Computer Assignment

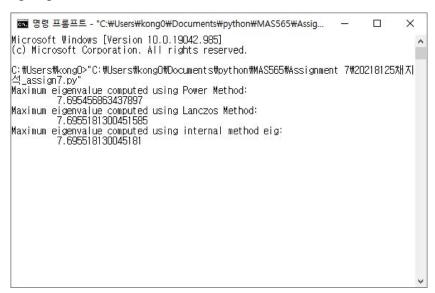
The program which does the required is submitted via KLMS along with this document. It performs two tasks, as required in the assignment sheet.

Firstly the program performs the Power method. The implementation details are straightforward, being a direct translation of the equations presented in Exercise 6.14. As in Assignment 5, the stop condition is set to be when the relative error is less than 10^{-6} .

After that the program performs the Lanczos method. In a broad sense the code is again a direct translation of equations (6.5.3.1a) and (6.5.3.1b) in the textbook. The iteration shall stop when a small value of $|\gamma_i|$ is met, and this threshold is again set to be 10^{-6} . Our goal is to compute the full tridiagonal matrix which is similar to A_7 , but the iteration may stop early when the number of iterations has reached the dimension of the Krylov space $K(q, A_7)$ where q is the initial vector. When $q = e_1$ as indicated in the assignment sheet, this dimension turns out to be 25. This fact can be verifed by using the symbolic computation module sympy and the Python code included in the appendix. Anyhow, if the Lanczos iteration terminates early, this means that the tridiagonal matrix we desire is actually a block diagonal matrix of tridiagonal matrices, and we have to find a new initial vector to initiate the iteration in order to compute the next block. Such initial vector must be orthogonal to all q_i 's computed up to that point, so we take the strategy of running through the canonical basis vectors and taking the orthogonal complement of the projections onto span(q_i) until we find a nonzero orthogonal complement.

Once we obtain all γ_i 's and δ_i 's the rest is simple: we pass the computed tridiagonal matrix into the function numpy.linalg.eig provided by the numpy package in order to compute the eigenvalues.

The computed maximum eigenvalues are presented as outputs, as we see below. As a comparison we also show the maximum eigenvalue obtained by using the numpy-provided function numpy.linalg.eig.



Power method show a relative error of 7.96×10^{-6} . However such seemingly large error is due to us setting the threshold to 10^{-6} . If we set a smaller threshold we see that the Power method actually produces a result much closer to the result of numpy.linalg.eig. The performance of the Lanczos method is much more interesting. The result coinsides with the true value up to 14 digits. Again, we cannot conclude hastily that the result of the Lanczos method agrees with that of the Power method, since we allowed the Power method to stop early with low accuracy. If we increase the accuracy of the Power method, we see an agreement.

It is also required to plot the eigenvalues of the computed tridiagonal matrix. The resulting plot is as the following figure.

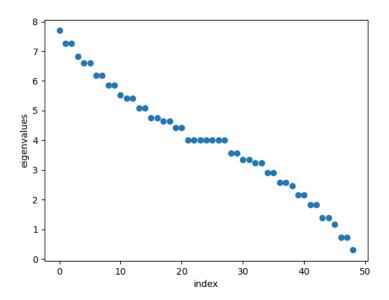


Figure 1: Scatter plot of the eigenvalues of A_7 , in decreasing order

Appendix

The promised code, which computes the dimension of the Krylov space $K(e_1, A_7)$, is as follows.

```
import numpy as np
import sympy as sp
def make_Tn(n):
   if n \le 0:
       raise KeyError
   off_main_diag = [-1 for _ in range(n-1)]
   upper = np.diag(off_main_diag, 1)
   lower = np.diag(off_main_diag, -1)
   return 4*np.eye(n) + upper + lower
def make_An(n):
   if n <= 0 :
       raise KeyError
   Tn = make_Tn(n)
   if n == 1 :
       return Tn
   I = np.eye(n)
   0 = np.zeros((n,n))
   res = np.block([Tn, -I] + [0] * (n-2))
   for i in range(1, n-1):
       tmp = np.block([0] * (i-1) + [-I, Tn, -I] + [0] * (n-2-i))
       res = np.vstack((res, tmp))
   tmp = tmp = np.block([0] * (n-2) + [-I, Tn])
   res = np.vstack((res, tmp))
   return res
```

```
n = 7
A = sp.Matrix(np.array(make_An(n), dtype = int))
e1 = sp.Matrix(np.array(np.hstack(([1], np.zeros(n**2-1))), dtype = int))

K = e1
for i in range(1, 49):
    e1 = A * e1
    K = K.col_insert(i, e1)

print(K.rank())
```