

Assignment 10 and optional part–Computation of homoclinic orbits of the equilibrium points

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1 Restricted Three-Body Problem

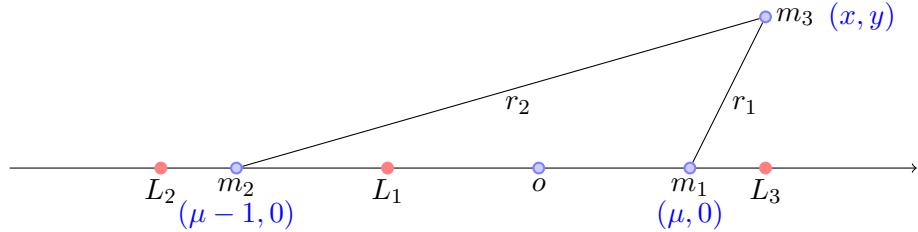


Figure 1: Restricted Three-Body Problem

The equations of motion are:

$$\begin{cases} x'' - 2y' = \Omega_x \\ y'' + 2x' = \Omega_y \end{cases} \quad (1)$$

And,

$$\Omega(x, y) = \frac{(x^2 + y^2)}{2} + \frac{1-\mu}{r_1} + \frac{\mu}{r_2} + \frac{\mu(1-\mu)}{2} \quad (2)$$

Here:

$$\begin{cases} \mu &= \frac{m_2}{m_1+m_2} \\ r_1 &= \sqrt{(x-\mu)^2 + y^2} \\ r_2 &= \sqrt{(x-\mu+1)^2 + y^2} \end{cases} \quad (3)$$

Let:

$$\begin{cases} x_1 &= x \\ x_2 &= y \\ x_3 &= x' \\ x_4 &= y' \end{cases} \quad (4)$$

The RTBP is expressed as:

$$\begin{cases} f_1 &= x'_1 = x_3 \\ f_2 &= x'_2 = x_4 \\ f_3 &= x'_3 = 2x_4 + \Omega_{x_1} \\ f_4 &= x'_4 = -2x_3 + \Omega_{x_2} \end{cases} \quad (5)$$

Where,

$$\begin{cases} \Omega_{x_1} &= x_1 - \frac{(1-\mu)(x_1-\mu)}{r_1^3} - \frac{\mu(x_1-\mu+1)}{r_2^3} \\ \Omega_{x_2} &= x_2(1 - \frac{1-\mu}{r_1^3} - \frac{\mu}{r_2^3}) \end{cases} \quad (6)$$

2 Equilibrium points— L_1 , L_2 , L_3

The position of L_1 is:

$$x_{L_1} = \mu - 1 + \xi \quad (7)$$

where,

$$\begin{aligned} f(\xi) &= \left(\frac{\mu(1-\xi)^2}{3 - 2\mu - \xi(3 - \mu - \xi)} \right)^{\frac{1}{3}} \\ \xi_0 &= \left(\frac{\mu}{3(1-\mu)} \right)^{\frac{1}{3}} \\ \xi_{n+1} &= f(\xi_n) \end{aligned}$$

The position of L_2 is:

$$x_{L_2} = \mu - 1 - \xi \quad (8)$$

where,

$$\begin{aligned} f(\xi) &= \left(\frac{\mu(1+\xi)^2}{3-2\mu+\xi(3-\mu+\xi)} \right)^{\frac{1}{3}} \\ \xi_0 &= \left(\frac{\mu}{3(1-\mu)} \right)^{\frac{1}{3}} \\ \xi_{n+1} &= f(\xi_n) \end{aligned}$$

The position of L_3 is:

$$x_{L_3} = \mu + \xi \quad (9)$$

where,

$$\begin{aligned} f(\xi) &= \left(\frac{(1-\mu)(1+\xi)^2}{1+2\mu+\xi(2+\mu+\xi)} \right)^{\frac{1}{3}} \\ \xi_0 &= 1 - \frac{7}{12}\mu \\ \xi_{n+1} &= f(\xi_n) \end{aligned}$$

3 result

$$\mu = 0.008, L_3 = 1.0033333053755920$$

The first crossing with Poincare section $y=0$, the point is:

$$\begin{pmatrix} 0.59826229280258836 \\ 2.9575936557792854^{-14} \\ -0.26780932584763739 \\ 0.80590085785342314 \end{pmatrix}$$

The unstable manifold of L_3 , with initial condition: $L_3 - 10^{-6} \cdot \vec{v}$.

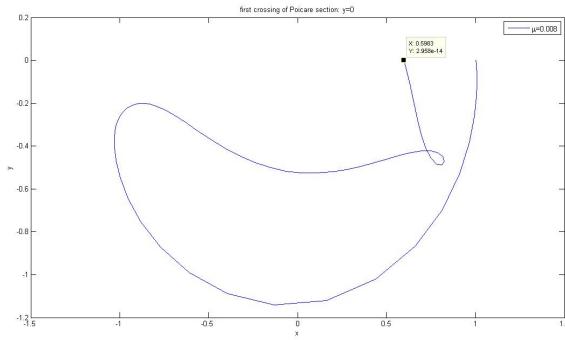


Figure 2: $\mu = 0.008$, 1st-crossing

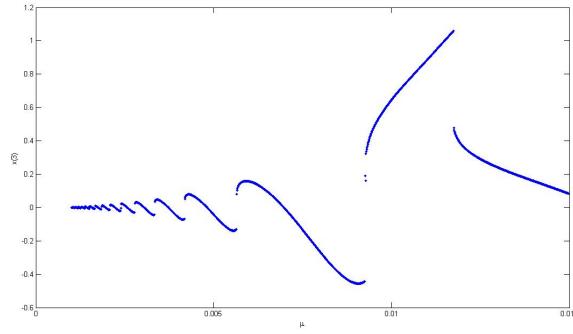


Figure 3: x' V.s μ in the interval $(0.001, 0.015)$

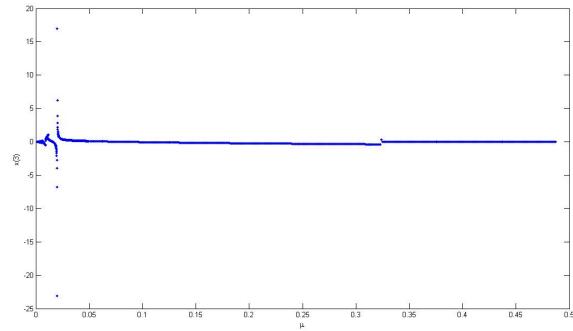


Figure 4: x' V.s μ in the interval $(0.001, 0.5)$

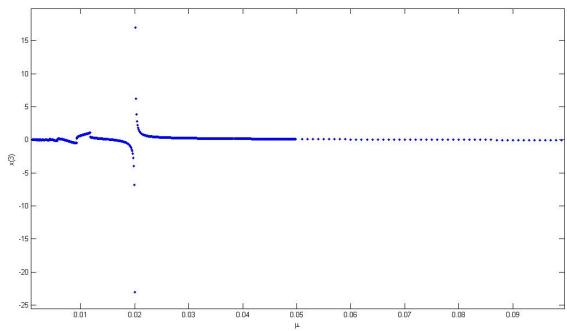


Figure 5: x' V.s μ in the interval $(0.001, 0.1)$

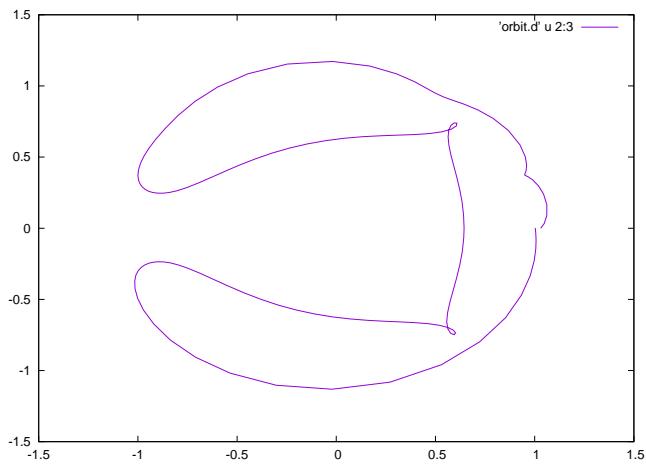


Figure 6: $\mu = 6.96499999^{-3}$, 2nd crossing

4 Code

Main function:

```
|| implicit real*8 (a-h,o-z)
|| parameter (n=4)
|| common/param/xmu
|| dimension x(n),oa(n,n),rr(n),ri(n),vr(n,n),vi(n,n)
|| open(10,file='orbit.d',status='unknown')
|| open(11,file='relation.d',status='unknown')
|| ti=0
|| T=0.3138977039438897d01
|| tmax=20.d0*T
|| np=10
|| p=-1
|| write(*,*) 'xincmu:'
|| read(*,*) xincmu
|| write(*,*) 'xmuLow:'
|| read(*,*) xmuLow
|| write(*,*) 'xmuUpp:'
|| read(*,*) xmuUpp
|| m=(xmuUpp-xmuLow)/xincmu-2
|| write (*,*) 'm:', m
|| do j=1,m
||   xmu=xmuLow+j*xincmu
||   call peq(xmu,xl1,xl2,xl3,cl1,cl2,cl3)
||   call jacobiA(xmu,xl3,n,oa)
||   call vapvep(oa,n,rr,ri,vr,vi)
||   x(1)=xl3
||   x(2)=0.d0
||   x(3)=0.d0
||   x(4)=0.d0
||   idir=dsign(1.d0,rr(3))
||   x(1)=p*vr(1,3)*1.d-6+x(1)
||   x(2)=p*vr(2,3)*1.d-6+x(2)
||   x(3)=p*vr(3,3)*1.d-6+x(3)
||   x(4)=p*vr(4,3)*1.d-6+x(4)
||   call poinci(n,x,tfinal,idir)
||   write(11,*) xmu,x(1),x(3)
|| enddo
|| end

C ****
C
```

```

C ****
C
SUBROUTINE POINC1(n,YI,tfinal,idirorig)
IMPLICIT REAL*8 (A-H,O-Z)
DIMENSION YI(n),YF(n),DGG(n),F(n)
icont=0
idir=idirorig
ti=0.d0
C DETERMINATION OF THE FIRST PASSAGE OF THE ORBIT THROUGH
y=0
CALL SECCIO(YI,GG,DGG)
IF(DABS(GG).LT.1.D-9)GG=0.d0
GA=GG
hab=.1e-16
hre=.1e-16
pabs=dlog10(hab)
prel=dlog10(hre)
istep=1
pas=5d0
ht=0.d0
t=ti
1   tmax=t+idir*pas
CALL taylor_f77_eq_rtbp_(t,yi,idir,istep,pabs,prel,
& tmax,ht,iordre,ifl)
CALL SECCIO(YI,GG,DGG)
IF(GG*GA.LT.0.D0)go to 22
write(10,*)t,(yi(ii),ii=1,4)
GA=GG
GO TO 1
C
C REFINEMENT OF THE INTERSECTION POINT YF(*) USING NEWTON
METHOD
C TO GET A ZERO OF THE FUNCTION GG (SEE SUBROUTINE SECCIO)
C
22   continue
icont=icont+1
if (icont.gt.20)then
    write(*,*)"problems finding the section"
    return
endif
CALL FIELD(T,YI,N,F)
P=0.D0
DO 3 I=1,N
3   P=P+F(I)*DGG(I)
H=-GG/P
if (h.ge.0.d0)idir=1

```

```

        if (h.lt.0.d0) idir=-1
        tmax=t+h

        CALL taylor_f77_eq_rtbp_(t,yi,idir,istep,pabs,prel,
& tmax,ht,iordre,ifl)
        CALL SECCIO(YI,GG,DGG)
        IF(DABS(GG).GT.1.D-13) GO TO 22
        DO 4 I=1,N
        YF(I)=YI(I)
        tfinal=t
        write(10,*)t,(yf(ii),ii=1,4)
        return
        end

C ****
C
C *
C      THE SURFACE g OF SECTION, IN THIS CASE
C          INPUT PARAMETERS:
C          Y(*)      POINT
C          OUTPUT PARAMETERS:
C          GG         FUNCTION THAT EQUATED TO 0 GIVES THE
C          SURFACE OF
C          SECTION
C          DGG(*)     GRADIENT OF FUNCTION GG
C

C *
C ****

SUBROUTINE SECCIO(Y,GG,DGG)
IMPLICIT REAL*8(A-H,O-Z)
DIMENSION Y(4),DGG(4)
GG=Y(2)
DO 1 I=1,4
1 DGG(I)=0.D0
DGG(2)=1.d0
RETURN
END

C ****

```

```

C
C      EQS OF MOTION IN synodical VARIABLES
C      X           TIME
C      Y(*)        POINT (Y(1),Y(2),...,Y(n))
C      NEQ         NUMBER OF EQUATIONS
C      OUTPUT PARAMETERS:
C      F(*)        VECTOR FIELD
C
C
C ****
C
C      subroutine field(t,x,neq,f)
C      implicit real*8 (a-h,o-z)
C      common/param/xmu
C      dimension x(neq),f(neq)
C      umu=1.-xmu
C      d1=x1-xmu
C      d2=x1+umu
C      r12=d1*d1+x2*x2
C      r22=d2*d2+x2*x2
C      r0=sqrt(r12)
C      r1=sqrt(r22)
C      r032=r12*r0
C      r132=r22*r1
C      omex=x1-(umu*(-xmu+x1)/r032)-(xmu*(x1+umu)/r132)
C      omey=x2*(1.-(umu/r032)-(xmu/r132))
C      f(1)=x(3)
C      f(2)=x(4)
C      f(3)=2*x(4)+omex
C      f(4)=-2*x(3)+omey
C      return
C      end
C
C ****
C
C ****
C
C      subroutine jacobiA(xmu,x,n,oa)
C      implicit real*8(a-h,o-z)
C      dimension oa(n,n)
C      umu=1.d0-xmu
C      x1=x
C      xd1=dabs(x1-xmu)
C      xd2=dabs(x1+umu)
C      r032=xd1**3

```

```

r132=xd2**3
r052=xd1**5
r152=xd2**5
xa=3*umu*(-xmu+x1)/r052
xb=3*xmu*(x1+umu)/r152
omexx=1.d0-umu/r032+xa*(-xmu+x1)-xmu/r132+xb*(x1+umu
    )
omeyy=1.d0-umu/r032-xmu/r132
oa(1,1)=0.d0
oa(1,2)=0.d0
oa(1,3)=1.d0
oa(1,4)=0.d0
oa(2,1)=0.d0
oa(2,2)=0.d0
oa(2,3)=0.d0
oa(2,4)=1.d0
oa(3,1)=omexx
oa(3,2)=0.d0
oa(3,3)=0.d0
oa(3,4)=2.d0
oa(4,1)=0.d0
oa(4,2)=omeyy
oa(4,3)=-2.d0
oa(4,4)=0.d0
return
end

```

5 optional part—some particular cases

5.1 The curve (μ, x)

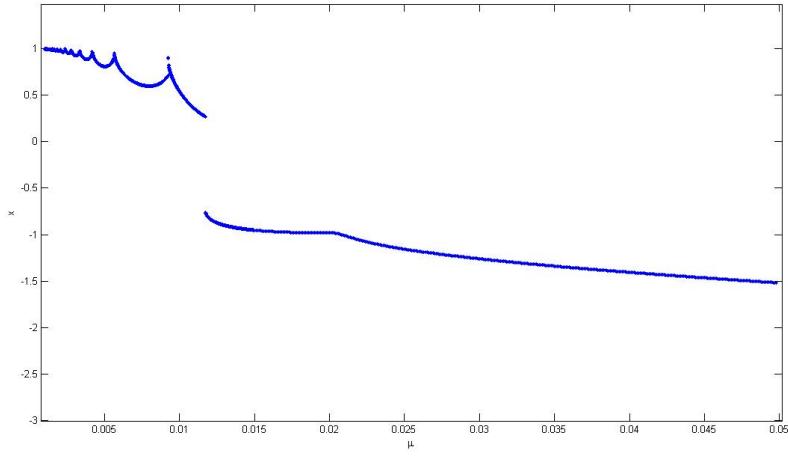


Figure 7: x V.s μ in the interval $(0.001, 0.015)$

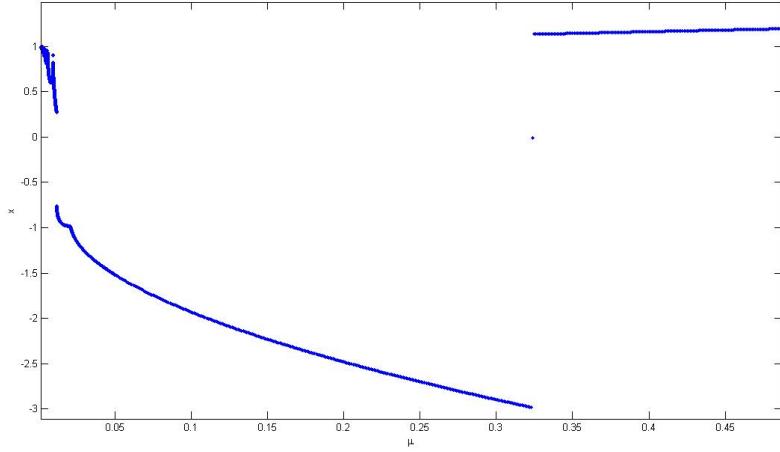


Figure 8: x V.s μ in the interval $(0.001, 0.5)$

5.2 Some fixed μ and comments

From the relationship between μ and x , i.e. figures 7 and 8, we can see in interval $\mu \in (0.001, 0.5)$, it is separated into three parts by $\mu \approx 0.01174$ and $\mu \approx 0.3245$. In the first part, the value of x is around 1, it means the first crossing is closed to the equilibrium point L_3 , which is showed in figures

9,?? and 11. The overall trend of first crossing point goes farther away from the L_3 as μ increasing, despite the oscillations at the beginning. When μ is bigger than 0.01174, it reaches the second part, where the first crossing is around -1 . We can see the examples from figures 12, 13 and 14. In the third part, the first crossing is around 1 again.

From the relationship between μ and x' , i.e. figures 3 and 4, we can see when μ is smaller than 0.01, there are several jumps in the x' . i.e. in one point(the first crossing), there are two different x' , which brings a small loop. And this loop is symmetric with respect to x-axis, as showed in figure 10b($\mu = 0.00567$, 2nd crossing). Besides, we can see, there is a singular point $\mu \approx 0.02$. It represents that the first-crossing collides with m_2 .

For the μ is equal to 0.0056, 0.00567, 0.0057 respectively.

For the μ is equal to 0.01179, 0.0159375, 0.083 respectively.

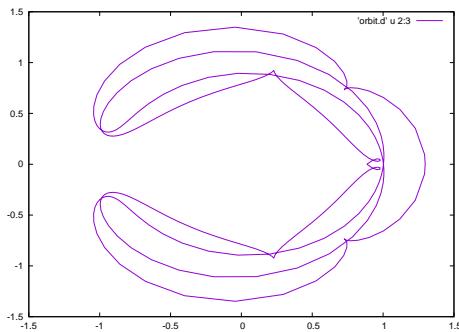


Figure 9: $\mu = 0.0056$

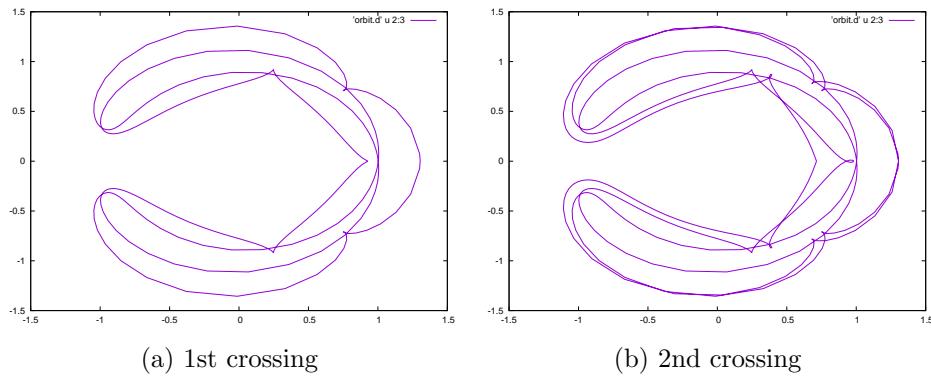


Figure 10: $\mu = 0.00567$

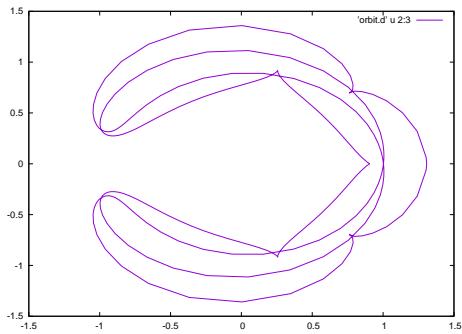


Figure 11: $\mu = 0.0057$

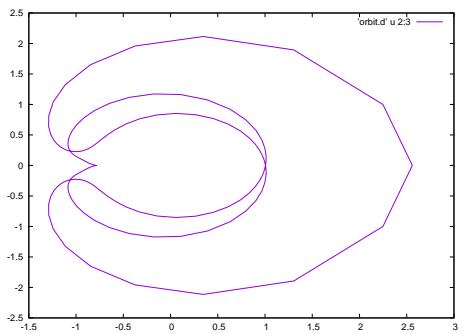


Figure 12: $\mu = 0.01179$

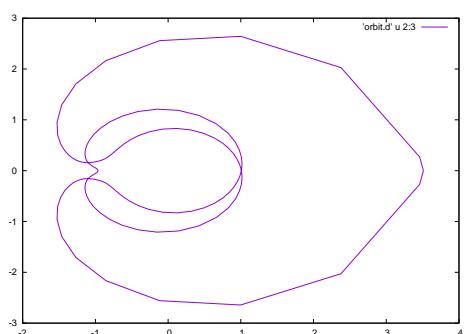


Figure 13: $\mu = 0.0159375$

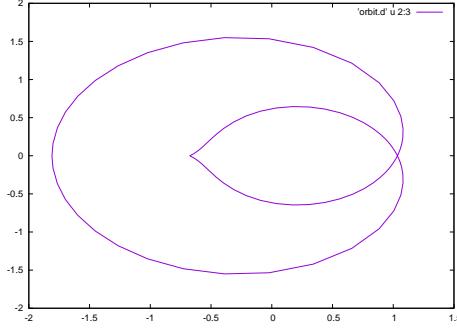


Figure 14: $\mu = 0.083$

6 Code

Main function:

```

|| implicit real*8 (a-h,o-z)
|| parameter (n=4)
|| common/param/xmu
|| dimension x(n),oa(n,n),rr(n),ri(n),vr(n,n),vi(n,n)
|| open(10,file='orbit.d',status='unknown')
|| open(11,file='relation.d',status='unknown')
|| ti=0
|| T=0.3138977039438897d01
|| tmax=20.d0*T
|| np=10
|| p=-1
|| write(*,*) 'mu: '
|| read(*,*) xmu

call peq(xmu,xl1,xl2,xl3,cl1,cl2,cl3)
call jacobIA(xmu,xl3,n,oa)
call vapvep(oa,n,rr,ri,vr,vi)
do k=1,2
  p=-1*p
  write(*,*)"p:", p
  do j=3,4
    x(1)=xl3
    x(2)=0.d0
    x(3)=0.d0
    x(4)=0.d0
    idir=dsign(1.d0,rr(j))
    write(*,*)"idir:", idir
    x(1)=p*vr(1,j)*1.d-6+x(1)
  end do
end do

```

```

x(2)=p*vr(2,j)*1.d-6+x(2)
x(3)=p*vr(3,j)*1.d-6+x(3)
x(4)=p*vr(4,j)*1.d-6+x(4)
write(*,*)'Initial t:'
write(*,*)ti
write(*,*)'Initial cond:'
write(*,*)(x(ii),ii=1,n)
write(*,*)'m times crossing'
read(*,*) m
do i=1,m
    call poinc1(n,x,tfinal,idir)
    t=tfinal+t
    write(*,*)'t:'
    write(*,*)t
enddo
write(10,90)
format()
enddo
enddo
end

```

90