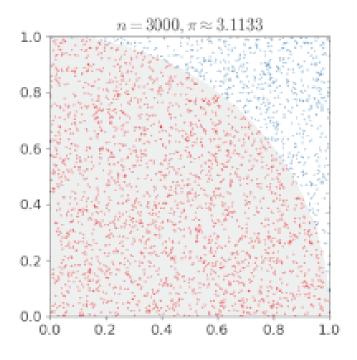
# **Lab Exercise**

- Monte Carlo Method
- Single Thread Version



# **Lab Exercise**

#### Making parallel version for Monte Carlo Method

- The parallel version get two command line arguments
  - 1) the number of threads
  - 2) the number of random points per thread
- Each thread must create random points and count the number of points that are inside the circle
- Each thread must add the number of points in the circle to the global variable "count".
- Once the threads have finished their work, the main thread uses the global variable to calculate the value of  $\pi$ .

# **Lab Exercise**

#### Running example

```
spl on ● SPL in spl/2023f/week14
□ ./w14 10000 10000
pi: 3.141585

spl on ● SPL in spl/2023f/week14
□ ./w14 10000 100000
pi: 3.141638

spl on ● SPL in spl/2023f/week14
□ ./w14 10000 1000000
pi: 3.141585

spl on ● SPL in spl/2023f/week14
□ ./w14 30000 10000000
pi: 3.141594
```

#### Exercise Hint

- Radius of the circle is always 1
- Generate random numbers between 0 and 1 using random\_r()

```
» struct random_data* random_state; // in data

» random_r(data->random_state, &x_int);

» x = (double)x_int / RAND_MAX;
21
```

#### **Exercise Submission**

- Submit your source code and Makefile
  - The make command should generate a w14 executable.
  - via iCampus
  - Bundle source code and Makefile with tar command
    - » tar.gz format
    - **\$ tar cvzf** [student\_id].tar.gz week14
  - We'll grade your submission with make
    - » If compilation fails, your points for this exercise will be zero