

# Randomness

What type of problems can we solve with the help of random numbers?

We can compute (potentially) complicated averages:

- How much my stock/option portfolio is going to be worth?
- What are my odds to win a certain competition?

## Random Number Generators

- Computers are deterministic - operations are reproducible
- How do we get random numbers out of a deterministic machine?

```
In [1]: import numpy as np
```

```
In [2]: np.random.rand(10)
```

```
Out[2]: array([0.19252561, 0.95875038, 0.34205705, 0.44060429, 0.28911005,  
              0.82233359, 0.29041589, 0.29592469, 0.69701575, 0.81690982])
```

```
In [3]: for x in range(0, 20):  
        numbers = np.random.rand(6)  
        #print(numbers)
```

They all seem random correct? Let's try to fix something called **seed** using `np.random.seed(10)`

What do you observe?

Let's see what this seed is...

## Pseudo-random Numbers

- Numbers and sequences appear random, but they are actually reproducible
- Great for algorithm developing and debugging
- How truly "random" are these numbers?

## Linear congruential generator

Given the parameters  $a$ ,  $c$ ,  $m$  and  $s$ , where  $s$  is the seed value, this algorithm will generate a sequence of pseudo-random numbers:

$$x_0 = s$$

$$x_{n+1} = (ax_n + c) \bmod(m)$$

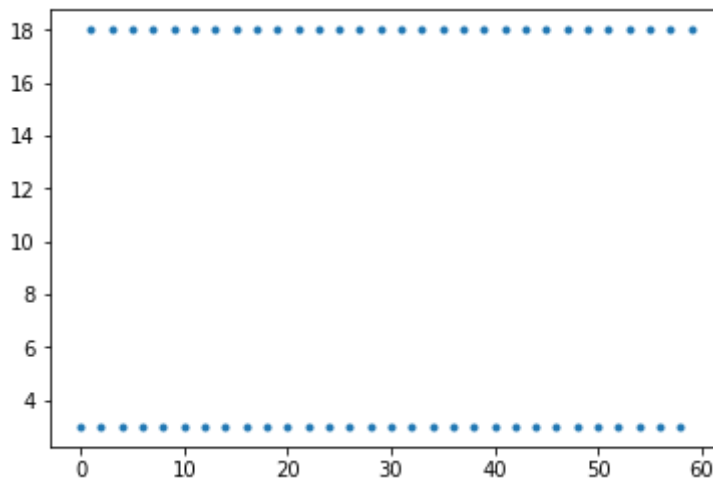
```
In [4]: s = 3 # seed
        a = 37 #10 # multiplier
        c = 2 # increment
        m = 19 # modulus
```

```
In [5]: n = 60
        x = np.zeros(n)
        x[0] = s
        for i in range(1,n):
            x[i] = (a * x[i-1] + c) % m
```

```
In [6]: import matplotlib.pyplot as plt
        %matplotlib inline
```

```
In [7]: plt.plot(x, '.')
```

```
Out[7]: [<matplotlib.lines.Line2D at 0x11227fd68>]
```



Notice there is a period, when numbers eventually start repeating. One of the advantages of the LCG is that by using appropriate choice for the parameters, we can obtain known and long periods.

Check here [https://en.wikipedia.org/wiki/Linear\\_congruential\\_generator](https://en.wikipedia.org/wiki/Linear_congruential_generator) ([https://en.wikipedia.org/wiki/Linear\\_congruential\\_generator](https://en.wikipedia.org/wiki/Linear_congruential_generator)) for a list of commonly used parameters of LCGs.

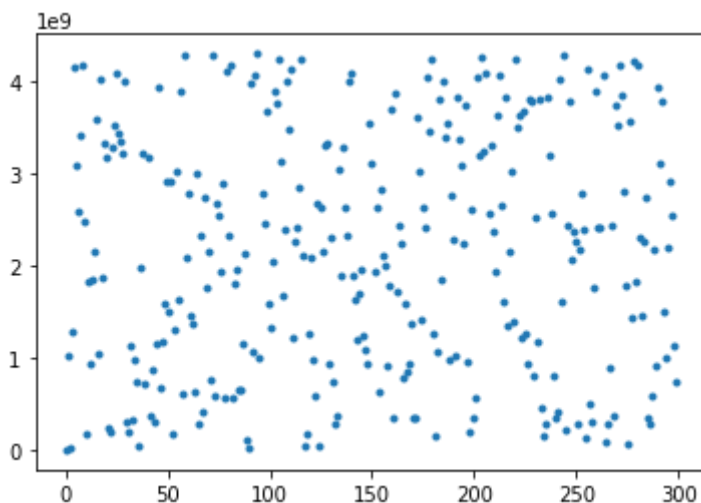
Let's try using the parameters from [Numerical recipes](https://en.wikipedia.org/wiki/Numerical_Recipes) ([https://en.wikipedia.org/wiki/Numerical\\_Recipes](https://en.wikipedia.org/wiki/Numerical_Recipes))

```
In [8]: s = 8
a = 1664525
c = 1013904223
m = 2**32

n = 300
x = np.zeros(n)
x[0] = s
for i in range(1,n):
    x[i] = (a * x[i-1] + c) % m

plt.plot(x, '.')
```

```
Out[8]: [<matplotlib.lines.Line2D at 0x11237add8>]
```



"Good" random number generators are efficient, have long periods and are portable.

## Random Variables

Think of a random variable  $X$  as a function that maps the outcome of an unpredictable (random) processes to numerical quantities.

For example:

- $X$  = the face of a bread when it falls on the ground. The random value can be the "battered" side or the "not battered" side
- $X$  = value that appears on top of dice after each roll

We don't have an exact number to represent these random processes, but we can get something that represents the **average** case. To do that, we need to know the likelihood of each individual value of  $X$ .

## Coin toss

Random variable  $X$ : result of a coin toss

In each toss, the random variable can take the values  $x_1 = 0$  (tail) and  $x_2 = 1$  (head), and each  $x_i$  has probability  $p_i = 0.5$ .

The **expected value** of a discrete random variable is defined as:

$$E(x) = \sum_{i=1}^m p_i x_i$$

Hence for a coin toss we have:

$$E(x) = 1(0.5) + 0(0.5) = 0.5$$

## Roll Dice

Random variable  $X$ : value that appears on top of the dice after each roll

In each toss, the random variable can take the values  $x_i = 1, 2, 3, \dots, 6$  and each  $x_i$  has probability  $p_i = 1/6$ .

The **expected value** of the discrete random variable is defined as:

```
In [9]: E = 0
        for i in range(6):
            E += (i+1)*(1/6)
        E
```

```
Out[9]: 3.5
```

# Monte Carlo Methods

Monte Carlo methods are algorithms that rely on repeated random sampling to approximate a desired quantity.

## Simulating a coin toss experiment

We want to find the probability of heads when tossing a coin. We know the expected value is 0.5. Using Monte Carlo with  $N$  samples (here tosses), our estimate of the expected value is:

$$E = \frac{1}{N} \sum_{i=1}^N f(x_i) = \frac{1}{N} \sum_{i=1}^N x_i$$

where  $x_i = 1$  if the toss gives head.

Let's toss a "fair" coin  $N$  times and record the results for each toss.

But first, how can we simulate one toss?

```
In [10]: toss = np.random.choice([0,1])  
print(toss)
```

```
0
```

```
In [11]: N = 30 # number of samples (tosses)
```

```
toss_list = []  
for i in range(N):  
    toss = np.random.choice([0,1])  
    toss_list.append(toss)
```

```
np.array(toss_list).sum()/N
```

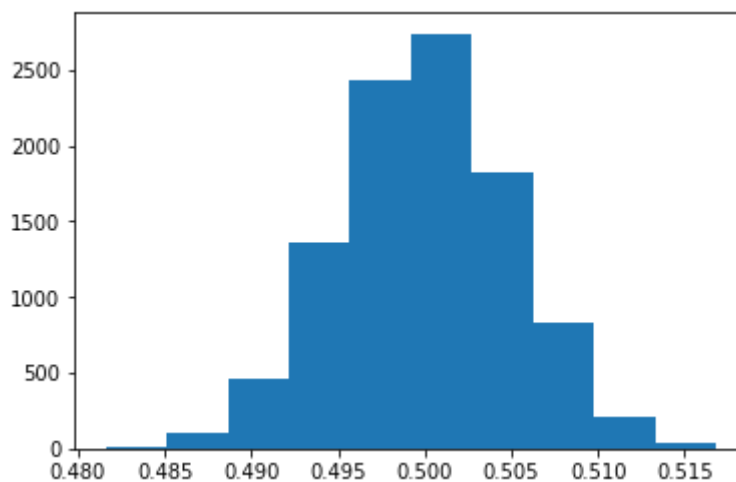
```
Out[11]: 0.5
```

Note that if we run the code snippet above again, it is likely we will get a different result. What if we run this many times?

```
In [37]: #clear
N = 10000 # number of tosses
M = 10000 # number of numerical experiments
nheads = []
for j in range(M):
    toss_list = []
    for i in range(N):
        toss_list.append(np.random.choice([0,1]))
    nheads.append( np.array(toss_list).sum()/N )
nheads = np.array(nheads)

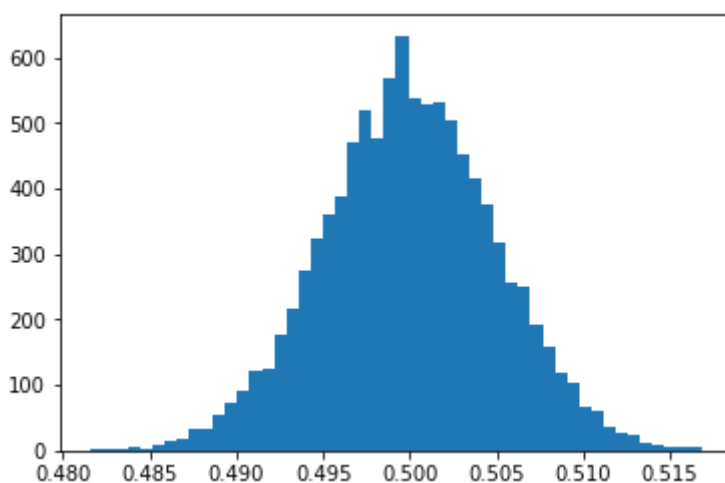
plt.hist(nheads)
```

```
Out[37]: (array([ 14., 109., 464., 1355., 2425., 2738., 1821., 826., 213.,
35.]),
array([0.4816 , 0.48512, 0.48864, 0.49216, 0.49568, 0.4992 , 0.50272,
0.50624, 0.50976, 0.51328, 0.5168 ]),
<a list of 10 Patch objects>)
```



```
In [38]: plt.hist(nheads, bins=50)
```

```
Out[38]: (array([ 2.,  1.,  3.,  5.,  3.,  8., 14., 19., 34., 34.,  5
 5.,
          72., 92., 121., 124., 176., 218., 275., 324., 362., 388., 47
 2.,
          519., 478., 568., 634., 539., 530., 531., 504., 451., 417., 37
 7.,
          319., 257., 251., 191., 160., 119., 105.,  66.,  60.,  35.,  2
 7.,
          25., 11., 10.,  5.,  4.,  5.]),
 array([0.4816 , 0.482304, 0.483008, 0.483712, 0.484416, 0.48512 ,
 0.485824, 0.486528, 0.487232, 0.487936, 0.48864 , 0.489344,
 0.490048, 0.490752, 0.491456, 0.49216 , 0.492864, 0.493568,
 0.494272, 0.494976, 0.49568 , 0.496384, 0.497088, 0.497792,
 0.498496, 0.4992 , 0.499904, 0.500608, 0.501312, 0.502016,
 0.50272 , 0.503424, 0.504128, 0.504832, 0.505536, 0.50624 ,
 0.506944, 0.507648, 0.508352, 0.509056, 0.50976 , 0.510464,
 0.511168, 0.511872, 0.512576, 0.51328 , 0.513984, 0.514688,
 0.515392, 0.516096, 0.5168  ]),
 <a list of 50 Patch objects>)
```



```
In [34]: print(nheads.mean(),nheads.std())
```

```
0.4996731 0.005044558096602714
```

What happens when we increase the number of numerical experiments?

## Approximating integrals

To approximate an integration

$$A = \int_{x_1}^{x_2} \int_{y_1}^{y_2} f(x, y) dx dy$$

we sample points uniformly inside a domain  $D = [x_1, x_2] \times [y_1, y_2]$ , i.e. we let  $X$  be a uniformly distributed random variable on  $D$ . Using Monte Carlo with  $N$  sample points, our estimate for the expected value (that a sample point is inside the circle) is:

$$S_N = \frac{1}{N} \sum_{i=1}^N f(X_i)$$

which gives the approximate for the integral:

$$A_N = (x_2 - x_1)(y_2 - y_1) \frac{1}{N} \sum_{i=1}^N f(X_i)$$

Law of large numbers:

as  $N \rightarrow \infty$ , the sample average  $S_N$  converges to the expected value  $E(X)$  and hence  $A_N \rightarrow A$

We will use Monte Carlo Method to approximate the area of a circle of radius  $R = 1$ .

Let's start with a uniform distribution on the unit square  $[0, 1] \times [0, 1]$ . Create a 2D array samples of shape  $(2, N)$ :

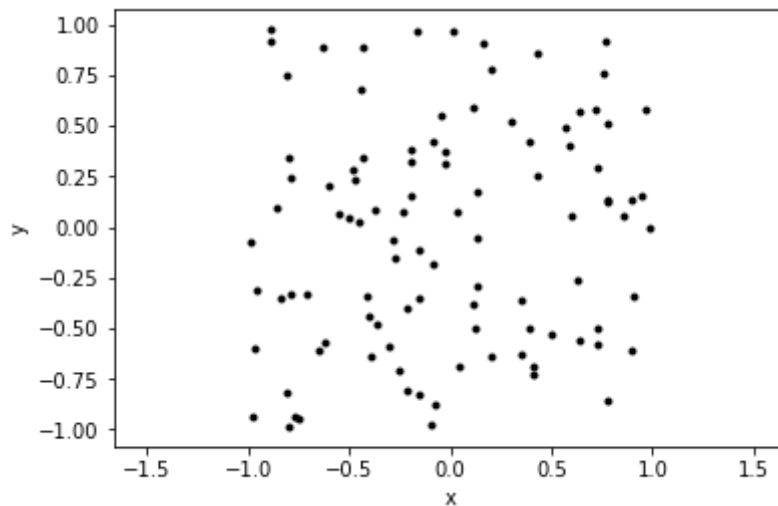
```
In [15]: N = 10**2
         samples = np.random.rand(2, N)
```

Scale the sample points "samples", so that we have a uniform distribution inside a square  $[-1, 1] \times [-1, 1]$ . Calculate the distance from each sample point to the origin  $(0, 0)$

```
In [16]: xy = samples * 2 - 1.0 # scale sample points
         r = np.sqrt(xy[0, :]**2 + xy[1, :]**2) # calculate radius
```



```
In [17]: plt.plot(xy[0,:], xy[1,:], 'k.')
plt.axis('equal')
plt.xlabel('x')
plt.ylabel('y');
```



We then count how many of these points are inside the circle centered at the origin.

```
In [18]: incircle = (r <= 1)
count_incircle = incircle.sum()
print(count_incircle)
```

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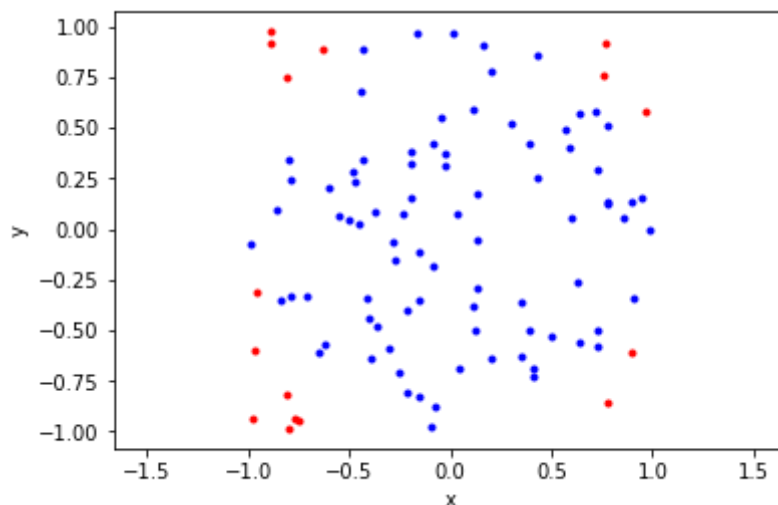
And the approximated value for the area is:

```
In [19]: A_approx = (2*2) * (count_incircle)/N
A_approx
```

Out[19]: 3.36

We can assign different colors to the points inside the circle and plot (just for visualization purposes).

```
In [20]: plt.plot(xy[0,np.where(incircle)[0]], xy[1,np.where(incircle)[0]], 'b.')
plt.plot(xy[0,np.where(incircle==False)[0]], xy[1,np.where(incircle==False)[0]], 'r.')
plt.axis('equal')
plt.xlabel('x')
plt.ylabel('y');
```



Combine all the relevant code above, so we can easily run this numerical experiment for different sample size  $N$ .

```
In [21]: #clear
N = 10**2
samples = np.random.rand(2, N)
xy = samples * 2 - 1.0 # scale sample points
r = np.sqrt(xy[0, :]**2 + xy[1, :]**2) # calculate radius
incircle = (r <= 1)
count_incircle = incircle.sum()
A_approx = (2*2) * (count_incircle)/N
print(A_approx)
```

3.0

Perform the same above, but now store the approximated area for different  $N$ , and plot:

```
In [22]: #clear
N = 10**6
samples = np.random.rand(2, N)
xy = samples * 2 - 1.0 # scale sample points
r = np.sqrt(xy[0, :]**2 + xy[1, :]**2) # calculate radius
incircle = (r <= 1)
N_samples = np.arange(1, N+1)
A_approx = 4 * incircle.cumsum() / N_samples
```

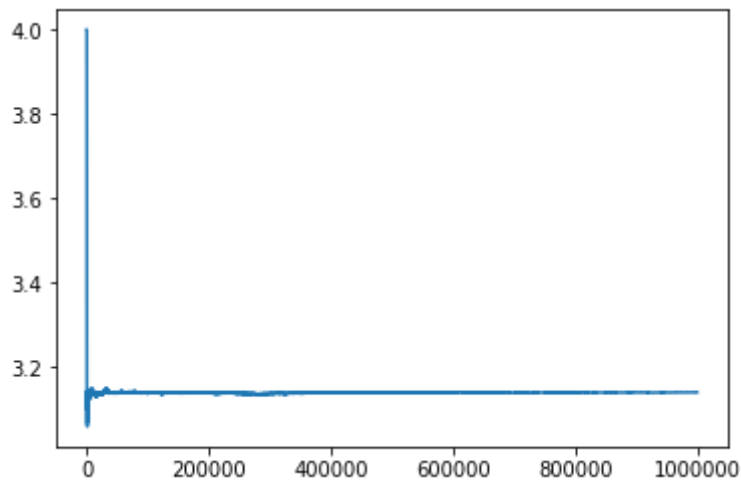
The approximated area is:

```
In [23]: #clear  
A_approx[-1]
```

```
Out[23]: 3.140908
```

```
In [24]: plt.plot(A_approx)
```

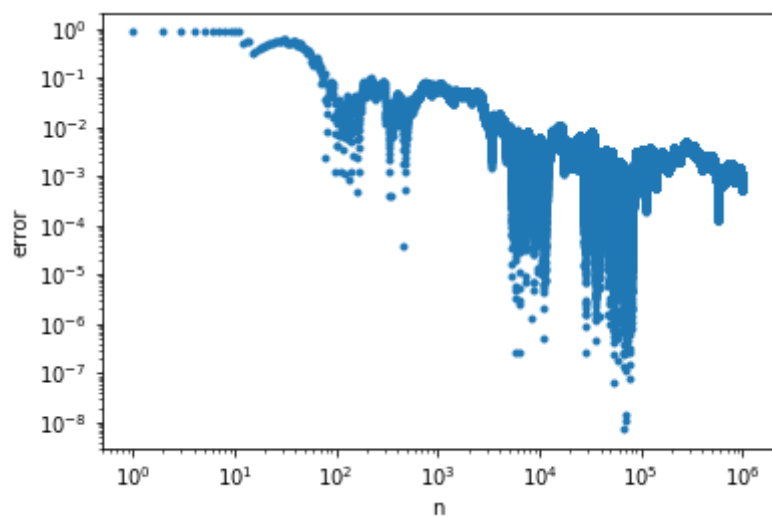
```
Out[24]: [<matplotlib.lines.Line2D at 0x11275f978>]
```



Which as expected gives an approximation for the number  $\pi$ , since the circle has radius 1. Let's plot the error of our approximation:

```
In [25]: plt.loglog(N_samples, np.abs(A_approx - np.pi), '.')  
plt.xlabel('n')  
plt.ylabel('error')
```

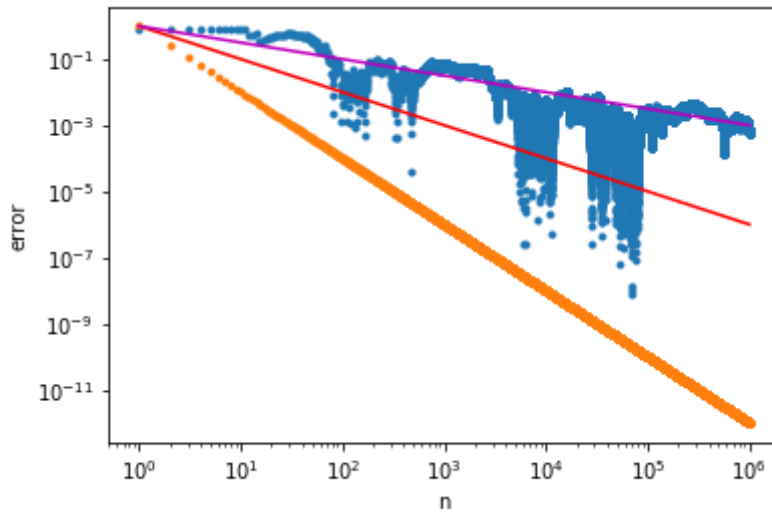
```
Out[25]: Text(0, 0.5, 'error')
```



```
In [26]: plt.loglog(N_samples, np.abs(A_approx - np.pi), '.')
plt.xlabel('n')
plt.ylabel('error')

plt.loglog(N_samples, 1/N_samples**2, '.')
plt.loglog(N_samples, 1/N_samples, 'r')
plt.loglog(N_samples, 1/np.sqrt(N_samples), 'm')
```

Out[26]: [



```
In [27]: N = 10**3
M = 1000

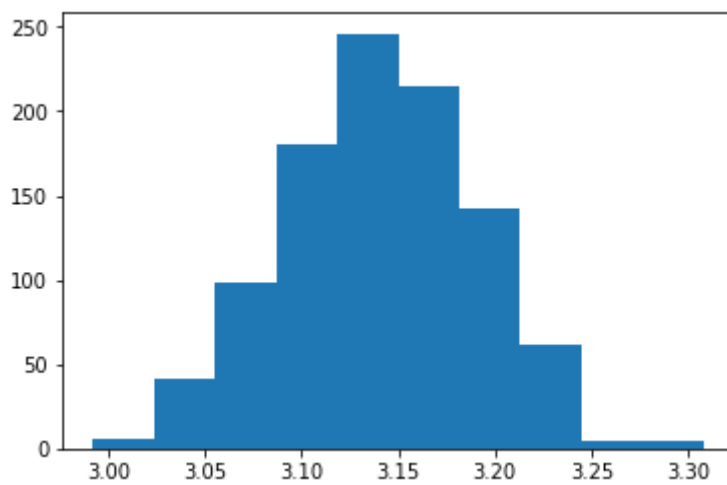
A_list = []

for i in range(M):
    samples = np.random.rand(2, N)
    xy = samples * 2 - 1.0 # scale sample points
    r = np.sqrt(xy[0, :]**2 + xy[1, :]**2) # calculate radius
    incircle = (r <= 1)
    count_incircle = incircle.sum()
    A_list.append( (2*2) * (count_incircle)/N )

A_array = np.array(A_list)

plt.hist(A_list)
```

```
Out[27]: (array([ 6., 41., 98., 180., 246., 215., 142., 62., 5., 5.]),
array([2.992 , 3.0236, 3.0552, 3.0868, 3.1184, 3.15 , 3.1816, 3.2132,
3.2448, 3.2764, 3.308 ]),
<a list of 10 Patch objects>)
```



## Approximating probabilities of a Poker game

What is the probability of winning a hand of Texas Holdem for a given starting hand?

Assumptions for this example:

- You are playing against only one player (the opponent)
- Both players stay until the end of a game (all 5 dealer cards), so players do not fold.

Monte Carlo simulation: for a given "starting hand" that we wish to obtain the winning probability, generate a set of N games and use the Texas Holdem rules to decide who wins (starting hand, opponent or there is a tie).

A game is a set of 7 random cards: 2 cards for opponent + 5 cards for the dealer.

For example, suppose the starting hand is 5 of clubs and 4 of diamonds. You perform N=1000 games, and counts 350 wins, 590 losses and 60 ties. Your numerical experiment estimated a probability of winning of 35%.

**This is your first MP!**