

# Linear Algebra and its Applications

## Assignment 5

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## Import all prerequisites

```
In [1]: from helper_functions import *
import time
import scipy as sp
import pandas as pd
import scipy.linalg
from scipy.linalg import cho_factor, cho_solve, cholesky
import matplotlib.pyplot as plt
```

## Initialization

Initialization of all data structures for storing the following:

- Running time of both solvers
- Solutions  $L, U, P, x$  or  $P, G, x$  and the chosen  $A$  and  $b$  for both solvers

```
In [2]: n_list=[2,5,9,10,12,13,14,15,16,30,50,90,100]
scipy_LU_rt={} #Running times by scipy's LU factorization algorithm
my_LU_rt_gepp={} # Running times by my implementation LU factorization algorithm
my_LU_rt_gerp={}
my_LU_rt_gecp={}
my_cholesky_rt={}
scipy_cholesky_rt={}

scipy_cholesky_rt={}
scipy_LU_subst_rt={} #Running times by scipy's Substitution algorithm
# Running times by my implementation of Substitution algorithm
my_subst_rt_gepp={}
my_subst_rt_gerp={}
my_subst_rt_gecp={}
my_cholesky_subst_rt={}

scipy_LU_solutions={}
scipy_cholesky_solutions={}
my_solutions_gepp={}
my_solutions_gerp={}
my_solutions_gecp={}
my_cholesky_solutions={}

```

# Execution

Execute the solvers and record the running times

```
In [5]: import warnings
warnings.filterwarnings("ignore")
for n in n_list:

    A=np.zeros((n,n))
    for i in range(n):
        for j in range(n):
            A[i,j]=1/(i+j+1)
    b=np.random.rand(n)

    #Partial Pivoting
    start_time = time.time()
    L,U,P,Q=LU_factorize_partial(A)
    my_LU_rt_gepp[n]=time.time()-start_time

    start_time=time.time()
    x=substitution(L,U,P,Q,b)
    my_subst_rt_gepp[n]=time.time()-start_time
    my_solutions_gepp[n]=(L,U,P,Q,x,A,b)

    #Rook Pivoting
    start_time = time.time()
    L,U,P,Q=LU_factorize_rook(A)
    my_LU_rt_gerp[n]=time.time()-start_time

    start_time=time.time()
    x=substitution(L,U,P,Q,b)
    my_subst_rt_gerp[n]=time.time()-start_time
    my_solutions_gerp[n]=(L,U,P,Q,x,A,b)

    #Complete Pivoting
    start_time = time.time()
    L,U,P,Q=LU_factorize_complete(A)
    my_LU_rt_gecp[n]=time.time()-start_time

    start_time=time.time()
    x=substitution(L,U,P,Q,b)
    my_subst_rt_gecp[n]=time.time()-start_time
    my_solutions_gecp[n]=(L,U,P,Q,x,A,b)

    #My Cholesky LU Solver
    start_time = time.time()
    G,P=my_cholesky(A)
    my_cholesky_rt[n]=time.time()-start_time
```

```

start_time=time.time()
x=substitution(G,G.transpose(),P,P.transpose(),b)
my_cholesky_subst_rt[n]=time.time()-start_time
my_cholesky_solutions[n]=(G,P,x,A,b)

#Scipy LU Solver
start_time=time.time()
p, l, u = sp.linalg.lu(A) # The decomposition is A=plu
scipy_LU_rt[n]=time.time()-start_time

lu, piv = sp.linalg.lu_factor(A)
start_time=time.time()
x2 = sp.linalg.lu_solve((lu, piv), b)
scipy_LU_subst_rt[n]=time.time()-start_time

scipy_LU_solutions[n]=(l,u,p,x2,A,b)

#Scipy Cholesky Solver
start_time=time.time()
try:
    L = cholesky(A, lower=True)
    scipy_cholesky_rt[n]=time.time()-start_time
except:
    L=np.NaN
try:
    c, low = cho_factor(A)
    x = cho_solve((c, low), b)
except:
    x=np.NaN

scipy_cholesky_solutions[n]=(L,x,A,b)

```

## Calculate Norms

Calculate the norms of  $PA - LU$  and  $Ax_0 - b$  using all solvers.

```

In [6]: my_PA_LU_norms_gepp={}
my_PA_LU_norms_gerp={}
my_PA_LU_norms_gecp={}
my_cholesky_GG_norms={}
scipy_PA_LU_norms={}
scipy_cholesky_GG_norms={}

my_Ax_b_norms_gepp={}
my_Ax_b_norms_gerp={}
my_Ax_b_norms_gecp={}
my_cholesky_AX_b_norms={}
scipy_Ax_b_norms={}

```

```

scipy_cholesky_AX_b_norms={}

for n in n_list:
    #Partial Pivoting Norms
    L,U,P,Q,x,A,b=my_solutions_gepp[n]
    my_PA_LU_norms_gepp[n]=np.linalg.norm(np.matmul(P,A@Q)-np.matmul(L,U),ord=2)
    my_Ax_b_norms_gepp[n]=np.linalg.norm(np.matmul(A,x)-b)

    #Rook Pivoting Norms
    L,U,P,Q,x,A,b=my_solutions_gerp[n]
    my_PA_LU_norms_gerp[n]=np.linalg.norm(np.matmul(P,A@Q)-np.matmul(L,U),ord=2)
    my_Ax_b_norms_gerp[n]=np.linalg.norm(np.matmul(A,x)-b,ord=2)

    #Complete Pivoting Norms
    L,U,P,Q,x,A,b=my_solutions_gecp[n]
    my_PA_LU_norms_gecp[n]=np.linalg.norm(np.matmul(P,A@Q)-np.matmul(L,U),ord=2)
    my_Ax_b_norms_gecp[n]=np.linalg.norm(np.matmul(A,x)-b,ord=2)

    #Cholesky Norms
    G,P,x,A,b=my_cholesky_solutions[n]
    try:
        my_cholesky_GG_norms[n]=np.linalg.norm(G@(G.transpose())-P@A@(P.transpose())
        my_cholesky_AX_b_norms[n]=np.linalg.norm(np.matmul(A,x)-b,ord=2)
    except:
        my_cholesky_GG_norms[n]=np.NaN
        my_cholesky_AX_b_norms[n]=np.NaN

    #Scipy LU Norms
    L,U,P,x,A,b=scipy_LU_solutions[n]
    scipy_PA_LU_norms[n]=np.linalg.norm(A-np.matmul(P,np.matmul(L,U)),ord=2)
    scipy_Ax_b_norms[n]=np.linalg.norm(np.matmul(A,x)-b,ord=2)

    #Scipy Cholesky Norm
    G,x,A,b=scipy_cholesky_solutions[n]
    try:
        scipy_cholesky_GG_norms[n]=np.linalg.norm(G@(G.transpose())-A,ord=2)
        scipy_cholesky_AX_b_norms[n]=np.linalg.norm(np.matmul(A,x)-b,ord=2)
    except:
        scipy_cholesky_GG_norms[n]=np.NaN
        scipy_cholesky_AX_b_norms[n]=np.NaN

```

## Norms Table

This table gives us the the matrix norms of  $PA - LU$ (or  $PAP^T - GG^T$ ) and  $Ax_0 - b$  for both solvers.

The columns labels are set as n for an  $n \times n$  matrix

```

In [7]: matrix_norms={}
col_names=["PA-LU gepp","PA-LU gerp","PA-LU gecp","Cholesky GG(t)-PAP(t)","PA-LU Sc
          "Ax-b gepp","Ax-b gerp","Ax-b gecp","Ax-b Cholesky","Ax-b SciPy(LU)","Ax
for n in n_list:
    matrix_norms[n]=[my_PA_LU_norms_gepp[n],my_PA_LU_norms_gerp[n],my_PA_LU_norms_g
                    scipy_PA_LU_norms[n],

```

```

        scipy_cholesky_GG_norms[n],my_Ax_b_norms_gepp[n],my_Ax_b_norms
        my_Ax_b_norms_gecp[n],my_cholesky_AX_b_norms[n],scipy_Ax_b_norms
        scipy_cholesky_AX_b_norms[n]]
norm_table=pd.DataFrame(matrix_norms, index=col_names)

```

```
In [8]: display(norm_table.iloc[:,6])
```

|                              |          | 2            | 5            | 9            | 10           | 12           | 13  |
|------------------------------|----------|--------------|--------------|--------------|--------------|--------------|-----|
| <b>PA-LU gepp</b>            | 0.000000 | 2.775558e-17 | 5.173899e-17 | 4.913042e-17 | 6.261478e-17 | 5.871185e-17 |     |
| <b>PA-LU gerp</b>            | 0.000000 | 2.775558e-17 | 5.173899e-17 | 4.913042e-17 | 6.261478e-17 | 5.871185e-17 |     |
| <b>PA-LU gecp</b>            | 0.000000 | 4.490946e-17 | 5.973831e-17 | 5.093450e-17 | 6.702340e-17 | 4.303850e-17 |     |
| <b>Cholesky GG(t)-PAP(t)</b> | 0.000000 | 1.340158e-16 | 6.700789e-17 | 1.190237e-16 | 2.262175e-16 |              | NaN |
| <b>PA-LU SciPy</b>           | 0.000000 | 2.775558e-17 | 4.033996e-17 | 1.962616e-17 | 5.160349e-17 | 5.722031e-17 |     |
| <b>Cholesky Scipy</b>        | 0.000000 | 0.000000e+00 | 2.504924e-17 | 2.715048e-17 | 3.359559e-17 | 3.301474e-17 |     |
| <b>Ax-b gepp</b>             | 0.000000 | 1.901657e-12 | 3.618864e-06 | 2.965031e-05 | 8.466475e-03 | 9.488734e-01 |     |
| <b>Ax-b gerp</b>             | 0.000000 | 1.901657e-12 | 3.618864e-06 | 2.965031e-05 | 8.466475e-03 | 9.488734e-01 |     |
| <b>Ax-b gecp</b>             | 0.000000 | 4.623118e-12 | 5.181760e-06 | 2.286102e-05 | 5.828233e-03 | 1.272643e+00 |     |
| <b>Ax-b Cholesky</b>         | 0.105849 | 8.178088e-02 | 2.998494e-01 | 6.461434e-01 | 7.405266e-01 |              | NaN |
| <b>Ax-b SciPy(LU)</b>        | 0.000000 | 3.631007e-12 | 6.229130e-06 | 2.720404e-05 | 6.879998e-03 | 7.823553e-01 |     |
| <b>Ax-b Scipy(Cholesky)</b>  | 0.000000 | 4.543342e-12 | 5.865915e-06 | 2.327277e-05 | 4.451865e-03 | 1.515401e+00 |     |

```
In [9]: display(norm_table.iloc[:,6:])
```

|                              | 14           | 15           | 16           | 30           | 50           |           |
|------------------------------|--------------|--------------|--------------|--------------|--------------|-----------|
| <b>PA-LU gepp</b>            | 5.330853e-17 | 6.062295e-17 | 5.949965e-17 | 6.744857e-17 | 7.113867e-17 | 7.323909  |
| <b>PA-LU gerp</b>            | 5.330853e-17 | 6.062295e-17 | 5.949965e-17 | 6.720375e-17 | 7.103216e-17 | 7.328106  |
| <b>PA-LU gecp</b>            | 4.908086e-17 | 4.865046e-17 | 5.244584e-17 | 7.322928e-17 | 7.438936e-17 | 8.073525  |
| <b>Cholesky GG(t)-PAP(t)</b> | NaN          | NaN          | NaN          | NaN          | NaN          |           |
| <b>PA-LU SciPy</b>           | 5.481817e-17 | 5.934849e-17 | 6.081652e-17 | 6.448746e-17 | 6.896413e-17 | 7.237879  |
| <b>Cholesky SciPy</b>        | 3.184169e-17 | NaN          | NaN          | NaN          | NaN          |           |
| <b>Ax-b gepp</b>             | 1.418014e+01 | 1.395668e+00 | 6.353899e+00 | 3.087226e+01 | 3.828336e+01 | 3.342325e |
| <b>Ax-b gerp</b>             | 1.418014e+01 | 1.395668e+00 | 6.353899e+00 | 3.991581e+01 | 3.926930e+01 | 4.062674e |
| <b>Ax-b gecp</b>             | 5.594126e+01 | 2.096839e+00 | 4.985956e+00 | 1.357073e+02 | 6.260920e+01 | 4.531907e |
| <b>Ax-b Cholesky</b>         | NaN          | NaN          | NaN          | NaN          | NaN          |           |
| <b>Ax-b SciPy(LU)</b>        | 4.907927e+00 | 1.332972e+00 | 2.507493e+00 | 3.268468e+02 | 1.179363e+02 | 6.532960e |
| <b>Ax-b SciPy(Cholesky)</b>  | 2.841543e+00 | NaN          | NaN          | NaN          | NaN          |           |

A is a Hilbert matrix and it ceases to be positive definite after  $n = 14$ . Thus, after  $n = 14$ , Cholesky decompositon doesn't work on A and we get *NaN* values.

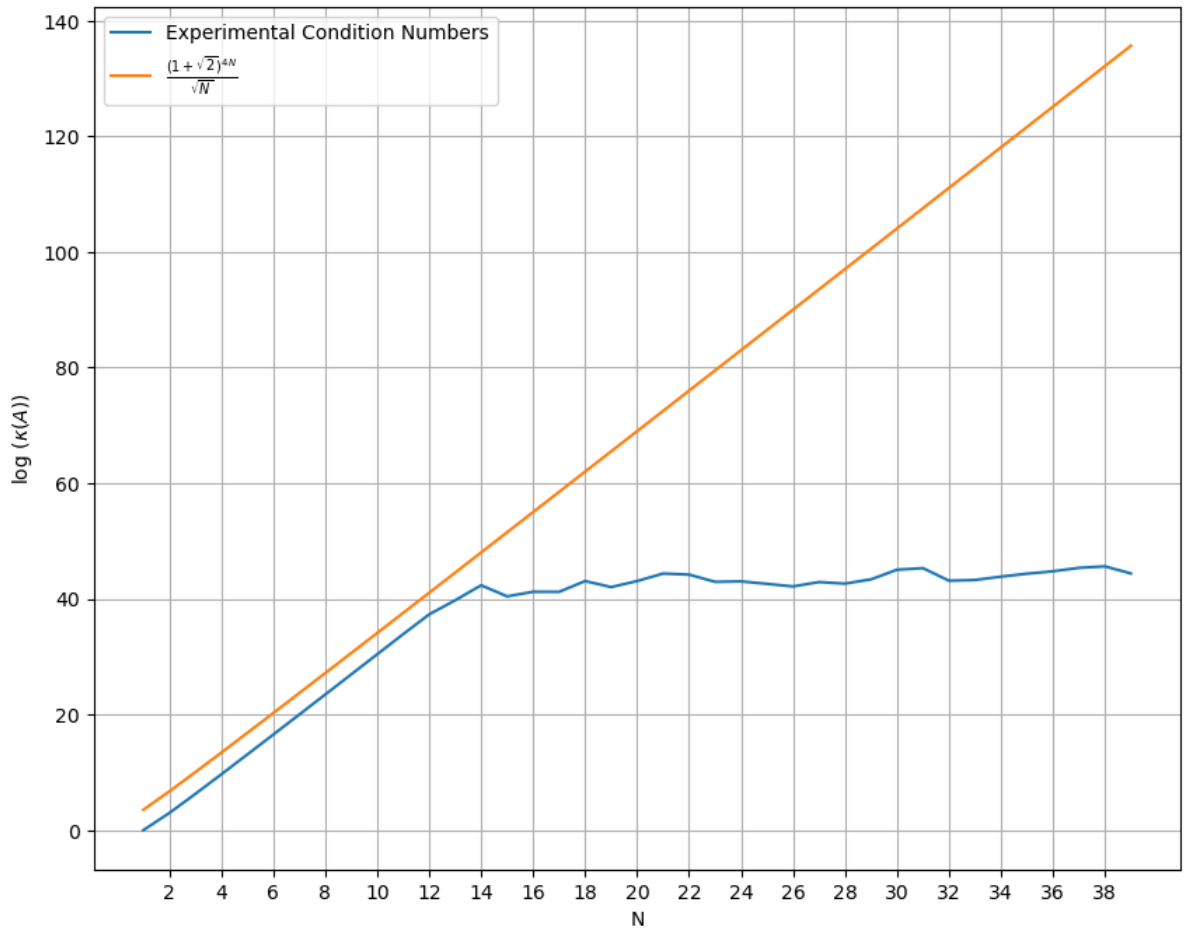
## Condition Number

```
In [10]: conds=[]
bounds=[]
n_list=np.arange(1,40)
for n in n_list:
    A=np.zeros((n,n))
    for i in range(n):
        for j in range(n):
            A[i,j]=1/(i+j+1)
    cond=np.linalg.norm(A)*np.linalg.norm(np.linalg.inv(A))
    conds.append(cond)
    bounds.append(((1+np.sqrt(2))**(4*n))/np.sqrt(n))

import matplotlib.pyplot as plt

plt.figure(figsize=(10, 8))

plt.plot(n_list,np.log(conds),label = 'Experimental Condition Numbers')
plt.plot(n_list,np.log(bounds),label = r'$\frac{(1+\sqrt{2})^{4N}}{\sqrt{N}}$')
plt.xlabel('N')
plt.ylabel('log ( $\kappa(A)$ )')
plt.xticks(2*n_list[:int(n/2)])
plt.legend()
plt.grid()
```



In the above graph, we plot the experimental condition number along with the value  $\frac{(1+\sqrt{2})^{4N}}{\sqrt{N}}$ .

We can say that for upto  $n = 14$ , the condition number has the bound  $O\left(\frac{(1+\sqrt{2})^{4N}}{\sqrt{N}}\right)$  but beyond  $n = 14$  calculating the condition numbers involves a lot of round off errors and thus stagnates.

Still we can say, that  $O\left(\frac{(1+\sqrt{2})^{4N}}{\sqrt{N}}\right)$  is a reasonable bound on the condition number as the experimental values are below the curve.

## References

1. Beckermann, B. (2000). The condition number of real Vandermonde, Krylov and positive definite Hankel matrices. *Numerische Mathematik*, 85(4), 553–577.  
doi:10.1007/pl00005392
2. Todd, J. "The condition of the finite segments of the Hilbert matrix, *NBS Appl. Math. Ser* 39 (1954): 109-116.