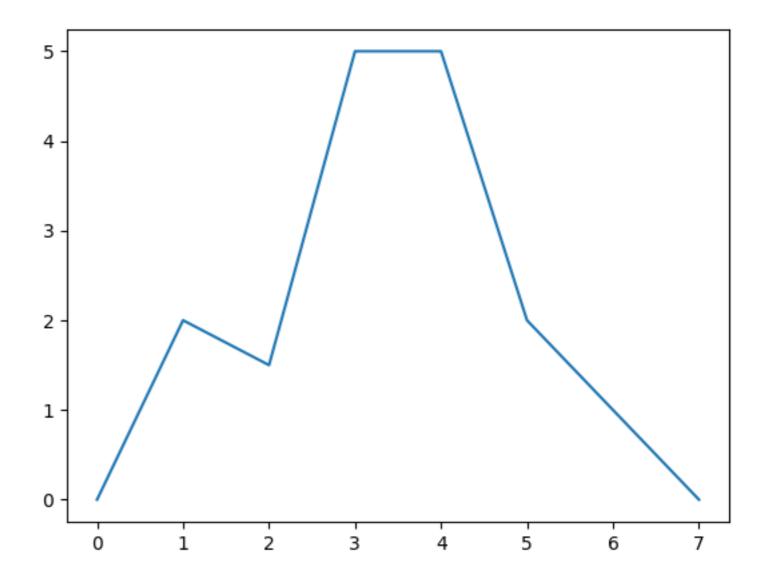
Introduction to Python

Lecture 4 Arul Lakshminarayan, 13/10/17

Basic plotting

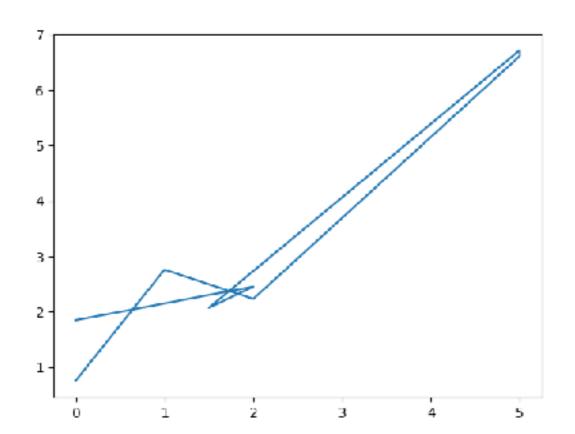
import pylab as plt import numpy as np

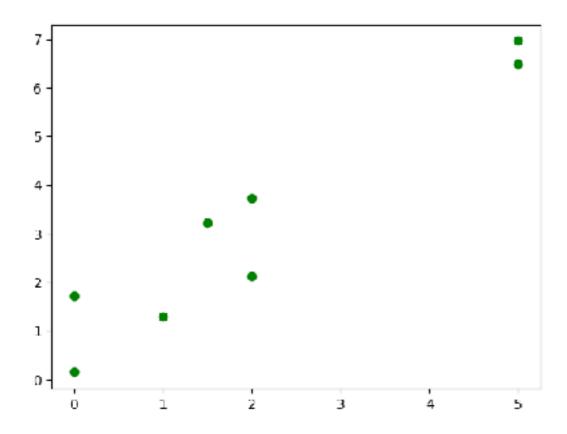
a=[.0,2.0,1.5,5,5,2,1,0] plt.plot(a) plt.show()

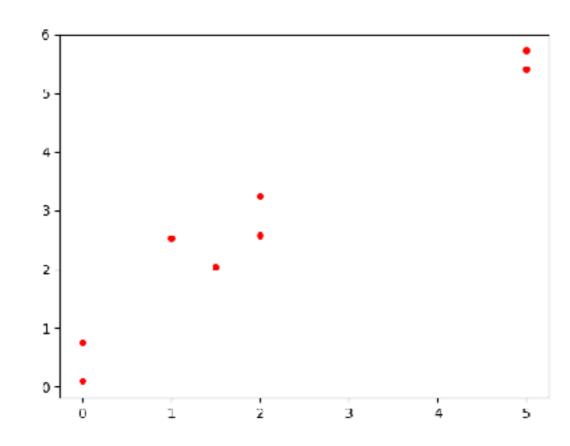


import pylab as plt import numpy as np

```
x=[.0,2.0,1.5,5,5,2,1,0]
y=x+2*np.random.random(len(x))
plt.plot(x,y)
#plt.plot(x,y,"go")
#plt.plot(x,y,"ro",markersize=4)
plt.show()
```







Example: A single spin in a magnetic field

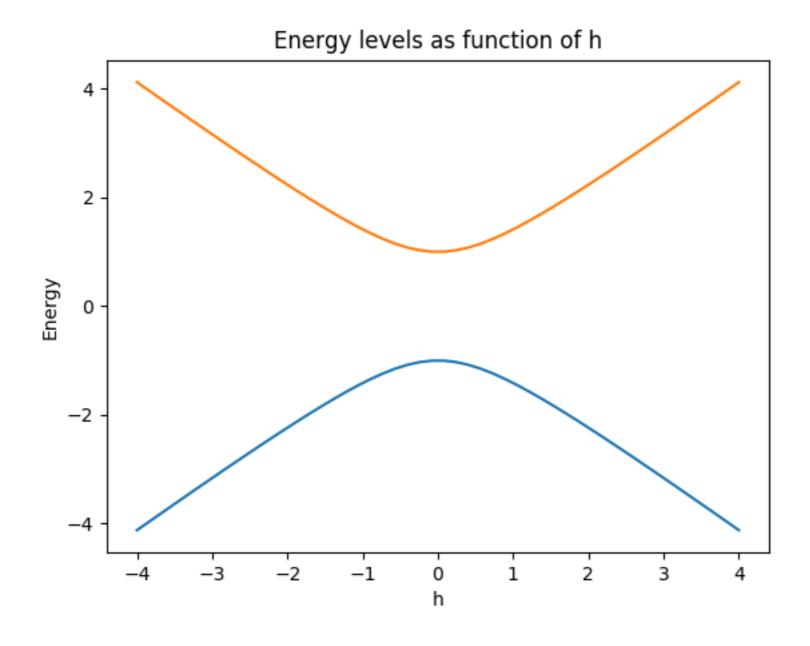
$$\sigma_z = \left(\begin{array}{cc} 1 & 0 \\ 0 & -1 \end{array}\right) \qquad \qquad \sigma_x = \left(\begin{array}{cc} 0 & 1 \\ 1 & 0 \end{array}\right)$$

$$H = \sigma_z + h \sigma_x = \begin{pmatrix} 1 & h \\ h & -1 \end{pmatrix}$$

The EIGENVALUEs of H are the energy levels

```
import numpy as np
import pylab as plt
sigmax=np.array([[0,1],[1,0]])
sigmaz=np.array([[1,0],[0,-1]])
evallist=[]
x=[]
y=[]
z=[]
for h in np.linspace(-4,4,50):
  hamil=sigmaz+h*sigmax
  eval1=np.linalg.eigvalsh(hamil)
  x.append(h)
  y.append(eval1[0])
  z.append(eval1[1])
  evallist.append([h,eval1[0],eval1[1]])
np.savetxt('evalszsx.dat',evallist,fmt="%2.5f")
plt.plot(x,y)
plt.plot(x,z)
plt.xlabel('h')
plt.ylabel('Energy')
plt.title('Energy levels as function of h')
plt.show()
```

```
Python — more evalszsx.dat — 80×24
-4.00000 -4.12311 4.12311
-3.83673 -3.96491 3.96491
-3.67347 -3.80715 3.80715
-3.51020 -3.64987 3.64987
-3.34694 -3.49314 3.49314
-3.18367 -3.33703 3.33703
-3.02041 -3.18165 3.18165
-2.85714 -3.02709 3.02709
-2.69388 -2.87350 2.87350
-2.53061 -2.72103 2.72103
-2.36735 -2.56989 2.56989
-2.20408 -2.42033 2.42033
-2.04082 -2.27265 2.27265
-1.87755 -2.12725 2.12725
-1.71429 -1.98463 1.98463
-1.55102 -1.84544 1.84544
-1.38776 -1.71052 1.71052
-1.22449 -1.58094 1.58094
-1.06122 -1.45815 1.45815
-0.89796 -1.34400 1.34400
-0.73469 -1.24088 1.24088
-0.57143 -1.15175 1.15175
-0.40816 -1.08009 1.08009
```



Doing it in 2 steps

```
!/usr/bin/env python3
# -*- coding: utf-8 -*-
Created on Fri Oct 6 11:44:00 2017
Hamiltonian: \sigma_z+ h* \sigma_x
@author: arul
11 11 11
import numpy as np
import pylab as plt
sigmax=np.array([[0,1],[1,0]])
sigmaz=np.array([[1,0],[0,-1]])
evallist=∏
for h in np.linspace(-4,4,50):
  hamil=sigmaz+h*sigmax
  eval1=np.linalg.eigvalsh(hamil)
  evallist.append([h,eval1[0],eval1[1]])
np.savetxt('evalszsx_1.dat',evallist,fmt="%2.5f")
```

Creates data in the form of 3 columns containing h, eval(0), and eval(1).

Saves the data in a file called 'evalszsx_1.dat'

Note the format "%2.5f"

loadtxt

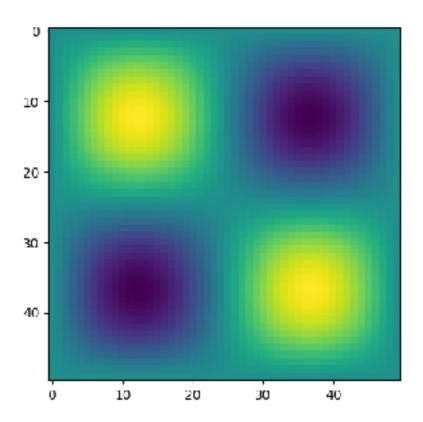
```
#!/usr/bin/env python3
# -*- coding: utf-8 -*-
Created on Fri Oct 6 11:44:00 2017
Hamiltonian: \sigma_z+ h* \sigma_x
@author: arul
import numpy as np
import pylab as plt
x,y,z=np.loadtxt('evalszsx_1.dat',unpack='True')
plt.plot(x,y)
plt.plot(x,z)
plt.xlabel('h')
plt.ylabel('Energy')
plt.title('Energy levels as function of h')
plt.show()
```

Without "unpacking"

```
#!/usr/bin/env python3
# -*- coding: utf-8 -*-
ПППП
Created on Fri Oct 6 11:44:00 2017
Hamiltonian: \sigma_z+ h* \sigma_x
@author: arul
import numpy as np
import pylab as plt
data=np.loadtxt('evalszsx_1.dat')
plt.plot(data[:,0],data[:,1])
plt.plot(data[:,0],data[:,2])
plt.xlabel('h')
plt.ylabel('Energy')
plt.title('Energy levels as function of h')
plt.show()
```

2-D density plot

```
##Plotting sin(x)*sin(y) as a density plot
##
import pylab as plt
import numpy as np
x=np.linspace(0,2*np.pi) #Defualt is 50 values
z=np.array([np.sin(xx)*np.sin(yy) for xx in x for yy in x])
z=z.reshape((50,50))
#z=∏
#for i in x:
   for j in x:
      z.append(np.sin(i)*np.sin(j))
#z=np.array(z).reshape((50,50))
plt.imshow(z)
plt.show()
```



Saving figures directly

```
import pylab as plt
import numpy as np

x=np.linspace(0,2*np.pi)

#y=x+2*np.random.random(len(x))

plt.plot(x,np.sin(x))

plt.plot(x,np.cos(x),"ro")

plt.xlabel('x')

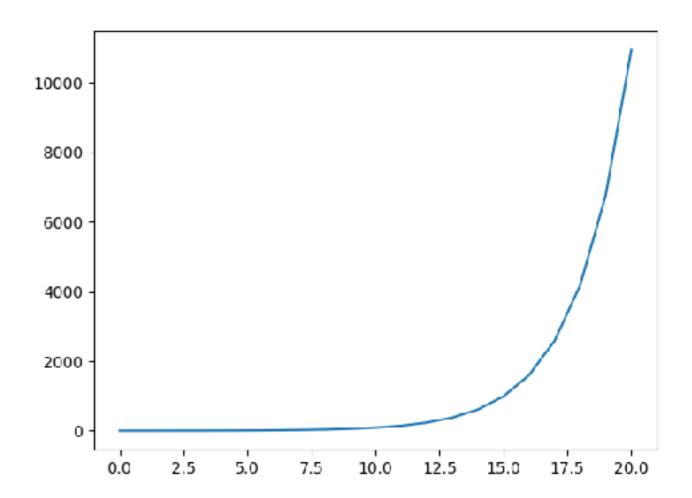
plt.ylabel('sin(x), cos(x)')

plt.savefig('foo.pdf')

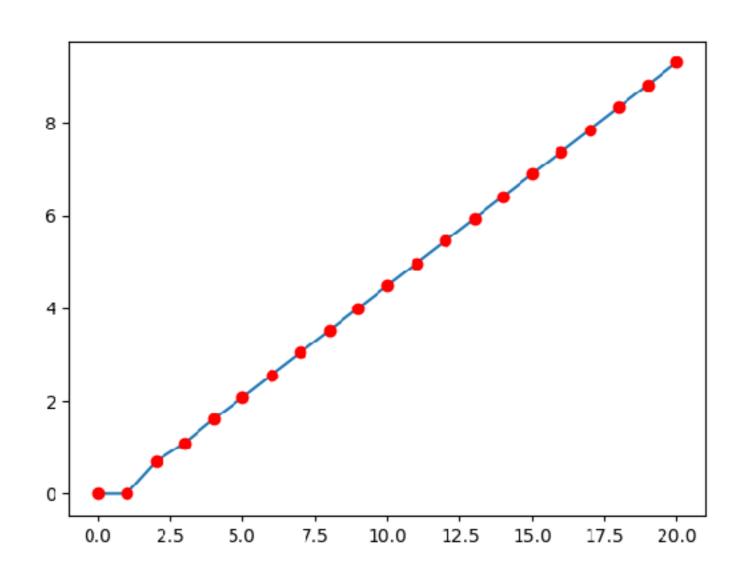
#plt.show()
```

Revisiting Fibonacci

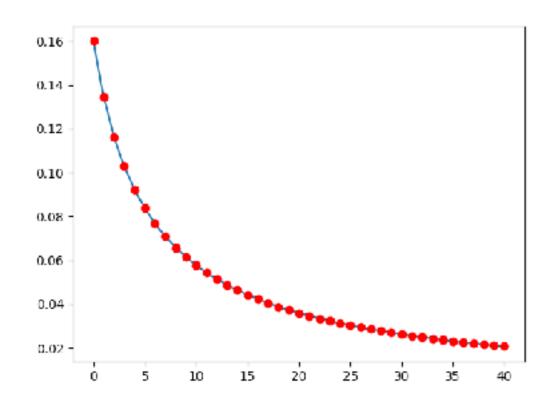
```
# Fibonacci numbers module, also plotting it
import numpy as np
import pylab as plt
def fib2(n): # return Fibonacci series up to n
  result = []
  a, b = 0, 1
  for i in range(n+1):
     result.append((i,b))
     a, b = b, a+b
  data=np.array(result)
  plt.plot(data[:,0],data[:,1])
  plt.show()
```

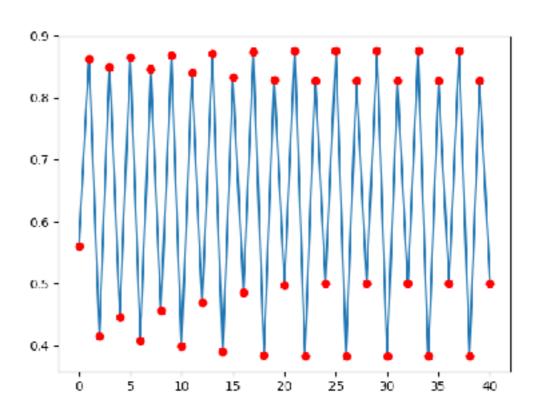


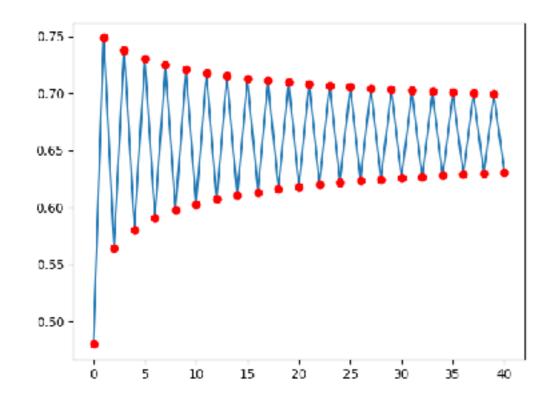
Revisiting Fibonacci: Exponential growth

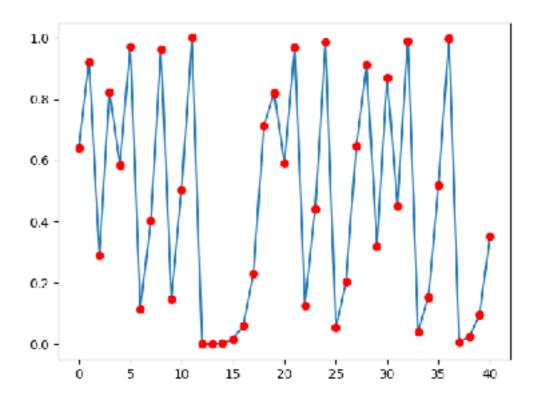


Revisiting Logistic map: r=1, 3, 3.5, 4









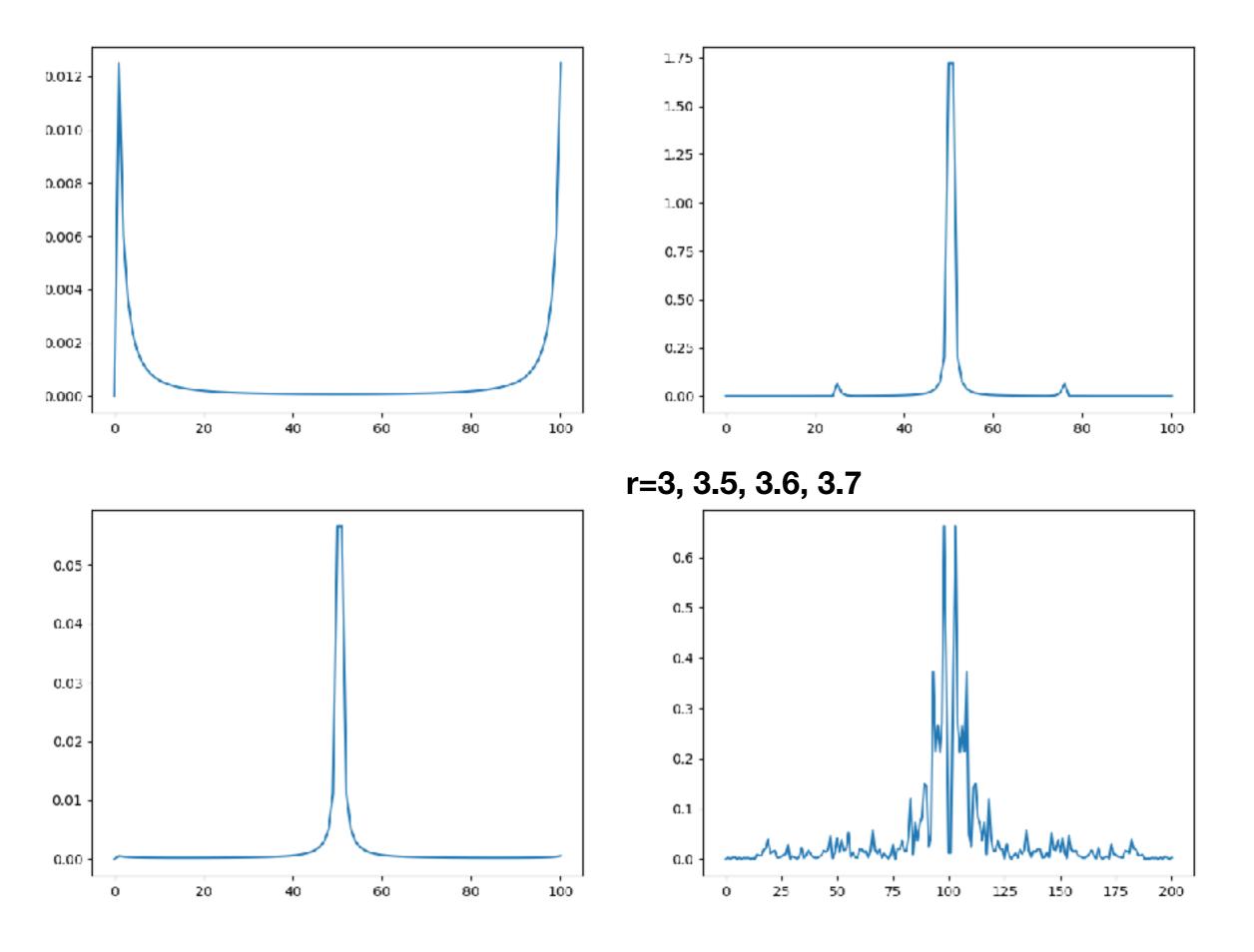
Fourier Transform DFT, and FFT

$$A_k = \sum_{m=0}^{n-1} a_m \exp\left[-2\pi i \frac{mk}{n}\right], \quad k = 0, \dots, n-1.$$

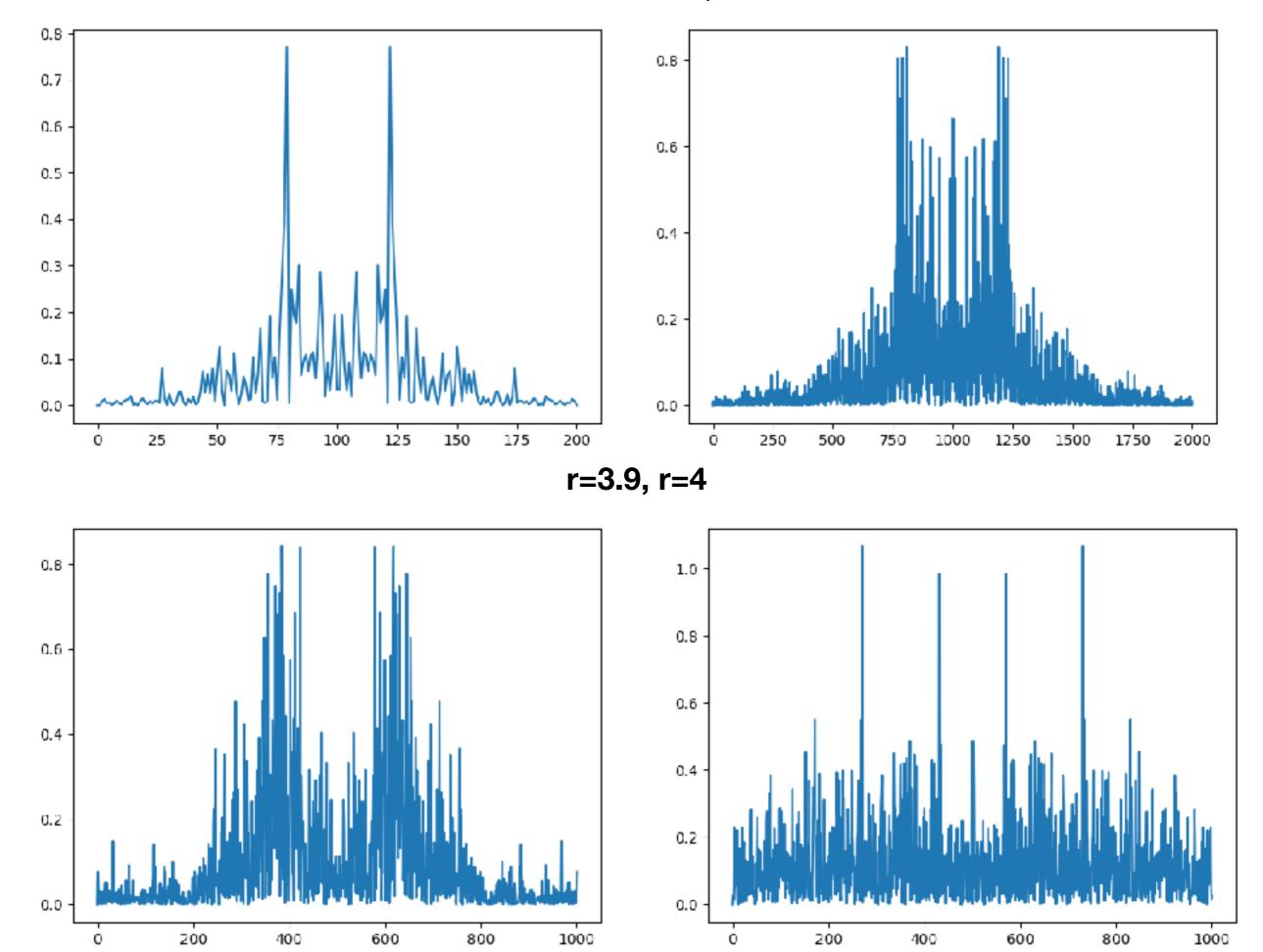
$$a_m = \frac{1}{n} \sum_{k=0}^{n-1} A_k \exp\left[2\pi i \frac{mk}{n}\right], \quad m = 0, \dots, n-1.$$

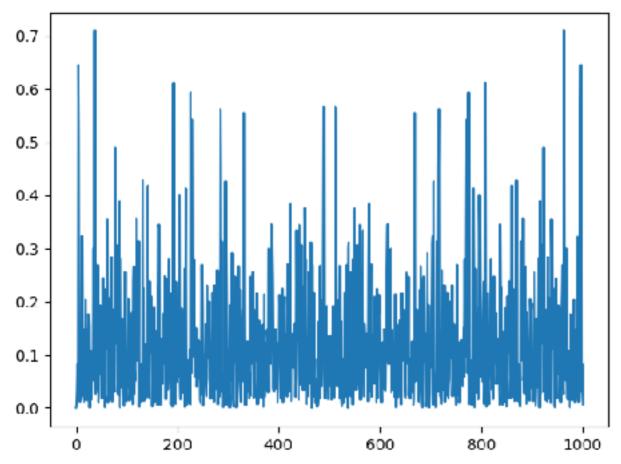
FFT of logistic map data

```
# FFT of logistic map
import numpy as np
import pylab as plt
def logistic_mapfft(r,x,n):
  xn=∏
  for i in range(n+1):
    #print(x,end=',')
    x=r^*x^*(1-x)
    xn.append(x)
  xna=np.array(xn)
  xn=xna-sum(xna)/len(xn) ##Remove "DC" component.
  xnf=np.fft.fft(xn,norm='ortho')
  plt.plot(np.abs(xnf)**2)
  #plt.plot(xn,"ro")
  plt.show()
```



r=3.8 for n=200, 2000





r=4, another orbit of length 1000

Random unif [0,1]

