

Longitudinal matching (Section 5.4, Figure 5.5)

Volker Ziemann, 211108, CC-BY-SA-4.0

This is essentially the same simulation as `LargeAmplitudeOscillations.mlx`, only the initial distribution is changed and the integration time increased to $100 \cdot T_s$. You can observe how the particles initially inside a square box eventually fill a circular region that essentially looks unchanged henceforth.

As before, the simulation is based on `pendulum_tracker()`, but changes the input distribution and the duration of the integration and increases the number of particles N .

```
clear all; close all
N=500;           % number of particles
Omegas=0.25;     % synchrotron frequency
Ts=2*pi/Omegas;  % the time for one small-amplitude oscillation
```

Here I set `dt` to `dt=Ts*100`.

```
dt=100*Ts;           % integration time=duration
```

and define the input distribution to be centered and wider in the horizontal (phase) direction.

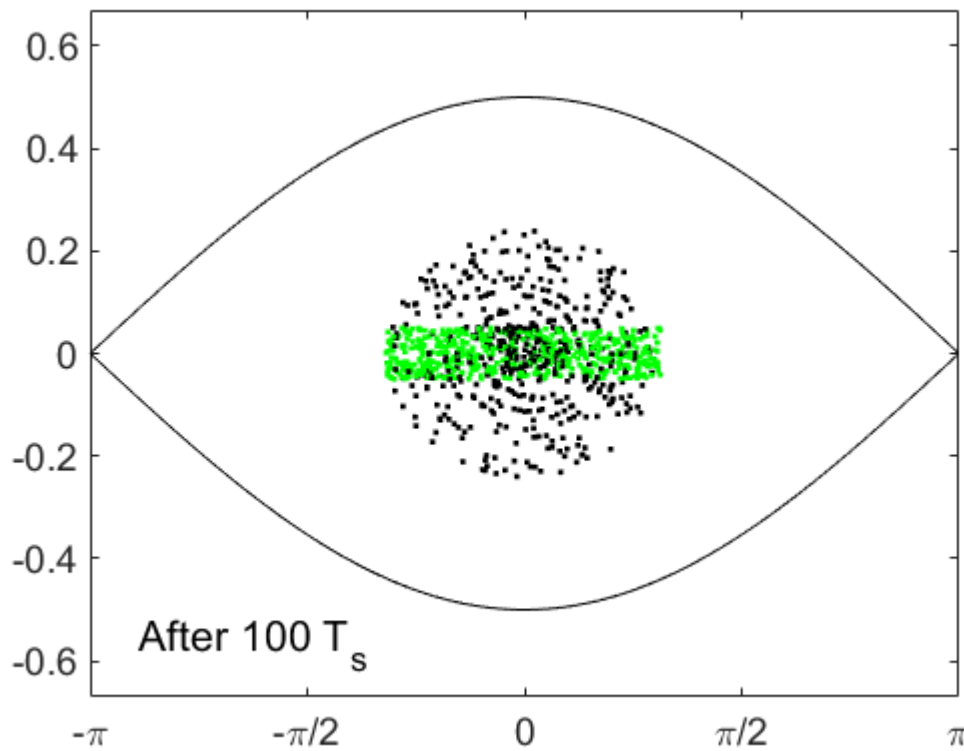
```
x0=[0.0,0.0];      % center of initial (green) distribution
dx=[2.0,0.1];      % spread of initial (green) distribution
```

Like before, we define the range of phases, plot the separatrix, and annotate the axes

```
phi=-pi:0.01:pi;   % range of angles
separatrix=2*Omegas*cos(0.5*phi); % plot the separatrix
hold off; plot(phi,separatrix,'k',phi,-separatrix,'k')
axis([-pi,pi,-0.67,0.67]); hold on; pause(0.001)
set(gca,'xtick',[-pi,-pi/2,0,pi/2,pi],'fontsize',14, ...
    'xticklabels',{'-\pi','-\pi/2','0','\pi/2','\pi'}) % xlabels
text(-2.8,-0.57,['After ',num2str(dt/Ts),' T_s'],'fontsize',16) % duration
```

and loop over the particles, where we make a green dot at the starting coordinates and a black dot at the final coordinates.

```
for k=1:N % loop over particles
    x=x0+(rand(1,2)-0.5).*dx;
    plot(x(1),x(2),'g. ');
    x1=pendulumtracker(x,Omegas,dt);
    plot(x1(1),x1(2),'k. ');
    pause(0.001)
end
```



And now you can go back to set $dt=3 \cdot T_s$ in order to generate the plot on the left-hand side in Figure 5.5 in the book.

That all, folks!

Appendix

The function `pendulumtracker()` receives the phase-space coordinates x at the start, the small-amplitude synchrotron frequency ω , and the integration time dt as input and returns the phase-space coordinates x_{out} . Internally, it integrates the equations of motion for a mathematical pendulum in closed form using Jacobi elliptic functions. This is much faster than numerically integration, especially for extremely large times, such as thousands or even millions of synchrotron periods. The coding closely follows Section 5.4, especially Equations 5.50 to 5.54.

```
function xout=pendulumtracker(x,omega,dt)
k2=(0.5*x(2)/omega)^2+sin(0.5*x(1))^2;           % just after eq. 5.45
k=sqrt(k2);
if (x(1)>pi) x(1)=x(1)-2*pi; end                 % map back into range [-pi,pi]
if (x(1)<-pi) x(1)=x(1)+2*pi; end
s=1; if (x(1)<0) s=-s; x(1)=-x(1); end           % keep track of quadrant
s1=1; if (x(2)<0) s1=-s1; end
if (k>1) % outside the separatrix
    kelf=ellipke(1/k2);                          % faster
    trev=2*kelf/(k*omega);
    t0=mod(dt,trev);
    tmp=s1*k*omega*t0+s*ellipticF(0.5*x(1),1/k2);
```

```

    % tmp=s1*k*omega*t0+s*elliptic12(0.5*x(1),1/k2);      % faster
    [sn,cn,dn]=ellipj(tmp,1/k2);
    if (abs(tmp) > kelf) sn=-sn; end
    xout(1)=2*asin(sn);                                  % eq. 5.52
    xout(2)=2*s1*omega*k*dn;
else            % inside the separatrix
    trev=4*ellipke(k2)/omega;                            % faster
    t0=mod(dt,trev);
    z0=asin(min(1,sin(0.5*x(1))/k));
    tmp=s1*omega*t0+s*ellipticF(z0,k2);
    % tmp=s1*omega*t0+s*elliptic12(z0,k2);              % faster
    [sn,cn,dn]=ellipj(tmp,k2);
    xout(1)=2*asin(k*sn);                                % eq. 5.54
    xout(2)=2*s1*omega*k*cn;
end
end

```

This version uses the built-in function `ellipticF()`, which is rather slow. They should be replaced by `elliptic12()` from the package `elliptic` from <https://github.com/moiseevigor/elliptic>, which speeds up the simulation dramatically.