1. 
$$det(A-2I)=0$$

A-2 $I = \begin{vmatrix} 0-2 & 3 \\ 4 & 4-2 \end{vmatrix}$ 
 $e(-2)(4-2)-12=0$ 
 $-42+2^2-12=0$ 
 $2=6,-2$ 

For  $A=6$ ,

 $A-6I)x=0$ 
 $\begin{bmatrix} A-6I)x=0 \\ 4 & -2 \end{bmatrix}\begin{bmatrix} x_1 \\ x_2 \end{bmatrix}=\begin{bmatrix} 0 \\ 0 \end{bmatrix}$ 
 $-6x_1+3x_2=0$ 
 $4x_1-2x_2=0$ .

solving, we get,  $x_1=\frac{x_2}{2}$ 
 $x_1=x_1\begin{bmatrix} 2 & 3 \\ 4 & 6 \end{bmatrix}\begin{bmatrix} x_1 \\ x_2 \end{bmatrix}=\begin{bmatrix} 0 \\ 0 \end{bmatrix}$ 
 $2x_1+3x_2=0$ 
 $4x_1+6x_2=0$ 
 $x=x_1\begin{bmatrix} -1 \\ 6 \end{bmatrix}$ 

(b) spectral rachus = 6.

We care about spectral rachus because it has we can impact on the convergence of matrix power sequences and series.

In nuclear agreeing, spectral rachus = 25

In nuclear agreeing, spectral rachus = 25

In nuclear agreeing, spectral rachus = 100

In nuclear agreeing, spectral rachus = 25

In nuclear agreeing Spectral ractions will be I. spectral rocki of the systems: (ii) 5.0/10.c = 0.5 (ii) 0.8/1.0 = 0.8 i) will converge most quickly and (ii) will converge most slowly. Smaller the spectral raches foster the convergence. (c) In the k-eigenvalue problem, le is a multiplication factor. Multiplication factor is the ratio of neutrons produced in one generation to the neutrons produced in the previous generation. for a self sustaining produced system the multiplication factor must be greater than or equal to one, factor

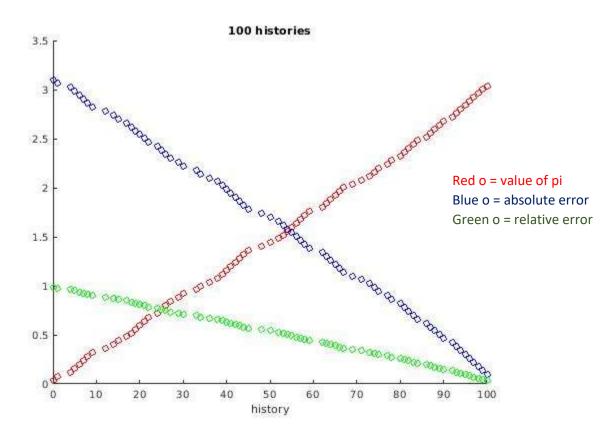
# **QUESTION 2**

#### For 100 histories:

Final value of pi=3.04

Absolute error= |pi-3.14159|=0.10159

Relative error= absolute error/3.14159 = 3.23E-2

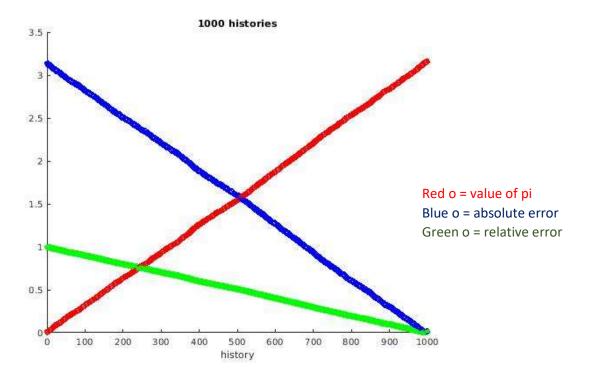


### For 1000 histories:

Final value of pi=3.1600

Absolute error= |pi-3.14159|=0.01841

Relative error= absolute error/3.14159 = 5.86E-3

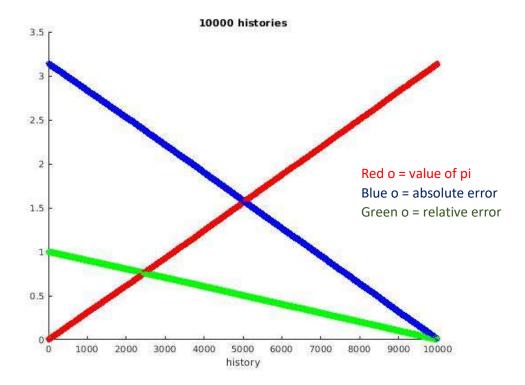


# For 10000 histories:

Final value of pi=3.1512

Absolute error= |pi-3.14159|=0.00961

Relative error= 3.059E-3



# MATIab code

```
clear
hold on;
counter = 0;
                %count thr number points falling inside a unit circle
history = 10000; %number of histories
for i=0:1:history
  rng('shuffle'); % calling for random number generation
  x = (1).*rand(); %generating random x values less than 1.0
  y = (1).*rand(); %generating random y values less than 1.0
   circle=((x*x)+(y*y))^0.5; %circle equation, calcuating the distance to the circumference
        if circle<1
                     %checking the coordinates falls within the unit circle
        counter=counter+1; %if falls within the circle, increases the points inside the circle by 1
        pi=(counter/history) * 4; %equation to calculate pi
        ab=abs(pi-3.14159); %calculating absolute error
        re=ab/3.14159; %calculating relative error
    plot(i,pi, '-ro'); %plotting value of pi in each iteration
    plot(i,ab, 'bo'); %plotting value of absolute error in each iteration
    plot(i,re, 'go'); %plotting value of relative rror in each iteration
    end
  end
pi_final=(counter/history) * 4; %final value of pi
```

Monta carlo is inherently parallel because the particle histories are independent of each other. It treats each physical process as a probabilistic process and randomly sample each process using an independent streetm of numbers. Home the particle can be smulated in parallel and sthe results can be summed up at the end of the process but the random number generator should be validated for seek parallelization. = 200 200+150+530 = 0.227 Probability of fission (f) = 51 = 530 200+150+530 = 0.602 P(inelostic scottering) = 20 . P(8) = 0.10x0.227 = 0.023 P (elastic scattering) = 180 . P(s) = 0.90x0.227 = 0.204