 sim.dt = P1/100
3 Nt = 1000
4 ts = np.linspace(0, 200000, Nt)
```

Finally let's do the integration refered to the central body reference system:

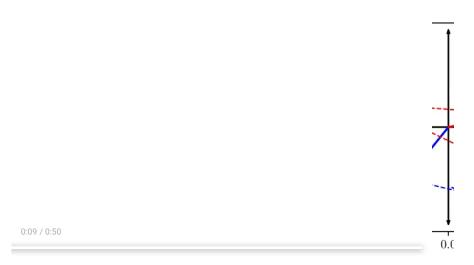
```
In [11]: 1 #Integration
         2 Es = np.zeros((2,Nt,3))
         3 for i,t in enumerate(tqdm(ts)):
               sim.integrate(t)
               sim.move_to_hel()
         6
               orbits = sim.calculate_orbits()
         7
               Es[0][i] = [orbits[0].a,
                           orbits[0].e,
                           np.mod(orbits[0].pomega,2*np.pi)]
         9
        10
               Es[1][i] = [orbits[1].a,
        11
                           orbits[1].e,
        12
                           np.mod(orbits[1].pomega,2*np.pi)]
         100%|
                                                      1000/1000 [00:48<00:00,
         20.56it/s]
```

With the simulation done, we can split the array with the results in individual arrays for the orbital elements of each particle:

Now we'll define a routine who given the semimajor axis, eccentricity and longitude of pericentre of the two main bodies at a certain time, it will plot the two osculating orbits and the vectors h_i and k_i :

```
In [13]:
                                 1 def vec_orbit_plot(a1, a2, e1, e2, pomega1, pomega2, L, i, Np=1000, s=1.1):
                                 2
                                                        #Excentricity vectors
                                 3
                                                        hvec = np.zeros(2)
                                 4
                                                        kvec = np.zeros(2)
                                 5
                                 6
                                                         #Components
                                 7
                                                        hvec[0] = e1[i]*np.sin(pomega1[i])
                                 8
                                                        hvec[1] = e2[i]*np.sin(pomega2[i])
                                 9
                               10
                                                        kvec[0] = e1[i]*np.cos(pomega1[i])
                              11
                                                        kvec[1] = e2[i]*np.cos(pomega2[i])
                              12
                              13
                                                         #Orbits
                              14
                                                        f = np.linspace(0, 2*np.pi, Np) #True anomaly
                              15
                                                        p1 = a1[i]*(1-e1[i]**2)
                                                                                                                                                                          #Semilatus rectum #1
                              16
                                                        p2 = a2[i]*(1-e2[i]**2)
                                                                                                                                                                         #Semilatus rectum #2
                              17
                              18
                                                         ##Polar ellipse equation #1
                              19
                                                        x1 = (p1/(1 + e1[i]*np.cos(f)))*np.cos(f)
                              20
                                                        y1 = (p1/(1 + e1[i]*np.cos(f)))*np.sin(f)
                              21
                              22
                                                         ##Polar ellipse equation #2
                              23
                                                        x2 = (p2/(1 + e2[i]*np.cos(f)))*np.cos(f)
                              24
                                                        y2 = (p2/(1 + e2[i]*np.cos(f)))*np.sin(f)
                              25
                              26
                                                        ##Rotated coordinates #1
                              27
                                                        x1_ = x1*np.cos(pomega1[i]) - y1*np.sin(pomega1[i])
                                                        y1_ = x1*np.sin(pomega1[i]) + y1*np.cos(pomega1[i])
                              28
                              29
                              30
                                                         ##Rotated coordinates #2
                              31
                                                        x2_ = x2*np.cos(pomega2[i]) - y2*np.sin(pomega2[i])
                                                        y2_ = x2*np.sin(pomega2[i]) + y2*np.cos(pomega2[i])
                              32
                              33
                              34
                                                        a = max([a1[i],a2[i]])
                              35
                                                        e = max(\lceil e1\lceil i \rceil, e2\lceil i \rceil \rceil)
                              36
                                                         #Complete Plot
                              37
                              38
                                                         ##Orbits
                                                        axs[0].text(-3.6, 3.3, r'$\varpi_1 =$ %.0f$^{\circ}$'%(pomega1[i]*rad))
                              39
                                                        axs[0].text(1.4, 3.3, r'$\varpi_2 =$ %.0f$^{\circ}$'%(pomega2[i]*rad))
                              40
                              41
                                                        axs[0].set_xlabel('x [AU]', fontsize=12)
                              42
                                                        axs[0].set_ylabel('y [AU]', fontsize=12)
                              43
                                                        axs[0].set_xlim((-a*(1+e)*s,a*(1+e)*s))
                              44
                                                        axs[0].set_ylim((-a*(1+e)*s,a*(1+e)*s))
                              45
                                                        axs[0].plot(x1_, y1_, 'k-', lw=1)
                              46
                                                        axs[0].plot(x2_, y2_, 'k-', lw=1)
                              47
                                                        axs[0].plot([0],[0], 'y*', ms=5)
                                                        axs[0].arrow(0, 0, a1[i]*(1-e1[i])*np.cos(pomega1[i]), a1[i]*(1-e1[i])*np.si
                              48
                              49
                                                        axs[0].arrow(0, 0, a2[i]*(1-e2[i])*np.cos(pomega2[i]), a2[i]*(1-e2[i])*np.si
                              50
                                                        axs[0].set_aspect('equal', 'box')
                              51
                                                        ##Excentricity Vectors
                              52
                                                        axs[1].set_xlim(-s*L,s*L)
                              53
                                                        axs[1].set_ylim(-s*L,s*L)
                              54
                                                        axs[1].text(1.15*hvec[0], 1.15*hvec[1], r'$\vec{h}$', fontsize=14)
                              55
                                                        axs[1].text(1.15*kvec[0], 1.15*kvec[1], r'$\vec{k}$', fontsize=14)
                              56
                                                        ####Axis
                                                        axs[1].arrow(0, 0, 0.9*s*L, 0, color='k', width=5e-4, head_length=0.005, head_length=0.
                              57
                              58
                                                        axs[1].arrow(0, 0, 0, 0.9*s*L, color='k', width=5e-4, head_length=0.005, head_length=0.
                              59
                                                        axs[1].arrow(0, 0, -0.9*s*L, 0, color='k', width=5e-4, head_length=0.005, he
                                                        axs[1].arrow(0, 0, 0, -0.9*s*L, color='k', width=5e-4, head_length=0.005, he
                              60
                              61
                                                        axs[1].arrow(0, 0, hvec[0], hvec[1], color='b')
                              62
                              63
                                                        axs[1].arrow(0, 0, kvec[0], kvec[1], color='r')
                              64
                                                        ####Curves
                                                        axs[1].plot(e1[:i]*np.sin(pomega1[:i]), e2[:i]*np.sin(pomega2[:i]), 'b--', label{eq:constraint} axs[1].plot(e1[:i]*np.sin(pomega1[:i]), 'b--', label{eq:constraint} axs[1].plot(e1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*np.sin(pomega1[:i]*
```

```
axs[1].plot(e1[:i]*np.cos(pomega1[:i]), e2[:i]*np.cos(pomega2[:i]), 'r--', ]
         66
          Finally let's make an animation of the time evolution of the osculating orbits and the
          excentricity vectors (with the path they follow):
In [14]:
          1 plt.ioff()
          2 fig, axs = plt.subplots(1, 2, dpi=150, figsize=(7,3))
          3 camera = cell.Camera(fig)
            for i in range(len(e1)):
          6
                 vec_orbit_plot(a1, a2, e1, e2, pomega1, pomega2, 0.1, i, s=1.3)
          7
                 camera.snap()
          8
          9 plt.close()
         10 plt.ion();
         11
         12 anim = camera.animate(interval=50)
         13 anim.save("figs/orbits_vectors.mp4")
          The animation is shown below:
In [15]:
          1 HTML(anim.to_html5_video())
```



As we can see, the osculating orbits of the two bodies rotate with different velocity, connecting with the path followed by the eccentricity vectors. Yo should have noticed that the path followed by these vectors is bounded, thus they oscillate between the first and third quadrant.