

Importing Numpy, Matplotlib and Scipy, also math

```
In [ ]: import numpy as np
import matplotlib.pyplot as plt
import scipy.linalg
import math
```

### Exercise 1

Creation of A in  $\mathbb{R}^n \times \mathbb{R}^n$  random matrix and  $x_{\text{true}}$  column vector of n ones

```
In [ ]: n=2
A = np.array([[1, 2], [3, 4]])
x_true = np.ones(n).reshape(-1,1)
```

*Computes the right-hand side of the linear system  $b = Ax_{\text{true}}$*

```
In [ ]: b = np.dot(A, x_true)
```

*Computes the condition number in 2-norm of the matrix A.*

```
In [ ]: condition_number_2 = np.linalg.cond(A, 2)
print(condition_number_2)
```

14.933034373659268

*It is ill-conditioned? **No, condition number is low***

*What if we use the  $\infty$ -norm instead of the 2-norm?*

```
In [ ]: condition_number_inf = np.linalg.cond(A, np.inf)
print(condition_number_inf)
```

20.999999999999993

**Still low condition number, so it is not conditioned**

*Solves the linear system  $Ax = b$  with the function `np.linalg.solve()`.*

```
In [ ]: x = np.linalg.solve(A, b)
```

*Computes the relative error between the solution computed before and the true solution  $x_{\text{true}}$*

```
In [ ]: rel_err = (np.linalg.norm(np.subtract(x, x_true), 2))/(np.linalg.norm(x_true, 2))
print(rel_err)
```

2.830524433501838e-16

*Plot a graph (using matplotlib.pyplot) with the relative errors as a function of n and (in a new window) the condition number in 2-norm  $K_2(A)$  and in  $\infty$ -norm, as a function of n*

```
In [ ]: def err_and_cond(n, vander = False, hilbert = False):
        # A = []
        # for i in range(0, n):
        #     A.append(np.arange((i*n)+1, n+(i*n)+1))
        #A= np.matrix(A)
        A = np.random.rand(n, n)
        if vander:
            A= np.vander(np.arange(1, n+1))
        if hilbert:
            A=scipy.linalg.hilbert(n)
        x_true = np.ones(n).reshape(-1,1)
        b = np.dot(A, x_true)
        condition_number_2 = np.linalg.cond(A, 2)
        condition_number_inf = np.linalg.cond(A, np.inf)
        x = np.linalg.solve(A, b)
        rel_err = (np.linalg.norm(np.subtract(x, x_true), 2))/(np.linalg.norm(x)
        return rel_err, condition_number_2, condition_number_inf
```

Getting values

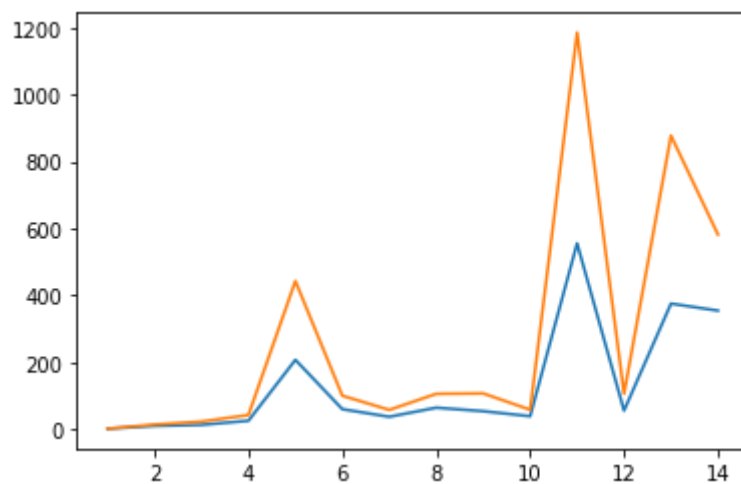
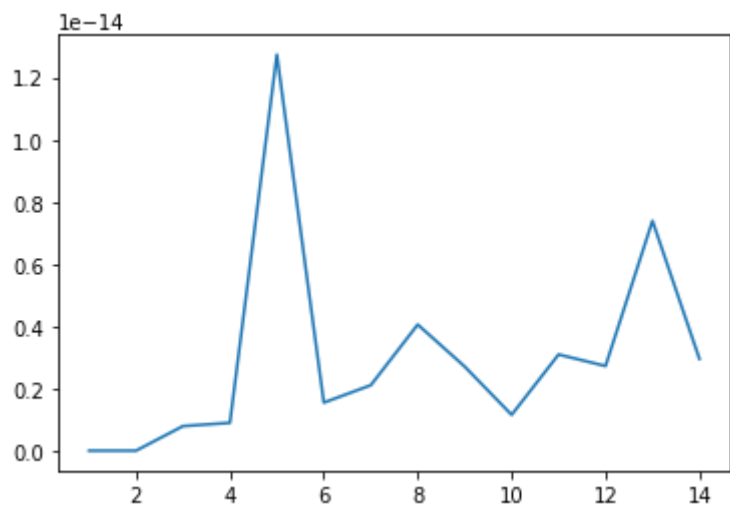
```
In [ ]: def get_values(start, end, step, vander= False, hilbert= False):
        indices = np.arange(start, end, step)
        rel_errs, conds_2, conds_inf = [], [], []
        for n in indices:
            rel_err, cond_2, cond_inf = err_and_cond(n, vander, hilbert)
            rel_errs.append(rel_err)
            conds_2.append(cond_2)
            conds_inf.append(cond_inf)
        return indices, rel_errs, conds_2, conds_inf
```

```
In [ ]: indices, rel_errs, conds_2, conds_inf = get_values(1, 15, 1)
```

Plotting

```
In [ ]: def my_plot(indices, rel_errs, conds_2, conds_inf):
        fig1 = plt.figure()
        plt.plot(indices, rel_errs)
        plt.show()
        fig2 = plt.figure()
        plt.plot(indices, conds_2)
        plt.plot(indices, conds_inf)
        plt.show()
```

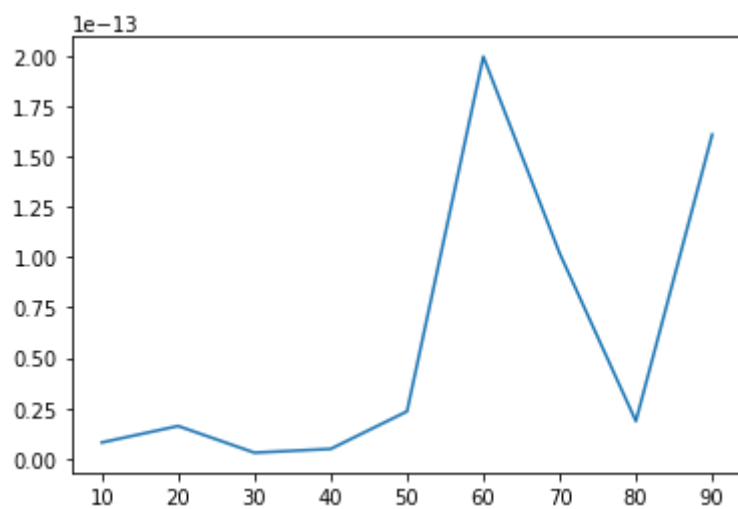
```
In [ ]: my_plot(indices, rel_errs, conds_2, conds_inf)
```

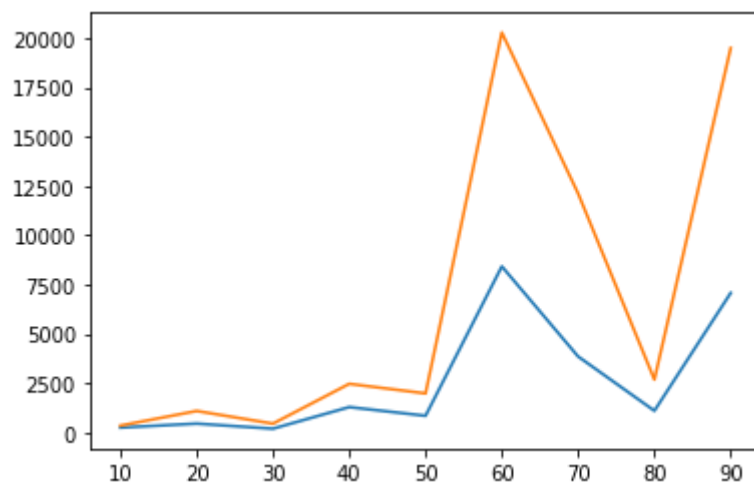


## Exercise 2

Random Matrix

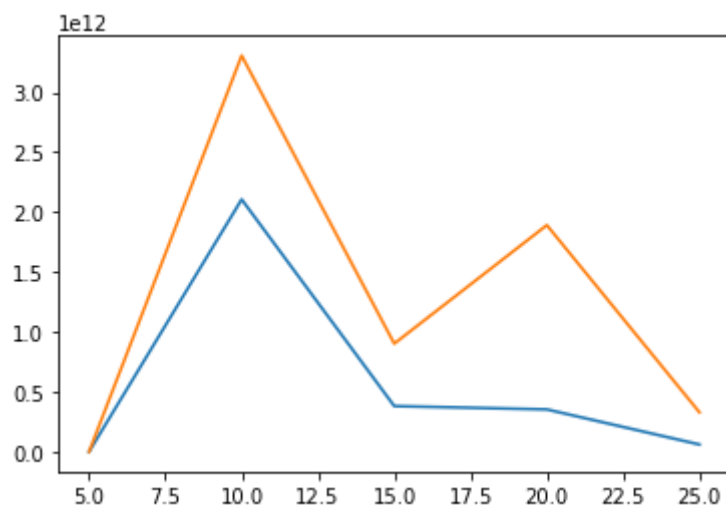
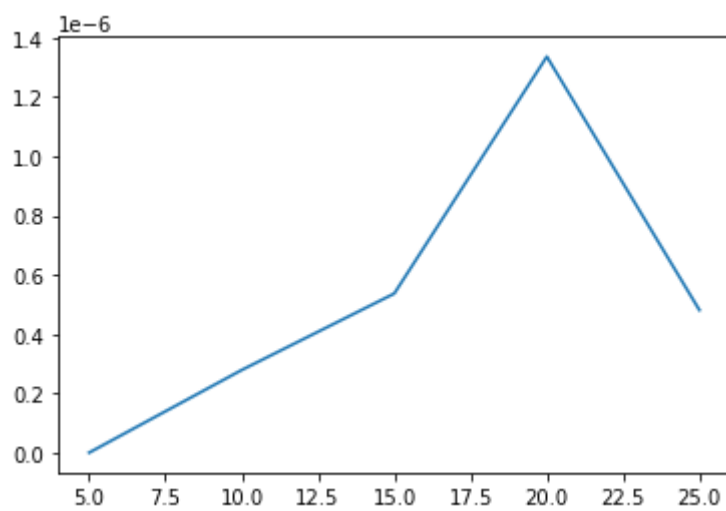
```
In [ ]: i, err, cond_2, cond_inf = get_values(10, 100, 10)
        my_plot(i, err, cond_2, cond_inf)
```





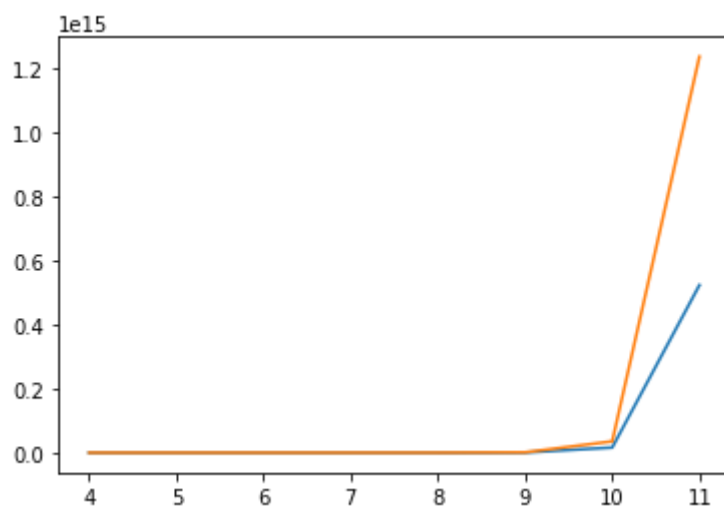
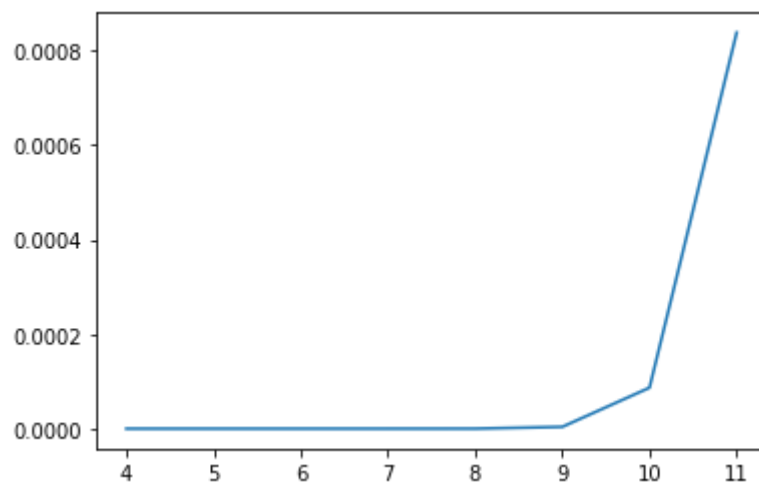
Vandermonde Matrix

```
In [ ]: i, err, cond_2, cond_inf = get_values(5, 30, 5, vander=True)
my_plot(i, err, cond_2, cond_inf)
```



Hilbert Matrix

```
In [ ]: i, err, cond_2, cond_inf = get_values(4, 12, 1, hilbert=True)
my_plot(i, err, cond_2, cond_inf)
```



### Exercise 3

Compute `machine_epsilon`, which is defined as the smallest floating point number such that it holds:  $\text{fl}(1 + \text{machine\_epsilon}) > 1$

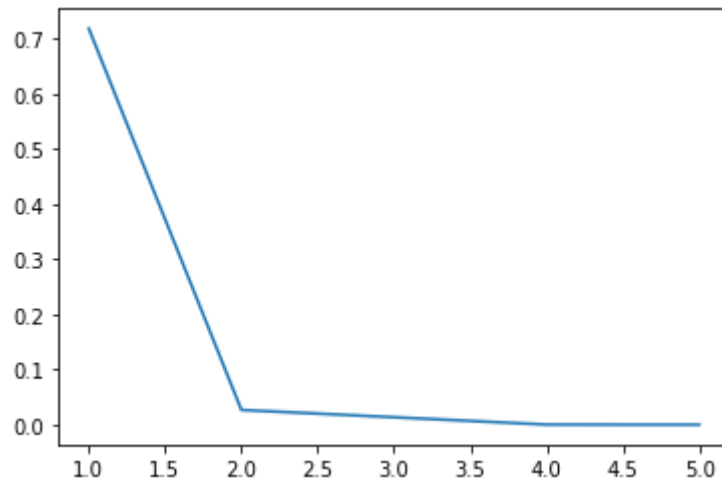
```
In [ ]: machine_epsilon = 1
while float(1+machine_epsilon) > 1:
    machine_epsilon = machine_epsilon/2
print(machine_epsilon)
```

1.1102230246251565e-16

Let's consider the sequence  $a_n = (1 + \frac{1}{n})^n$ . It is well known that:  $\lim_{n \rightarrow \infty} a_n = e$  where  $e$  is the Euler constant. Choose different values for  $n$ , compute  $a_n$  and compare it to the real value of the Euler constant. What happens if you choose a large value of  $n$ ? Guess the reason

```
In [ ]: def a_n(n):
    return (1+(1/n))**n
```

```
In [ ]: ns = [1, 50, 100, 10000, 999999999]
ecs = []
for n in ns:
    ecs.append(a_n(n))
abs_errs = []
for ec in ecs:
    abs_errs.append(math.e - ec)
plt.figure()
plt.plot(np.arange(1, len(ns)+1), abs_errs)
plt.show()
```



Let's consider the matrices:  $A = \begin{bmatrix} 4 & 2 \\ 1 & 3 \end{bmatrix}$ ,  $B = \begin{bmatrix} 4 & 2 \\ 2 & 1 \end{bmatrix}$  Compute the rank of  $A$  and  $B$  and their eigenvalues. Are  $A$  and  $B$  full-rank matrices? Can you infer some relationship between the values of the eigenvalues and the full-rank condition? Please, corroborate your deduction with other examples.

```
In [ ]: A = np.matrix([[4, 2], [1, 3]])
B = np.matrix([[4, 2], [2, 1]])
A_rank = np.linalg.matrix_rank(A)
B_rank = np.linalg.matrix_rank(B)
A_eig = np.linalg.eig(A)
B_eig = np.linalg.eig(B)
print(A_rank, B_rank)
print(A_eig[0])
print(B_eig[0])
```

```
2 1
[5. 2.]
[5. 0.]
```

$A$  is a full rank matrix, while  $B$  is not (is rank 1). If there is eigval 0, the  $\det = 0$ , so the matrix is not full ranked.

Example

```
In [ ]: #Full rank and not a eigval = 0
C = np.matrix([[1, 2, 3], [6, 12, 9], [9, 2, 1]])
print(np.linalg.matrix_rank(C))
print(np.linalg.eig(C)[0])

#Not full rank and not a eigval = 0
D = np.matrix([[1, 2, 3], [4, 5, 6], [7, 8, 9]])
print(np.linalg.matrix_rank(D))
print(np.linalg.eig(D)[0])

#Not full rank because a eigval = 0
E = np.matrix([[1, 0, 3], [2, 0, 4], [5, 0, 9]])
print(np.linalg.matrix_rank(E))
print(np.linalg.eig(E)[0])

3
[15.46698572 -3.8716714  2.40468568]
2
[ 1.61168440e+01 -1.11684397e+00 -1.30367773e-15]
2
[ 0.          -0.56776436 10.56776436]
```