MAT 167 - Project 1

Emily Ng, Honghui Li, Steve Sliva, Xinran Jiang, Yuchen Liu

August 2023

Problem 1

Given:

Matrices can be partitioned into several submatrices known as "blocks." Using a blocked approach, matrix-matrix multiplication can be performed with a divide and conquer approach. Supposed we have two matrices,

a) Show that C can be written as

$$C = \begin{pmatrix} M_1 + M_4 + M_5 - M_7 & M_3 + M_5 \\ M_2 + M_4 & M_1 - M_2 + M_3 + M_6 \end{pmatrix}$$
$$M1 = (A_{11} + A_{22})(B_{11} + B_{22})$$
$$M2 = (A_{21} + A_{22})B_{11}$$

 $M3 = A_{11}(B_{12} - B_{22})$

 $M4 = A_{22}(B_{21} - B_{11})$

 $M5 = (A_{11} + A_{12})B_{22}$

 $M6 = (A_{21} - A_{11})(B_{11} + B_{12})$

 $M7 = (A_{12} - A_{22})(B_{21} + B_{22})$

This works for any even shape matrix, so no matter if the A_{xy} is a number or sub-matrix, it follows the rule that we have $A_{xy}(A_{ab} + A_{cd}) = A_{xy} \cdot A_{ab} + A_{xy} \cdot A_{cd}$ (Distributive Property)

1-1) for the first column and first row

$$M_1 + M_4 + M_5 - M_7 =$$

 $A_{11}B_{22} + A_{22}B_{11} + A_{22}B_{22} - A_{22}B_{21} - A_{22}B_{11} + A_{11}B_{22} + A_{12}B_{22} + A_{12}B_{21} + B_{22}A_{12} - A_{22}B_{21} - A_{22}B_{22} - A_{22}B_{22} - A_{22}B_{21} - A_{22}B_{22} - A_{22}B_{22} - A_{22}B_{21} - A_{22}B_{22} - A_{22}B_{21} - A_{22}B_{22} - A_{22}B_{22} - A_{22}B_{21} - A_{22}B_{22} - A_{22}B_{21} - A_{22}B_{22} - A_{22}B_{21} - A_{22}B_{22} - A_{22}B_{21} - A_{22}B_{22} - A_{22}B_{22} - A_{22}B_{21} - A_{22}B_{22} - A_{22}B_{21} - A_{22}B_{22} - A_{22}B_{22} - A_{22}B_{21} - A_{22}B_{22} - A_{22}B_{22} - A_{22}B_{21} - A_{22}B_{22} - A_{22}B_{22} - A_{22}B_{22} - A_{22}B_{21} - A_{22}B_{22} - A_{22}B_{22} - A_{22}B_{21} - A_{22}B_{22} - A_{2$

$$= A_{11}B_{11} + A_{12}B_{21}$$

1-2) for the first column and second row

$$M3 + M5 =$$

$$A_{11}B_{12} - A_{11}B_{22} + A_{11}B_{22} + A_{12}B_{22}$$

= $A_{11}B_{12} + A_{12}B_{22}$

2-1) for the second column and first row

$$M2 + M4 =$$

$$A_{21}B_{11} + A_{22}B_{11} + A_{22}B_{21} - A_{22}B_{11}$$

$$= A_{21}B_{11} + A_{22}B_{21}$$

2-2) for the second column and second row

$$M1 - M2 + M3 + M6 =$$

$$\frac{A_{11}B_{22}+A_{22}B_{11}+A_{22}B_{22}-A_{22}B_{21}-A_{21}B_{11}-A_{22}B_{11}+A_{11}B_{12}-A_{11}B_{22}+A_{21}B_{11}+A_{21}B_{12}-A_{11}B_{11}-A_{22}B_{21}-A_{21}B_{22}+A_{21}B_{21}+A_{21}B_{22}+A_{21}B_{21}+A_{21}B_{22}+A$$

 $=A_{21}B_{12}+A_{22}B_{22}$

So we can have:

$$C = \begin{pmatrix} A_{11}B_{11} + A_{12}B_{21} & A_{11}B_{12} + A_{12}B_{22} \\ A_{21}B_{11} + A_{22}B_{21} & A_{21}B_{12} + A_{22}B_{22} \end{pmatrix}$$

b) Implement this method of blocked matrix multiplication and verify your code works by comparing it to the results of your programming language's built-in matrix product routine.

```
1000 import numpy as np
   def matrix_multiply(A, B):
        A11, A12 = A[0]
       A21, A22 = A[1]

B11, B12 = B[0]

B21, B22 = B[1]
       M1 = (A11 + A22) * (B11 + B22)
1008
       M2 = (A21 + A22) * B11
       M3 = A11 * (B12 - B22)
       M4 = A22 * (B21 - B11)
       M5 = (A11 + A12) * B22
       M6 = (A21 - A11) * (B11 + B12)
       M7 = (A12 - A22) * (B21 + B22)
        C11 = M1 + M4 - M5 + M7
        C12 = M3 + M5
        C21 = M2 + M4
        C22 = M1 - M2 + M3 + M6
        result = np.array([[C11, C12], [C21, C22]])
        return result
_{1024}ig|\# Test the function with example matrices A and B
C = matrix_multiply(A, B)
1028
   print("Matrix A:")
   print (A)
   print ("Matrix B:")
1032 print (B)
   print("Resultant Matrix C:")
1034 print (C)
   np.dot(A, B)
```

Listing 1: Python code For question 1b

By this code it should be able to test this function and compare with the built-in function of 'numpy' The out put is

```
Matrix A:
[[1 2]
[3 4]]
Matrix B:
[[5 6]
[7 8]]
Resultant Matrix C:
[[19 22]
[43 50]]
array([[19, 22],
[43, 50]])
```

C) Implement this method recursively such that it keeps subdividing the problem until the blocks of A and B are just scalars

```
import numpy as np
   def matrix_multiply_recursive(A, B):
        if len(A) == 1:
            return A * B
        n = len(A)
        half_n = n // 2
        A11 = A[:half_n, :half_n]
        A12 = A[:half_n, half_n:]
        A21 = A[half_n:, :half_n]
A22 = A[half_n:, half_n:]
        B11 = B[:half_n, :half_n]
        B12 = B[:half_n, half_n:]
        B21 = B[half_n:, :half_n]
B22 = B[half_n:, half_n:]
       M1 = matrix_multiply_recursive(A11 + A22, B11 + B22)
        M2 = matrix_multiply_recursive(A21 + A22, B11)
        M3 = matrix_multiply_recursive(A11, B12 - B22)
       M4 = matrix_multiply_recursive(A22, B21 - B11)
        M5 = matrix_multiply_recursive(A11 + A12, B22)
       M6 = matrix_multiply_recursive(A21 - A11, B11 + B12)
        M7 = matrix_multiply_recursive(A12 - A22, B21 + B22)
        C11 = M1 + M4 - M5 + M7
        C12 = M3 + M5
        C21 = M2 + M4
        C22 = M1 - M2 + M3 + M6
       C = np.vstack((np.hstack((C11, C12)), np.hstack((C21, C22))))
   # Test the function with example matrices A and B
1036 \, A = \text{np.array} ([[1, 2, 3, 4], [5, 6, 7, 8], [9, 10, 11, 12], [13, 14, 15, 16]])
   B = \text{np.array} ([[16, 15, 14, 13], [12, 11, 10, 9], [8, 7, 6, 5], [4, 3, 2, 1]])
C = matrix_multiply_recursive(A, B)
   print("Matrix A:")
1040
   print (A)
1042 print ("Matrix B:")
   print (B)
1044 print ("Resultant Matrix C:")
   print (C)
1046 # Perform matrix multiplication using NumPy's built-in function for comparison
   C_{\text{numpy}} = \text{np.dot}(A, B)
   print (C_numpy)
   # Check if the results are close
   print ("Are the results equal?", np. allclose (C, C_numpy))
```

Listing 2: Python code For question 1c

This code will allows recursive run the same process until they converge into 2x2 matrix, Then we can use a random matrix A and random matrix b to test whether if the function will gives the same answer as the build in function for numpy package By running the code above we get the output as:

```
Matrix A:
[[ 1 2 3 4]
  [ 5 6 7 8]
  [ 9 10 11 12]
  [13 14 15 16]]
```

```
Matrix B:

[[16 15 14 13]

[12 11 10 9]

[ 8 7 6 5]

[ 4 3 2 1]]

Resultant Matrix C:

[[ 80 70 60 50]

[240 214 188 162]

[400 358 316 274]

[560 502 444 386]]

[[ 80 70 60 50]

[240 214 188 162]

[400 358 316 274]

[560 502 444 386]]

Are the results equal? True
```

It shows out output is identical as the build in package , which means that our method works as we expected

We can actually try a larger 64x64 matrix,

```
# Generate random 64x64 matrices A and B
A = np.random.randint(1, 10, (64, 64))
B = np.random.randint(1, 10, (64, 64))

# Perform matrix multiplication using the recursive function
C_recursive = matrix_multiply_recursive(A, B)

# Perform matrix multiplication using NumPy's built—in function for comparison
C_numpy = np.dot(A, B)

# Check if the results are close
print("Are the results equal?", np.allclose(C_recursive, C_numpy))
```

Are the results equal? True

D) Numerically compare the asymptotic work for this method and naive matrix multiplication. Here is an exmaple code snippet of very naive matrix multiplication

```
import numpy as np

def matrix_multiply_naive(A, B):
    if A.shape[1] != B.shape[0]:
        raise ValueError("Number of columns in matrix A must be equal")

result = np.zeros((A.shape[0], B.shape[1]))

for i in range(A.shape[0]):
        for j in range(B.shape[1]):
            for k in range(A.shape[1]):
                result[i, j] += A[i, k] * B[k, j]

return result
```

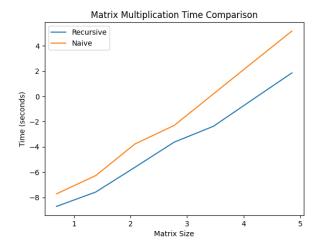


Figure 1: Time of matrix mutiple by two methods

```
import time
   import matplotlib.pyplot as plt
   # Compare execution times for different matrix sizes
   sizes = [2, 4, 8, 16, 32, 64, 128]
   recursive_times = []
   naive\_times = []
   for size in sizes:
        A = np.random.rand(size, size)
1008
       B = np.random.rand(size, size)
        start_time = time.time()
        matrix_multiply_recursive(A, B)
        end_time = time.time()
        recursive_times.append(end_time - start_time)
        start_time = time.time()
        matrix_multiply_naive(A, B)
        end_time = time.time()
1018
        naive_times.append(end_time - start_time)
   # Take the natural logarithm of the execution times
   ln_recursive_times = np.log(recursive_times)
   ln_naive_times = np.log(naive_times)
  | in_size=np.log(sizes)
   # Plot the results
   plt.plot(in_size , ln_recursive_times , label='Recursive')
plt.plot(in_size , ln_naive_times , label='Naive')
plt xlabel ('Matrix Size')
   plt.ylabel ('Time (seconds)')
   plt.title('Matrix Multiplication Time Comparison')
1030
   plt.legend()
   plt.show()
   # Estimate the slope using linear regression (numpy.polyfit)
| slope_recursive , _ = np.polyfit (in_size , ln_recursive_times , 1)
   slope_naive, _ = np.polyfit(in_size, ln_naive_times, 1)
   print(slope_recursive)
   print(slope_naive)
```

we get the slope, which is the big O of the two method equal to

- 2.841099045760702
- 3.637288115717954

Problem 2

In class we saw (or will see) that the classical Gram-Schmidt process for generating an orthonormal basis isn't numerically stable, so the "modified" Gram-Schmidt method is used instead.

a) Write a routine that takes in a matrix, A and returns a set of orthonormal vectors, Q using classical Gram-Schmidt.

```
# CGS
   def proj(v1, v2):
        return (numpy.dot(v2, v1) / numpy.dot(v1, v1)) * v1
1002
   def normalize(i):
1004
        normalized_vector = i / numpy.linalg.norm(i)
        return normalized_vector
   def cgs(A):
1008
       # Q_t is the transpose of final matrix / basis Q
       # store col vectors as row vectors in Q for easy modification
        Q_t = []
        for i in range(len(A)):
           # v_i
            a_i = A[i]
            v_i = A[i]
           # v_pt: previous v's transpose (row vector)
1018
            for v_pt in Q_t:
                 \# v_p: previous v's (col vector)
                v_p = numpy.transpose(v_pt)
                proj_vec = proj(v_p, a_i)
                v_i = v_i - proj_vec
           # append current v_i
            Q_t.append(numpy.transpose(v_i))
1028
       # print(Q_t)
1030
        for i in range(len(Q_t)):
            Q_t[i] = normalize(Q_t[i])
        Q = numpy.transpose(Q_t)
        return Q
```

Listing 3: Python code For question 2a

b) Write a routine that takes in a matrix, A and returns a set of orthonormal vectors, Q using modified Gram-Schmidt.

```
# MGS
    import numpy
   def proj(v1, v2):
        \texttt{return (numpy.dot(v2, v1) / numpy.dot(v1, v1))} * v1
1004
   def normalize(i):
        normalized_vector = i / numpy.linalg.norm(i)
        return normalized_vector
1010 def mgs(A):
       # Q-t is the transpose of final matrix / basis Q
       # use transpose of Q so each col vector becoms row vector and easier to access
        Q_t = numpy.transpose(A)
        for i in range(len(Q_t)):
            Q_t[i] = normalize(Q_t[i])
            q_i = numpy.transpose(Q_t[i])
1018
           \# subtract q_i's axis from rest q_k, k > i
            for k in range(i + 1, len(Q_-t)):
                q_k = numpy.transpose(Q_t[k])
                # subtract projection, already normlized
                q_k = q_k - numpy.dot(q_i, q_k) * q_i
                Q_t[k] = numpy.transpose(q_k)
       Q = numpy.transpose(Q_t)
        return Q
```

Listing 4: Python code For question 2b

- c) Recall that the generated matrix, Q is a normal matrix, that is, in exact arithmetic we get, $Q^T Q = I$. Take your two routines and check how accurate this is, i.e., compute $||Q^TQ I||$ and see how close it is to zero. Do this for the matrices,
- c-1) A = np.random.random((n, n))
- c-2) and A = 0.00001 * np.eye(n) + scipy.linalg.hilbert(n) for a few sizes, n.

Result in table format $(n \in 3, 5, 7, 9)$:

```
A = np.random.random((n, n))
          cgs result
                                 mgs result
______
      1.343782689223772e-15 | 5.822058658345461e-16
3 |
      5.079327213521194e-15 | 1.6425403990754258e-15
5 I
       1.29751252117977e-14 | 5.166999309333753e-15
7 I
      9.295267406282773e-14 |
                             3.2555245376247796e-14
A = 0.00001 * np.eye(n) + scipy.linalg.hilbert(n)
          cgs result
                     | mgs result
3 |
      7.080997315971274e-14
                               6.63317609727746e-15
5 I
      3.165790514913122e-08 |
                              8.512887360371612e-13
7 |
      8.243819382900969e-06 |
                             2.8665923273128248e-12
9 |
       0.000335415914015717
                             1.3617858663174204e-11
```

Code:

```
1000 # helper function
    def print_result(N, Error):
        print(" n | print("-" * 60)
                            cgs result
                                                        mgs result")
1004
        for i in range(len(Error)):
            print(f"{N[i]:2} | {Error[i][0]:24} | {Error[i][1]:24}")
1008 # main
   def main():
        \# random matrix, Q_-t*Q - I error test
       N = [3, 5, 7, 9]
        Errors_1 = []
        Errors_2 = []
        for n in N:
            # Random matrix of size n
            \ddot{A} = numpy.random.random((n, n))
1018
            # cgs() and mgs() returens result Q matrix
            result_cgs = cgs(A)
            result_mgs = mgs(A)
            \# \ |\,|\,QtQ\!\!=\!\!I\,|\,| \ \text{for cgs}
            error_cgs = numpy.linalg.norm( numpy.dot(numpy.transpose(result_cgs),
1024
        result_cgs) - numpy.identity(n))
            # ||QtQ=I|| for mgs
            error_mgs = numpy.linalg.norm( numpy.dot(numpy.transpose(result_mgs),
        result_mgs) - numpy.identity(n))
            Errors_1.append([error_cgs, error_mgs])
            # Random matrix of size n following the given equation
1030
            A = 0.00001 * numpy.eye(n) + scipy.linalg.hilbert(n)
            result_cgs = cgs(A)
            result_mgs = mgs(A)
            # || QtQ=I || for cgs
            error_cgs = numpy.linalg.norm( numpy.dot(numpy.transpose(result_cgs),
1036
        result_cgs) - numpy. identity(n))
            \# ||QtQ=I|| for mgs
            error_mgs = numpy.linalg.norm( numpy.dot(numpy.transpose(result_mgs),
        result_mgs) - numpy.identity(n))
            Errors_2.append([error_cgs, error_mgs])
1040
        print("A = np.random.random((n, n))")
        print_result(N, Errors_1)
        print("\nA = 0.00001 * np.eye(n) + scipy.linalg.hilbert(n)")
        print_result(N, Errors_2)
1046
1048 # call main function
   main()
```

Listing 5: Python code For question 2c

Problem 3

Suppose we have factored $A = QR \in \mathbb{R}^{n \times n}$, where Q is unitary, R is upper triangular. Let $u, v \in \mathbb{R}^n$.

a) Derive a method for computing the QR factorization of $\hat{A} = A + uv^T$ that doesn't involve forming \hat{A} and factoring it. *Hint*: multiply on the left by QT.

Let
$$\hat{A} = A + uv^T$$
, $A = QR \in \mathbb{R}^{nxn}$, $Q^TQ = I$, and $uv \in \mathbb{R}^n$.

$$\hat{A} = A + uv^T$$

$$Q^T \hat{A} = Q^T A + Q^T uv^T$$

$$Q^T \hat{A} = R + Q^T uv^T$$

$$Q^T \hat{A} = R + (Q^T u)v^T$$

Let $Q^T u = w$, since $Q^T u$ is an inner product, then we have w as col-vector ($O(n^2)$ to get w)

$$Q^T \hat{A} = R + wv^T$$

Apply rotations (denoted G_1) to w to rotate from $\begin{bmatrix} w_1 \\ w_2 \\ \dots \\ w_n \end{bmatrix}$ to $\begin{bmatrix} ||w|| \\ 0 \\ \dots \\ 0 \end{bmatrix}$ (O(1) to get each rotation in G_1)

$$G_1 Q^T \hat{A} = G_1 (R + wv^T)$$

$$G_1 Q^T \hat{A} = G_1 R + G_1 wv^T$$

$$G_1 Q^T \hat{A} = G_1 R + ||w|| e_1 v^T$$

($O(n^2)$ to apply rotation since they reduced to 2n row operation)

 G_1 R will produce an upper Hessenberg matrix.

Now let us apply n times Givens rotations (denoted G_2) to the G_1 R to ensure that it's upper triangular.

$$\hat{G}_{1}(Q^{T}\hat{A}) = \hat{G}_{1}R + ||w||e_{1}v^{T}$$

$$\hat{G}_{2}(\hat{G}_{1}(Q^{T}\hat{A})) = \hat{G}_{2}(\hat{G}_{1}R + ||w||e_{1}v^{T})$$

$$(\hat{G}_{2}\hat{G}_{1}Q^{T})\hat{A} = \hat{G}_{2}(\hat{G}_{1}R + ||w||e_{1}v^{T})$$

$$\hat{Q}^{T}\hat{A} = \hat{R}$$

$$\hat{A} = \hat{Q}\hat{R}$$

($O(n^2)$ to apply rotation since they reduced to 2n row operation)

From above we get:

$$\hat{Q}^T = \hat{G}_2 \hat{G} 1 Q^T$$

$$\hat{Q} = (\hat{G}_2 \hat{G} 1 Q^T)^T$$

$$\hat{R} = \hat{G}_2 (\hat{G}_1 R + ||w|| e_1 v^T)$$

From the above analyze, this method will have overall complexity of $O(n^2)$ witch is less than the naive approach with $O(n^3)$

9

b) Write a routine that does this and verify that it works by comparing it to naively computing a new QR of \hat{A} . Note that you might be off by a sign. This is fine because a QR decomposition is not unique. Also don't worry about having the most optimized code, just explain in (a) why the update is faster.

Code:

```
1000 import numpy as np
   import math
                     = helper functions =
1004
   def random_invertible_matrix(rank, min_val = -20, max_val = 20):
        while True:
           # Generate a random matrix of size n x n
            random_matrix = np.random.randint(min_val, max_val + 1, size=(rank, rank))
           # Check if the matrix is invertible
            if np.linalg.matrix_rank(random_matrix) == rank:
                return random_matrix
   def random_vector(rank, min_val = -20, max_val = 20):
        return np.random.randint(-20, 20 + 1, \text{ size} = (\text{rank}, 1)), np.random.randint(-20, 20, 20)
        +1, size = (rank, 1)
   def apply_rotation(matrix, cos, sin, axis_pos):
       # axis_pos used zero-outed row position
1018
        row_a = matrix[axis_pos - 1].copy()
       row_b = matrix[axis_pos].copy()
       # calculate new rows
          | CS
       # | -S C |
                   for rotating clockwise ans zero out 2nd term
        matrix[axis_pos - 1] = cos * row_a + sin * row_b
        matrix[axis_pos] = -sin * row_a + cos * row_b
1028
   def apply_givens_rotation_zero_out(matrix, row, col):
        a = matrix[row - 1][col]
        b = matrix [row][col]
        r = math.sqrt(a*a + b*b)
        \cos = a/r
        \sin = b/r
        row_a = matrix[row - 1].copy()
        row_b = matrix [row].copy()
        matrix[row - 1] = cos * row_a + sin * row_b
        matrix[row] = -sin * row_a + cos * row_b
        return cos, sin
                 generate matrix and vector =
1048 | rank = 4
   A = random_invertible_matrix(rank)
   u, v = random_vector(rank)
   v_t = np.transpose(v)
print ("Random Matrix A:")
   print (A)
1054 print ()
1056 # Get original QR decom.
   Q, R = np.linalg.qr(A)
1058
1060 # =
                main solution body =
```

```
# calculate (Q^T)u as w
1062
   Q_T = np. transpose(Q)
   w = np.matmul(Q_T, u)
1064
   # part 1 rotation, rotate u to be [[r] [0] ... [0]]
   rank = 4
1066
    Q_ht = Q_T.copy()
   R_h = R. copy()
1068
    for i in range (rank - 1, 0, -1):
        \# i is the row position of the term currently zeroing out
1070
        a = w[i - 1][0]
        \mathbf{b} = \mathbf{w}[\mathbf{i}][0]
        r = math.sqrt(a*a + b*b)
        \cos = a/r
        \sin = b/r
        # apply G (rotation matrix) to w (making sure rotation is correct
1078
        # apply_rotation(w, cos, sin, i)
        # print(np.transpose(w)), tested working
1080
        # rotate w
        w[i - 1][0] = r
        \mathbf{w}[\mathbf{i}][0] = 0
        # apply G (rotation matrix) to R-hat
1086
        apply_rotation(R_h, cos, sin, i)
        apply_rotation(Q_ht, cos, sin, i)
1088
   print("G_1*R Will be upper hesssenberg")
    print (R_h)
1092 print ()
1094 # add norm(w)elv_t to R_h
    e1 = np.zeros((rank, 1))
    e1[0][0] = 1
   R_h = R_h + np.matmul(w[0][0] * e1, v_t)
1098
   # part 2 rotation, zero out the R_ht to be upper triangular
   for i in range(1, rank):
        # zero out one term in R_h and save the rotation parameter c,s
        \cos, \sin = apply\_givens\_rotation\_zero\_out(R_h, i, i-1)
1104
        # apply same rotation to Q_ht
        apply_rotation(Q_ht, cos, sin, i)
|Q_h| = np.transpose(Q_ht)
1110
   |# =
                    = check result =
    # print result
print ("Result:")
   print ("Q_hat")
   print (Q_h)
    print ("R-hat")
1116 print (R_h)
A_hat_result = np.matmul(Q_h, R_h)
1120 \mid A_hat = A + np.matmul(u, v_t)
1122 print ()
    print("Original A_hat = A + u*v_t")
   print (A_hat)
    print("Re-composed A_hat = Q_hat * R_hat")
1126 print (A_hat_result)
```

Listing 6: Python code For question 3b

Output:

```
Random Matrix A:
[[-12 -9 -13 -15]
[-12 -17 0 9]
[ -8 10 -1 11]
[ 13 -11 19 13]]
G_1*R Will be upper hessenberg
[[ 18.41706898 -0.93508297 8.740993
                                     -7.19607331]
[-13.48375208 -7.87773959 -18.54205917 -21.17589946]
 [ -0.
            -22.97970493 10.43947085
                                     9.77154179]
 [ 0.
                         1.34332851 -0.56105418]]
Result:
Q_hat
[[ 0.05910791  0.00967877 -0.75042688  0.65823406]
[ 0.56574718 -0.69099139 -0.2760349 -0.35533895]
 [ 0.74307093  0.19514135  0.45552741  0.44973351]
 [-0.35253649 -0.69595767 0.39135535 0.48805935]]
R_hat
[-0.00000000e+00 2.33345973e+01 -1.42740098e+01 -1.36299233e+01]
 [ 0.00000000e+00 0.00000000e+00 2.08640562e+01 2.09346515e+01]
 [ 0.00000000e+00 0.00000000e+00 -2.22044605e-16 -1.91289693e+00]]
Original A_hat = A + u*v_t
[[ 28 25 -25 -21]
[ 268 221 -84 -33]
[ 352 316 -109 -43]
 [-167 -164 73 40]]
Re-composed A_hat = Q_hat * R_hat
[[ 28. 25. -25. -21.]
[ 268. 221. -84. -33.]
 [ 352. 316. -109. -43.]
[-167. -164. 73. 40.]]
```

Problem 4

The nth Legendre polynomial $P_n(x)$ is given by:

$$P_n(x) = \frac{1}{2^n} \sum_{k=0}^{\lfloor n/2 \rfloor} (-1)^k \binom{n}{k} (x^{n-2k})$$

a) Recall the classical way to compute the eigenvalues of a matrix is to look at the roots of the characteristic polynomial. Find a matrix who's characteristic polynomial is the n-th Legendre polynomial. Hint: The matrix should be upper Hessenberg.

$$A = \begin{bmatrix} 0 & 0 & 0 & 0 & \dots & 0 & -c_1 \\ 1 & 0 & 0 & 0 & \dots & 0 & -c_2 \\ 0 & 1 & 0 & 0 & \dots & 0 & -c_3 \\ \vdots & \vdots & \vdots & \vdots & \ddots & \vdots & \vdots \\ 0 & 0 & 0 & 0 & \dots & 0 & -c_{n-1} \\ 0 & 0 & 0 & 0 & \dots & 1 & -c_n \end{bmatrix}$$

where c_n is the n^{th} terms of the Legendre polynomial here we took the first 5^{th} Legendre polynomial as example The first five Legendre polynomials are given by:

$$\begin{aligned} P_0(x) &= 1 \\ P_1(x) &= x \\ P_2(x) &= \frac{1}{2}(3x^2 - 1) \\ P_3(x) &= \frac{1}{2}(5x^3 - 3x) \\ P_4(x) &= \frac{1}{8}(35x^4 - 30x^2 + 3) \\ P_5(x) &= \frac{1}{8}(63x^5 - 70x^3 + 15x) \end{aligned}$$

where for **2x2 matrix** there are $P_2(x) = \frac{1}{2}(3x^2 - 1) \propto x^2 - \frac{2}{3}$

$$A_{2\times 2} = \begin{bmatrix} 0 & 1\\ \frac{2}{3} & 0 \end{bmatrix}$$

For 3x3 matrix there are $P_3(x) = \frac{1}{2}(5x^3 - 3x) \propto x^3 - \frac{6}{5}$

$$A_{3\times3} = \begin{bmatrix} 0 & 1 & 0 \\ 0 & 0 & 1 \\ 0 & \frac{6}{5} & 0 \end{bmatrix}$$

For 4x4 matrix there are $P_4(x) = \frac{1}{8}(35x^4 - 30x^2 + 3) \propto x^4 - \frac{48}{7}x^2\frac{24}{35}$

$$A_{4\times4} = \begin{bmatrix} 0 & 1 & 0 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 1 \\ 0 & \frac{48}{7} & 0 & \frac{24}{35} \end{bmatrix}$$

For 5x5 matrix there are $P_5(x) = \frac{1}{8}(63x^5 - 70x^3 + 15x) \propto x^5 - \frac{560}{63}x^3 + \frac{40}{21}x^3$

$$A_{5\times5} = \begin{pmatrix} 0 & 1 & 0 & 0 & 0 \\ 0 & 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 0 & 1 \\ 0 & \frac{40}{21} & 0 & \frac{560}{63} & 0 \end{pmatrix}$$

During our research, we found ther are another idea presented by (Golub & Welsch, 1969b), where Any set of orthogonal polynomials, pj(.x)¥-i, satisfies a three term recurrence relationship:

$$P_i(x) = (a_i x + b_j) P_{j-i}(x) - C_j P_{j-2}(x)$$

where

$$j = 1, 2, \dots n; p - i(x) = 0, p0(x) = l,$$

then we have

$$xp(x) = Tp(x) + \frac{l}{a_N}p_N(x)e_N$$

Where T represents the tridiagonal matrix and $\mathbf{e}^T = (0, 0, \dots, 0, 1)\mathbf{r}$. Therefore, $p_n(\xi) = 0$ if and only if $\xi - p(\xi) = T(\xi)$ where ξ is an eigenvalue of the tridiagonal matrix T. In reference [12], it is demonstrated that T becomes symmetric when the polynomials are orthonormal. If T is not symmetric, then we can carry out a diagonal similarity transformation that will result in a symmetric tridiagonal matrix J.

$$J = \begin{bmatrix} a_1 & b_1 & 0 & \cdots & 0 \\ b_1 & a_2 & b_2 & \cdots & 0 \\ 0 & b_2 & a_3 & \ddots & \vdots \\ \vdots & \vdots & \ddots & \ddots & b_{n-1} \\ 0 & 0 & \cdots & b_{n-1} & a_n \end{bmatrix}$$

A Jacobi matrix, which is a symmetric tridiagonal matrix, corresponds to the characteristic polynomial that is a monic version (with a leading coefficient of 1) of the specific set of orthogonal polynomials being studied. For instance, if we consider the Legendre polynomials, we can start by noting that the monic Legendre polynomials follow a two-term recurrence relation:

$$\hat{P}_{n+1}(x) = x\hat{P}_n(x) - \frac{n^2}{4n^2 - 1}\hat{P}_{n-1}(x)$$

where $\hat{P}_n(x) = \frac{(n!)^2 2^n}{(2n)!} P_n(x)$ is the monic Legendre polynomial.

From this, we can derive an explicit expression for the corresponding Jacobi matrix (here I give the 5-by-5 case):

$$\begin{pmatrix}
0 & \frac{1}{\sqrt{3}} & 0 & 0 & 0\\
\frac{1}{\sqrt{3}} & 0 & \frac{2}{\sqrt{15}} & 0 & 0\\
0 & \frac{2}{\sqrt{15}} & 0 & \frac{3}{\sqrt{35}} & 0\\
0 & 0 & \frac{3}{\sqrt{35}} & 0 & \frac{4}{\sqrt{63}}\\
0 & 0 & 0 & \frac{4}{\sqrt{63}} & 0
\end{pmatrix}$$

 $\frac{n}{\sqrt{4n^2-1}}$ in the (n,n+1) and (n+1,n) positions, and 0 elsewhere. the eigenvector of these two different kinds of matrix will be somewhat identical to each other b) Write a routine to compute the eigenvalues of a upper Hessenberg matrix using the shifted QR method. It doesn't matter which shift you use but be wary that the Wilkonson shift because it might temporarily become complex before the method converges.

```
def QR_algorithm_Wilkinson_shift(A):
          "The QR algorithm with Wilkinson shift for finding eigenvalues.
           code-block:: matlab
            function [myeig tmvec] = QRwilkinson(A)
1006
           %———— Phase 1 -
           T = mytridiag(A);
                   - Phase 2
           m = length(A);
            myeig = zeros(m,1);
           % Size of tmvec is not known in advance.
           % Estimating an initial size.
1014
            tmvec = zeros(m*8,1);
           mu = 0;
            counter = 0;
1018
            while (m > 1)
                counter = counter + 1;
1020
                muMat = diag(mu * ones(m, 1));
                [Q,R] = myQR(T - muMat);
                T = R*Q + muMat;
                tmvec(counter) = abs(T(m,m-1));
                delta = (T(m-1,m-1) - T(m,m)) / 2;
                mu = T(m,m) - sign(delta) * T(m,m-1) / \dots
                        (abs(delta) + norm([delta T(m,m-1)]));
                if (tmvec(counter) < 1e-8)
1028
                    myeig(m) = T(m,m);
                    m = m - 1;
                    T = T(1:m, 1:m);
                end
            end
1034
            myeig(1) = T;
            tmvec = tmvec(1:counter);
            end
1038
```

c) Use your routine from (b) to compute the roots the Legendre polynomials for $n=1,\,2,\,3,\,4,\,5$. Compare your approximate roots to the actual roots (they can be found easily online).

Here we tried n=5

```
1000
           # Define the matrix elements
     elements = [
             \begin{bmatrix} 0 \ , \ 1/\text{np.sqrt}(3) \ , \ 0 \ , \ 0 \ , \ 0 \end{bmatrix} , \\ [1/\text{np.sqrt}(3) \ , \ 0 \ , \ 2/\text{np.sqrt}(15) \ , \ 0 \ , \ 0 ] \ , 
            [0, 2/\text{np.sqrt}(15), 0, 3/\text{np.sqrt}(35), 0],
1004
            [0, 0, 3/np.sqrt(35), 0, 4/np.sqrt(63)],
            [0, 0, 0, 4/\text{np.sqrt}(63), 0]
1006
     matrix = np.array([
            [0, 1, 0, 0, 0],
             \begin{bmatrix} 0 & 0 & 1 & 0 & 0 \end{bmatrix}, \\ \begin{bmatrix} 0 & 0 & 1 & 0 & 0 \end{bmatrix}, \\ \begin{bmatrix} 0 & 0 & 0 & 1 & 0 \end{bmatrix}, \\ \begin{bmatrix} 0 & 0 & 0 & 0 & 0 \end{bmatrix}, \\ \end{bmatrix} 
            [0, 40/21, 0, 560/63, 0]
     # Convert the list of lists to a NumPy matrix
      A_matrix = np.array(elements)
1018
     B_matrix = np.array(elements)
     eigenvalues_shifted_qr = QR_algorithm_Wilkinson_shift(A_matrix)
1020
     print("Eigenvalues from Shifted QR Method:", eigenvalues_shifted_qr)
     eigenvalues_shifted_qr = QR_algorithm_Wilkinson_shift(B_matrix)
     print ("Eigenvalues from Shifted QR Method:", eigenvalues_shifted_qr)
```

Eigenvalues from Shifted QR Method: [-9.06179846e-01 -5.38469310e-01 1.78162900e-17 9.06179846e-01 5.38469310e-01]

Eigenvalues from Shifted QR Method: [-9.06179846e-01 -5.38469310e-01 1.78162900e-17 9.06179846e-01 5.38469310e-01]

We can compare this with the roots online:

Number of points, <i>n</i>	Points, x_i		Weights, w _i	
1	0		2	
2	$\pm \frac{1}{\sqrt{3}}$	±0.57735	1	
3	0		$\frac{8}{9}$	0.888889
	$\pm\sqrt{rac{3}{5}}$	±0.774597	$\frac{5}{9}$	0.55556
4	$\pm\sqrt{\frac{3}{7}-\frac{2}{7}\sqrt{\frac{6}{5}}}$	±0.339981	$\frac{18+\sqrt{30}}{36}$	0.652145
	$\pm\sqrt{\frac{3}{7}+\frac{2}{7}\sqrt{\frac{6}{5}}}$	±0.861136	$\frac{18-\sqrt{30}}{36}$	0.347855
5	0		$\frac{128}{225}$	0.568889
	$\pm\frac{1}{3}\sqrt{5-2\sqrt{\frac{10}{7}}}$	±0.538469	$\frac{322+13\sqrt{70}}{900}$	0.478629
	$\pm\frac{1}{3}\sqrt{5+2\sqrt{\frac{10}{7}}}$	±0.90618	$\frac{322 - 13\sqrt{70}}{900}$	0.236927

Figure 2: Gauss–Legendre quadrature

References

[1] Hbldh. (n.d.). $b2ac/b2ac/eigenmethods/qr_algorithm.py$ at $master \cdot hbldh/b2ac$. GitHub. https://github.com/hbldh/b2ac/blob/master/b2ac/eigenmethods/qr_algorithm.py

[2] Wikipedia contributors. (2023). Gauss-Legendre quadrature. Wikipedia. $\verb|https://en.wikipedia.org/wiki/Gauss\%E2\%80\%93Legendre_quadrature|$

[3] Golub, G. H., & Welsch, J. H. (1969). Calculation of Gauss quadrature Rules. *Mathematics of Computation*, 23 (106), 221.

https://doi.org/10.2307/2004418