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## Homework IV

In this project I deal with the continuous Hopfield Model. In the lecture and in this project we see that there is huge similarities between binary and continuous Hopfield network. Therefore in this project I look at the first project which I already done.

In this project it is given that there will be two neurons model with  $T = \begin{bmatrix} 0 & 1 \\ 1 & 0 \end{bmatrix}$ . Also non-linear function is given us and we calculate the inverse of this non-linear function.

In the code I find the equilibrium points, trajectories and both 2D and 3D energy contour maps. I begin the my project with the initializing the values and components that determined before the algorithm. I put the code for this part is given below.

```
r1=1;
r2=1;
c1=1;
c2=1;
lambda=2; %scaling factor
T=[0 1;1 0]; % T-matrix
func=@(x) 2/pi*atan(lambda*pi*x/2); %non-linear function
inv_func=@(x) 2/(lambda*pi)*tan(pi*x/2); %inverse of non-linear function
energy= @(x,y) -1/2*(x*y+y*x)+1/lambda*integral(inv_func,0,x)/r1+1/lambda*integral(inv_func,0,y)/r2; % energy
function
```

This part is obviously defines the parameters that chosen and the functions that will be used in the project.

After this part I find the first input and output values. First input is in 4th quadrant and it takes random values.

```
epoch=1; %number of iteration
x(epoch,:)=rand -rand; %initial input values
y(epoch,:)=func(x(epoch,:));
e(epoch)=energy(y(1),y(2));
```

From this point I begin the my main algorithm. In this part I calculated  $\delta(x)$  values,  $x$  values from new deltas and  $e$  values. Also I control the convergence of energy in here.

```
while (1) %%algorithm part
    epoch=epoch+1;
    delta=[0 0];
    delta(1,1)=r1*y(epoch-1,2)*T(2,1)-r1*c1*x(epoch-1,1);
```

```

delta(1,2)=r2*y(epoch-1,1)*T(1,2)-r2*c2*x(epoch-1,2);
x(epoch,1)=x(epoch-1,1)+delta(1,1); %%%update the x matrix
x(epoch,2)=x(epoch-1,2)+delta(1,2);
y(epoch,:)=func(x(epoch,:)); %%%update the y matrix
e(epoch)=energy(y(epoch,1),y(epoch,2));
if abs(e(epoch)-e(epoch-1))<0.0001
    break
end
end
end

```

From now on, I work on the plotting part. First of all I draw the convergence of trajectories and stable equilibrium points. I draw beginning point with blue “o” and I draw the stable equilibrium points with black “x”.

```

figure(1);
plot(y(1,1),y(1,2),'ro',y(:,1),y(:,2)); %%%plot of the trajectories
hold on;
l=length(e);
plot(y(l,1),y(l,2),'kx',y(l-1,1),y(l-1,2),'kx'); %%%plot of the stable equilibrium points

```

After plotting trajectories and equilibrium points I draw the energy contour map. For that first of all I create energyMatrix and I plot 2D and 3D contours.

```

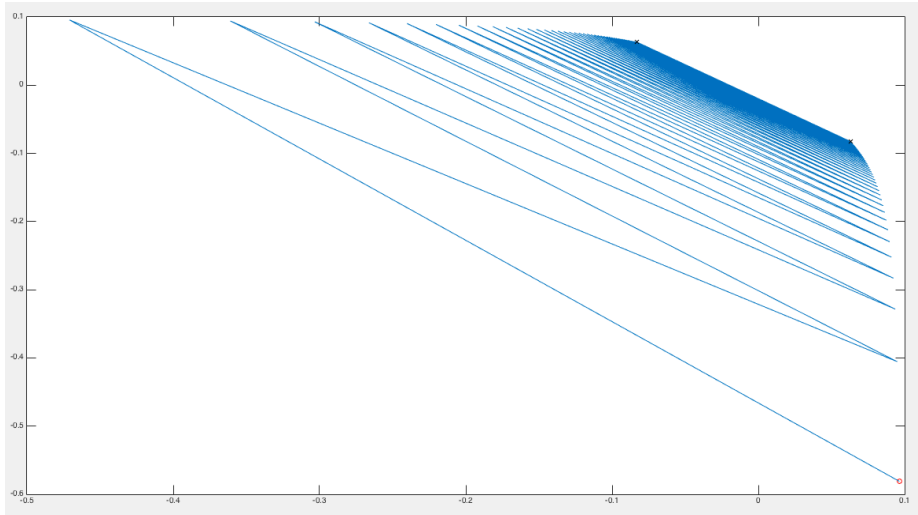
energyMatrix = zeros(100,100);
v1 = linspace(-1,1,100);
v2 = linspace(-1,1,100);
for i=1:100
    for j=1:100
        if(i==j)
            continue;
        end
        energyMatrix(i,j)=energy(v1(i),v2(j));
    end
end
v=0:0.25:1;
figure(2)
contour(v1,v2,energyMatrix,v,'ShowText','on');
figure(3)
contour3(v1,v2,energyMatrix,'ShowText','on');

```

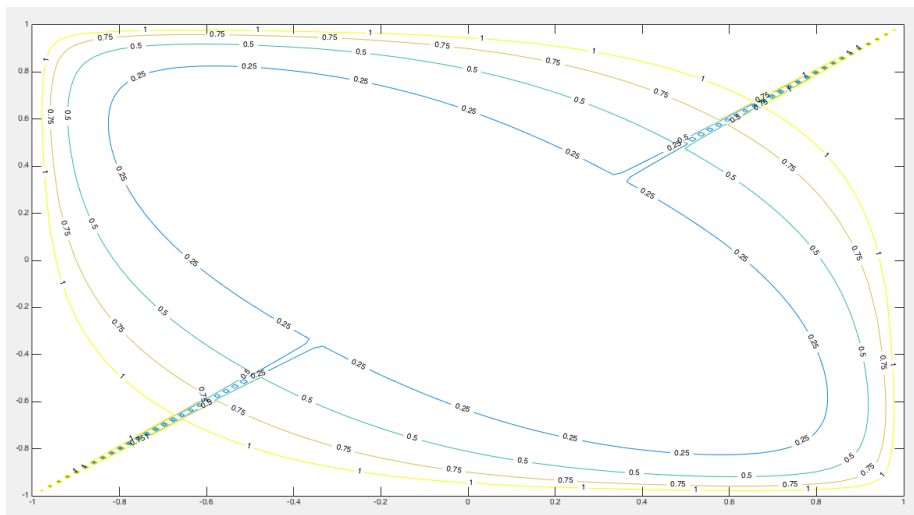
This is the whole of my code. Now I put the outputs of the system for different lambdas.

For  $\lambda=1$ ;

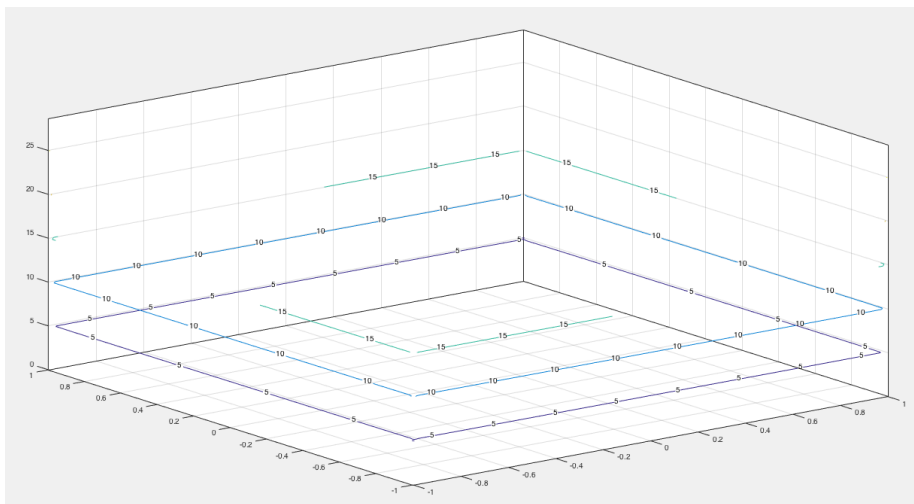
The red point is the initial random starting point. From here the trajectories go to black equilibrium points.



In this graph stable equilibrium points are  $(0.063, -0.083)$  and  $(-0.083, 0.063)$ .



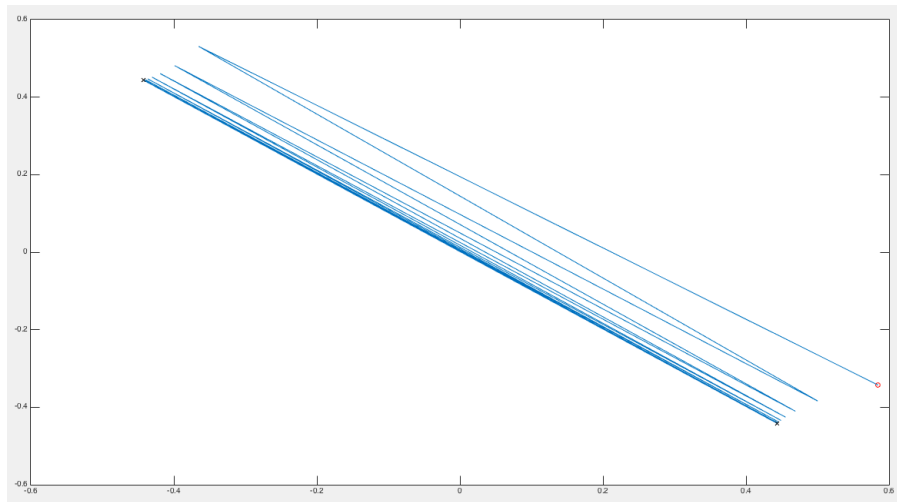
This is energy contour maps for this  $\lambda$ .



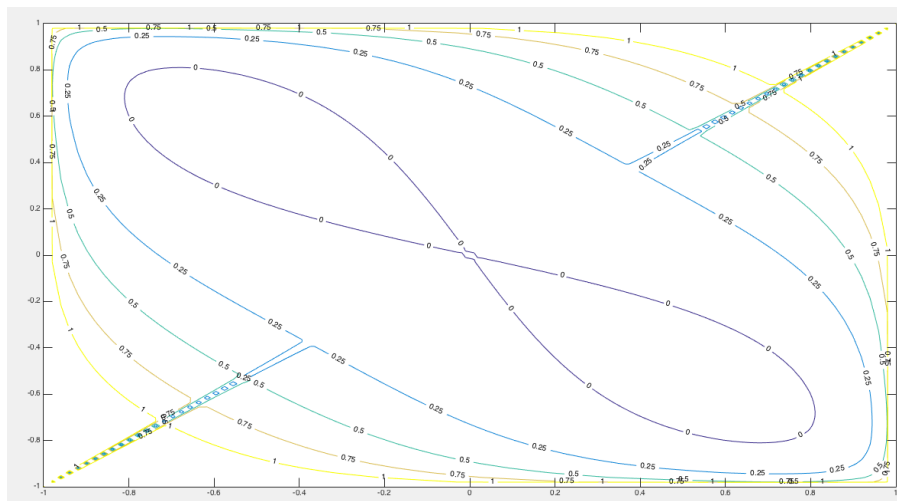
This is the 3D energy contour maps for this  $\lambda$ .

For  $\lambda=1.2$ ;

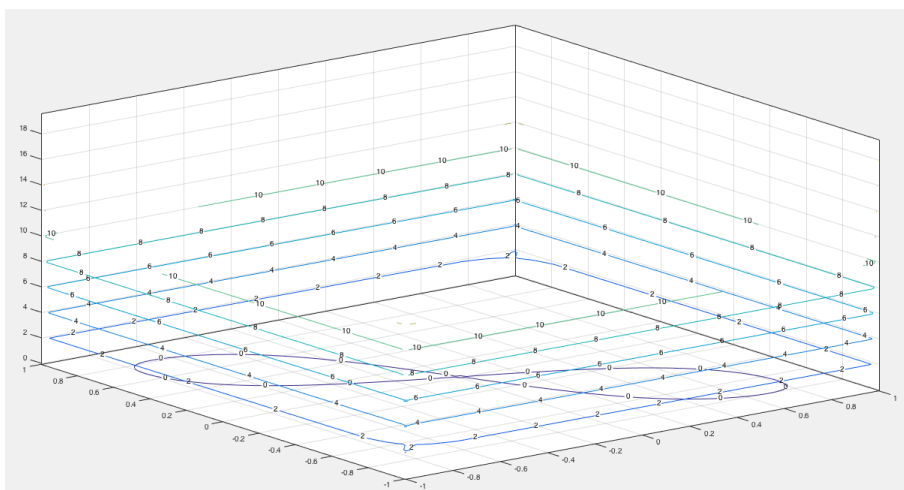
The red point is the initial random starting point. From here the trajectories go to black equilibrium points.



In this graph  
stable  
equilibrium  
points are  
(-0.442, 0.443)  
and  
(0.443, -0.442).



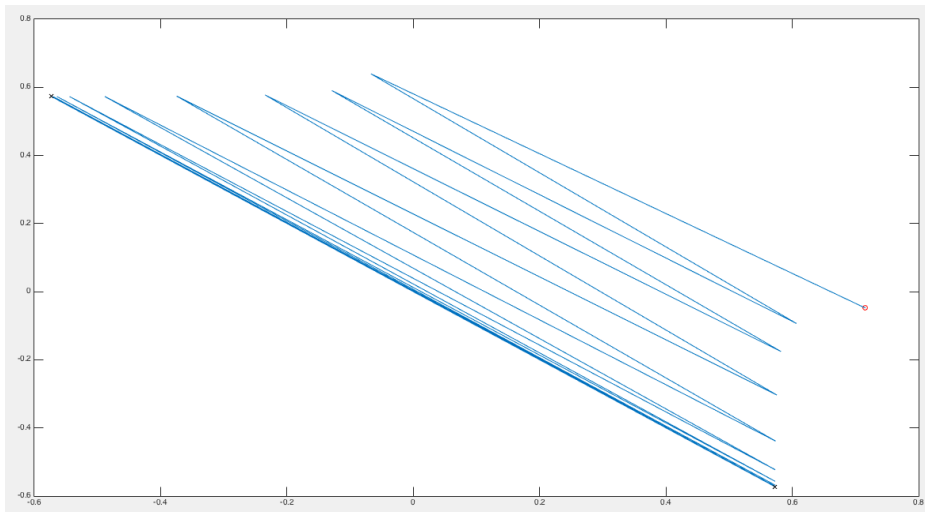
This energy  
contour for  
 $\lambda=1.2$ .



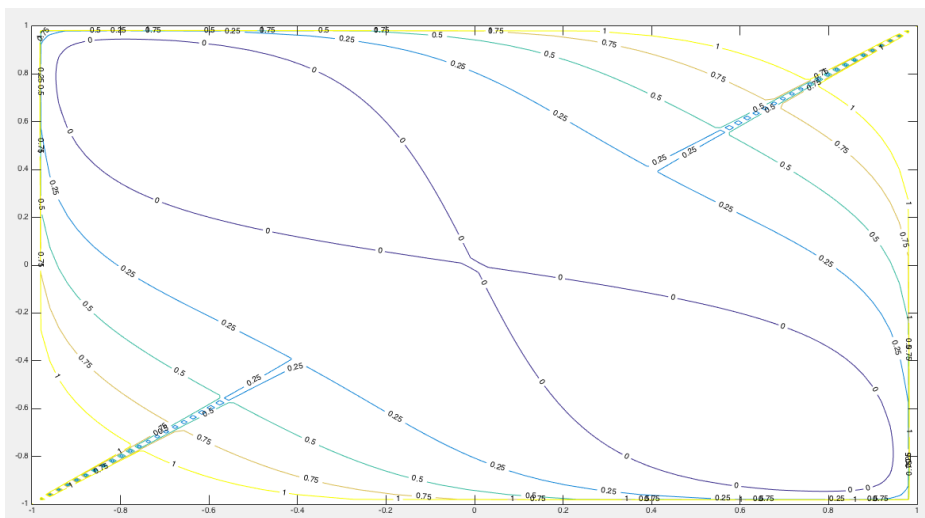
3D energy  
contour for the  
system with  
 $\lambda=1.2$

For  $\lambda=1.4$ ;

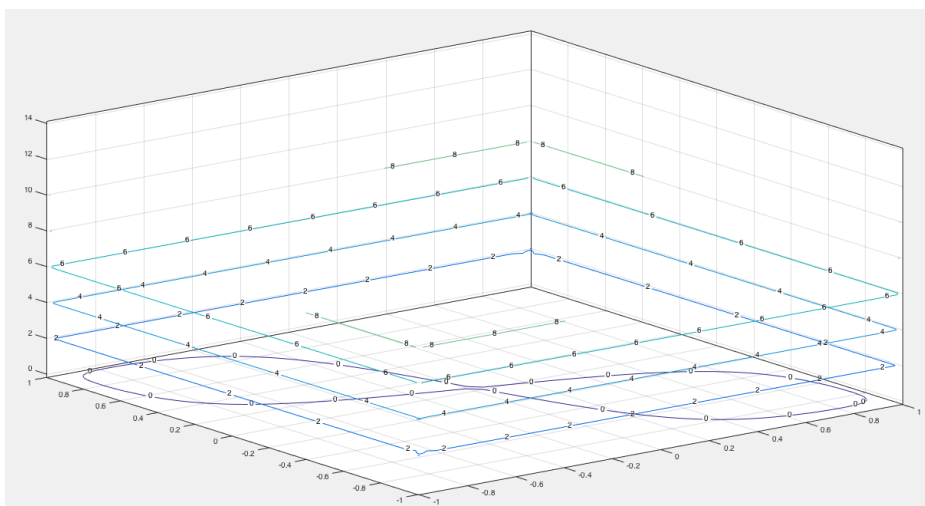
The red point is the initial random starting point. From here the trajectories go to black equilibrium points.



In here stable equilibrium points are  $(0.573, -0.573)$  and  $(-0.573, 0.573)$ .



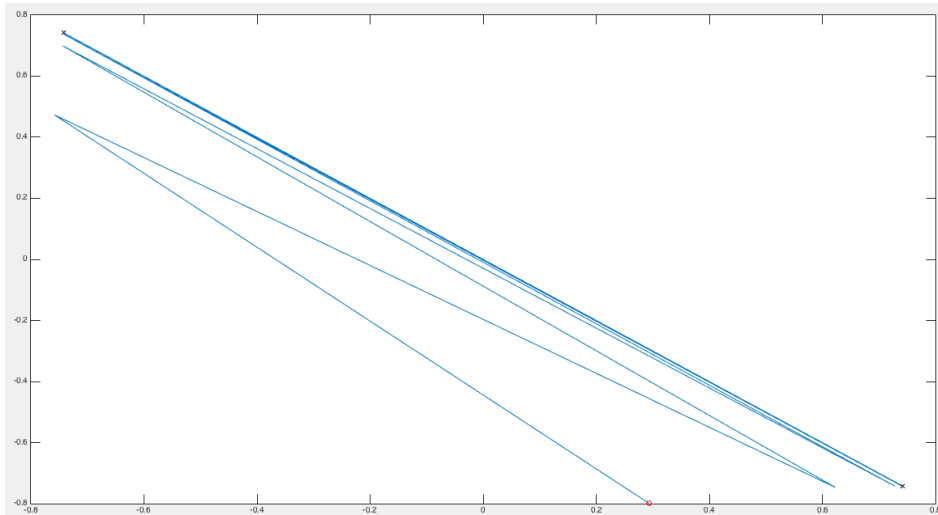
Energy contour maps for  $\lambda=1.4$ .



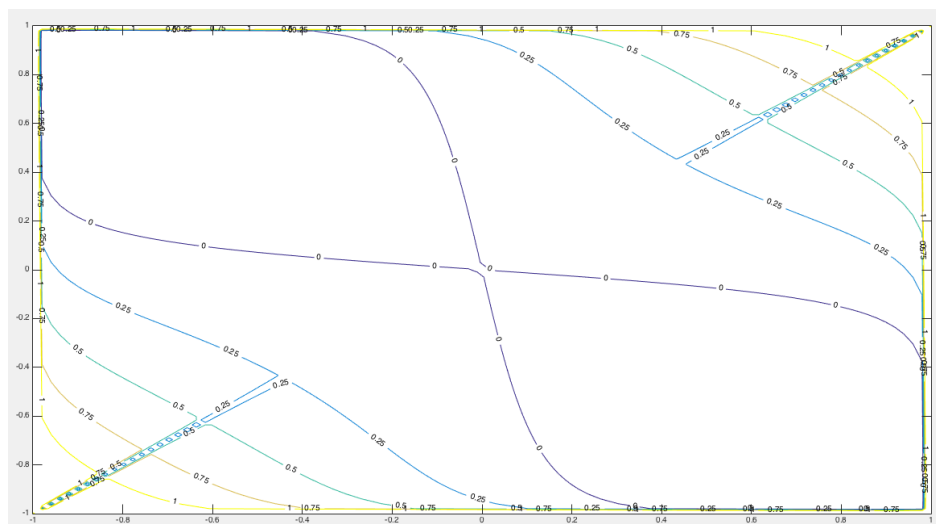
3D energy contour maps of system with  $\lambda=1.4$ .

For  $\lambda=2$ ;

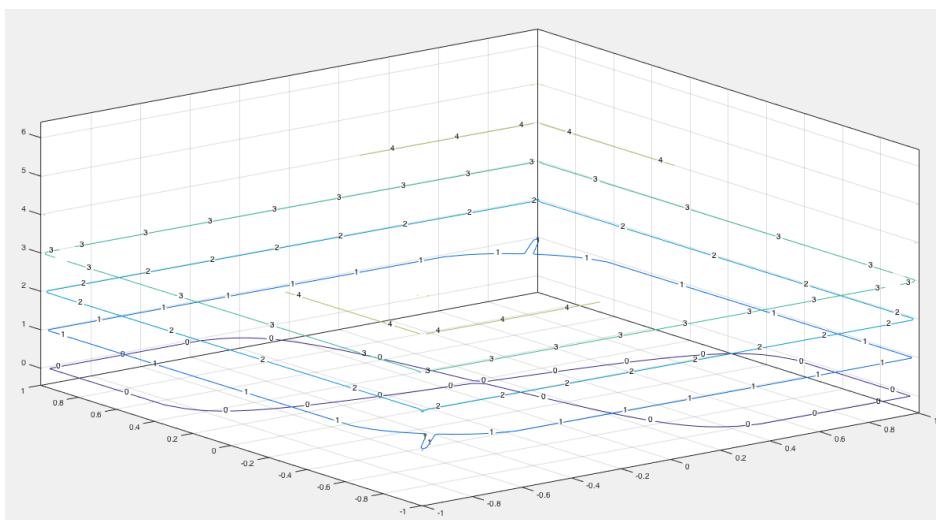
The red point is the initial random starting point. From here the trajectories go to black equilibrium points.



The stable equilibrium points of this graph is that  $(0.742, -0.742)$  and  $(-0.742, 0.742)$ .



Energy contour of system for  $\lambda=2$ .



3D energy contour of system.

We can see that when  $\lambda$  is small our equilibrium points converges to zeros however when  $\lambda$  is high our equilibrium points goes through the corners of the system.