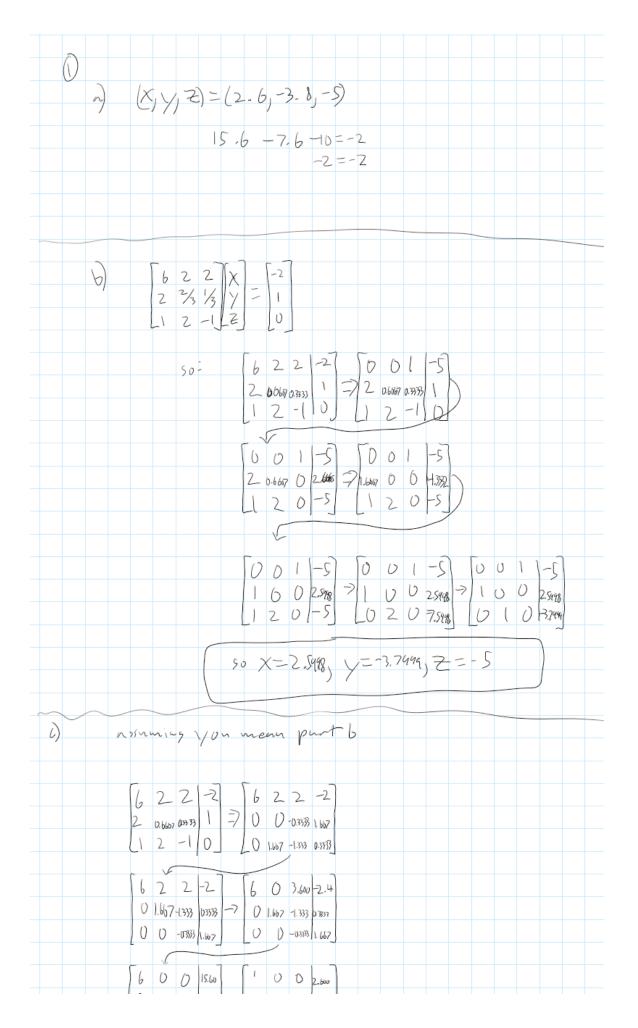
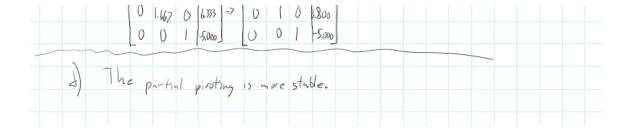
Homeowork 12

Problem 1

a)

```
In [1]:
       import numpy as np
       import scipy as scipy
       import householder as hh
In [2]: a = 1ambda x, y, z: 6*x+2*y+2*z
       b= lambda x,y,z: 2*x + (2/3)*y + (1/3)*z
       c = lambda x,y,z: x+2*y-z
In [3]: a1=a(2.6,-3.8,-5)
       b1=b(2.6,-3.8,-5)
       c1=c(2.6,-3.8,-5)
       print("Equation 1(x,y,z)=",a1)
       print("Equation 2(x,y,z)=",b1)
       print("Equation 3(x,y,z)=",c1)
      Equation 2(x,y,z) = 1.00000000000000000
      Equation 3(x,y,z)=0.0
       a-d)
```





Problem 2

```
In [4]: A = np.array(([12,10,4],[10,8,-5],[4,-5,3]))
        print(A)
       [[12 10 4]
        [10 8 -5]
        [ 4 -5 3]]
In [5]: print(np.linalg.eigvals(A))
       [-5.1984251 20.1984251 8.
                                       ]
In [6]: D,C = hh.householder(A)
        A3 = np.diag(D)+np.diag(C,1)+np.diag(C,-1)
        print(A3)
        print(np.linalg.eigvals(A3))
       [10 4]
       [[ 12 -10
                  0]
       [-10
              3
                  5]
       [ 0
              5
                  7]]
       [-4.9774314 19.11607184 7.86135957]
```

The above finding for the householder matrix supplies eigenvalues close to the original. I believe this is due to some numerical errors in the calculation.

Problem 3

```
lam = (x.T@A@x)/(x.T@x)
        if np.abs(lam - lam_store) < tol:</pre>
            break
        lam_store = lam
    return lam, x, i
def powerNoN(A):
    x = np.ones(len(A)).T
    tol = 1e-6
    max_iter = 100
    lam_store = 0
    for i in range(max_iter):
        x = A@x / np.linalg.norm(A@x)
        lam = (x.T@A@x)/(x.T@x)
        if np.abs(lam - lam_store) < tol:</pre>
            break
        lam_store = lam
    return lam, x, i
def powersmall(N):
    B=np.zeros((N,N))
    for i in range(0,N):
        for j in range(0,N):
            B[i][j]=1/(i+j+1)
    lambig,xbig,iterbig = power(N)
    A = B-lambig*np.identity(N)
    x = np.ones(N).T
    tol = 1e-6
    max_iter = 100
    lam_store = 0
    for i in range(max_iter):
        x = A@x / np.linalg.norm(A@x)
        lam = (x.T@A@x)/(x.T@x)
        if np.abs(lam - lam_store) < tol:</pre>
            break
        lam_store = lam
    return lam+lambig, x, i
```

```
In [17]: | lam1,x1,iter1 = power(4)
         print(lam1)
         print(x1)
         print(iter1)
        0.9999996889870862
                     0.70953094 0.53214854 0.46193483]
        [0.
        11
In [18]: lam2, x2, iter2 = power(20)
         print(lam2)
         print(x2)
         print(iter2)
        0.9999996302430362
                     0.43410165 0.32557626 0.28261831 0.25753438 0.24035505
         0.22752285 0.21740095 0.20911282 0.20213897 0.19614823 0.19091758
         0.18629022 \ 0.18215206 \ 0.1784177 \quad 0.17502169 \ 0.17191282 \ 0.16905038
         0.16640147 0.16393916]
        13
         b)
In [36]:
         lamsmall,xsmall,itersmall = powersmall(16)
         print(lamsmall)
         N=16
         B=np.zeros((N,N))
         for i in range(0,N):
              for j in range(0,N):
                  B[i][j]=1/(i+j+1)
         lamList = np.linalg.eigvals(B)
         print(min(lamList)-lamsmall)
        0.004951309919023927
        (-0.004951309919023933-9.227487176840956e-19j)
         This method is accurate to 3 decimal places.
         c)
```

```
In [45]: E = 0.01*np.ones((N,N))

C = B+E

lamPertMin = (np.min(np.linalg.eig(C)[0]))

print("Difference in smallest eigenvalues:",lamsmall-lamPertMin)
print("2-Norm of E:",np.linalg.norm(E))
```

Difference in smallest eigenvalues: 0.004951309919023941 2-Norm of E: 0.16

Therefore this is consistent with the Bauer-Fike Theorem.

d)

```
In [57]: F = np.array([[1,0,0],[0,(np.sqrt(2)/2)*(1-1j),0],[0,0,(np.sqrt(2)/2)*(1+1j)]])
         print(F)
        [[1.
                    +0.j
                                 0.
                                           +0.j
                                                                   +0.j
                                                                               ]
         [0.
                                 0.70710678-0.70710678j 0.
                    +0.j
                                                                  +0.j
         [0.
                    +0.j
                                                        0.70710678+0.70710678j]]
                                           +0.j
In [58]: print(powerNoN(F))
        ((0.804737854124365+5.204480025290235e-18j), array([ 0.57735027+0.00000000e+00j,
        -0.57735027-6.00627088e-17j,
               -0.57735027+6.00627088e-17j]), 99)
```

Here $|\lambda|=1$ for all eigenvalues, and is the maximum, therefore the power method will break down as $\frac{|\lambda_2|}{|\lambda_1|}$ is 1 and no eigenvalue is approached.