HW5

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0.1 ### HW 5

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I worked by myself on this assignment.

```
[1]: import numpy as np
from scipy.io import mmread
import timeit
import warnings

# I was getting a RuntimeWarning: overflow encountered in true_divide in

→ jacobi's method
warnings.filterwarnings('ignore')
```

```
[2]:
    read_matrix
        file_name: file where the matrix resides

Reads the matrix in from memory using numpy

def read_matrix(file_name):
    # https://numpy.org/doc/stable/reference/generated/numpy.loadtxt.html
    ext = file_name[-3:]
    if ext == "txt":
        return np.loadtxt(file_name,delimiter=',')
    elif ext == "mtx":
        return mmread(file_name).A
```

```
[3]:  
Question #1 functions:

generate_ones
    matrix: a `n x m` matrix stored in a np.ndarray
    -----
    RETURNS: a `n x 1` np.ndarray of all ones using numpy
```

```
def generate_ones(matrix):
         # https://numpy.orq/doc/stable/reference/qenerated/numpy.ones.html
         return np.ones((matrix.shape[0], 1), dtype=type(matrix[0][0]))
[4]: \ \ '''
     Question #3 functions:
     lu
         A: an `n x m` matrix stored in a np.ndarray
         RETURNS: a tuple containing the L and U matrices obtained from performing \Box
      \hookrightarrow LU decomp.
     lu_solve
         A: an `n x m` matrix stored in a np.ndarray
         b: an `n x 1` matrix stored in a np.ndarray
         RETURNS: an `n x 1` matrix stored in a np.ndarray containing a solution x_{\sqcup}
      \hookrightarrow to the equation Ax=b
     cholesky
         A: an `n x m` matrix stored in a np.ndarray
         RETURNS: the L matrix obtained from performing cholesky decomp.
     cholesky_solve
         A: an `n x m` matrix stored in a np.ndarray
         b: an `n x 1` matrix stored in a np.ndarray
         RETURNS: an `n x 1` matrix stored in a np.ndarray containing a solution x_{\sqcup}
      \hookrightarrow to the equation Ax=b
     AS FOR:
         forward_sub
         back\_sub
         Although I roughly understood what these algorithms were doing going into \Box
      \hookrightarrow the assignment,
         https://johnfoster.pqe.utexas.edu/numerical-methods-book/LinearAlgebra_LU.
      \hookrightarrow html was incredibly
          useful\ because\ I\ didn't\ want\ to\ spend\ considerable\ time\ implementing\ these
      ⇔functions because we could've
         simply used a library function to solve.
     111
     def p3_solve(A,b):
```

a = np.array(A, dtype=np.float128)

```
if is_symmetric(a):
        return cholesky_solve(a,b)
    return lu_solve(a,b)
def lu(A):
    # n = number of rows
    n = A.shape[0]
   U = A.copy()
    L = np.eye(n, dtype=np.double)
    #Loop over rows
    for i in range(n):
        #Eliminate entries below i with row operations
        #on U and reverse the row operations to
        #manipulate L
        factor = U[i+1:, i] / U[i, i]
        L[i+1:, i] = factor
        U[i+1:] -= factor[:, np.newaxis] * U[i]
    return L, U
def lu_solve(A, b):
   L, U = lu(A)
    # solve Ly=b for y using forward sub
    y = forward_sub(L, b)
    # then solve Ux=y for x using back sub
    return back_sub(U, y)
    # seemed rather consistent with the scipy method so I kept the forward and
\hookrightarrow backward that I have.
    # return solve_triangular(U, b)
def cholesky(A):
   # n = number of rows
    n = A.shape[0]
   L = np.zeros((n, n), dtype=np.double)
    for k in range(n):
        L[k, k] = np.sqrt(A[k, k] - np.sum(L[k, :] ** 2))
        L[(k+1):, k] = (A[(k+1):, k] - L[(k+1):, :] @ L[:, k]) / L[k, k]
    return L
def cholesky_solve(A, b):
   L = cholesky(A)
    # just special case of LU decomp where U = L.T
    y = forward_sub(L, b)
    return back_sub(L.T, y)
def forward_sub(L, b):
    # n = number of rows
   n = L.shape[0]
```

```
# allocating space for the solution vector
   y = np.zeros_like(b, dtype=np.double);
    #Here we perform the forward-substitution.
    #Initializing with the first row.
   y[0] = b[0] / L[0, 0]
    #Looping over rows in reverse (from the bottom up),
   #starting with the second to last row, because the
   #last row solve was completed in the last step.
   for i in range(1, n):
       y[i] = (b[i] - np.dot(L[i,:i], y[:i])) / L[i,i]
   return v
def back_sub(U, y):
   # n = number of rows
   n = U.shape[0]
   # allocating space for the solution vector
   x = np.zeros_like(y, dtype=np.double);
    #Here we perform the back-substitution.
   #Initializing with the last row.
   x[-1] = y[-1] / U[-1, -1]
   #Looping over rows in reverse (from the bottom up),
   #starting with the second to last row, because the
   #last row solve was completed in the last step.
   for i in range(n-2, -1, -1):
        x[i] = (y[i] - np.dot(U[i,i:], x[i:])) / U[i,i]
   return x
```

```
[5]: '''
     Question #4 functions:
     jacobi solve
         A: an `n x m` matrix stored in a np.ndarray
         b: an `n x 1` matrix stored in a np.ndarray
         max\_iters=25: can change the number of max iterations the method will use,
      \hookrightarrow25 by default
         x=None: can supply an inital quess for the solution x, I initialize x to be_{\square}
      \rightarrowall ones based on A when x==None
         RETURNS: an `n x 1` matrix stored in a np.ndarray containing a solution x_{\sqcup}
      \hookrightarrow to the equation Ax=b using
                   jacobi's iterative method
     111
     def jacobi_solve(A, b, max_iters=25, x=None):
          """Solves the equation Ax=b via the Jacobi iterative method."""
         # Create an initial quess if needed
                                                                                            Ш
                                                                                            Ш
```

```
if x is None:
    x = generate ones(A)
# Create a vector of the diagonal elements of A
# and subtract them from A
                                                                             1.1
                                                                             Ш
D = np.diag(A)
R = A - np.diagflat(D)
try:
    # Iterate for max iters times
    for i in range(max_iters):
        temp = x
        x = (b - np.dot(R,x)) / D
# p5 giving me problems
except np.linalg.LinAlgError:
    return temp
return x
```

```
[6]: '''
     Question #5 functions:
     gauss_seidel_solve
         A: an `n x m` matrix stored in a np.ndarray
         b: an `n x 1` matrix stored in a np.ndarray
         max\_iters=25: can change the number of max iterations the method will use, \sqcup
      \hookrightarrow25 by default
          x=None: can supply an inital quess for the solution x, I initialize x to be_{\sqcup}
      \rightarrowall ones based on A when x==None
         RETURNS: an `n x 1` matrix stored in a np.ndarray containing a solution x_{\sqcup}
      \hookrightarrow to the equation Ax=b using
                   gauss/seidel's iterative method
     I I I
     def gauss_seidel_solve(A, b, num_iters=25, x=None):
         # Create an initial guess if needed
         if x is None:
              x = np.ones((A.shape[0], 1))
         L = np.tril(A)
         U = A - L
         try:
```

```
# Iterate for num_iters times
for i in range(num_iters):
    temp = x
    x = np.dot(np.linalg.inv(L), b - np.dot(U, x))
# p5 giving me problems
except np.linalg.LinAlgError:
    return temp
return x
```

```
[7]: '''
     OTHER RELEVANT HELPER FUNCTIONS:
         relative_error
             truth: `n x m` matrix that is claiming to be the absolute correct values
             sol: `n x m` matrix that is any arbitrary solution obtained
             RETURNS: the relative error of the arbitrary solution to the true\sqcup
      \hookrightarrow solution
         is_symmetric
             matrix: `n x m` matrix
             RETURNS: true if the matrix is symmetric, and false otherwise
     def relative_error(truth, sol):
         return np.linalg.norm(sol - truth) / np.linalg.norm(sol)
     def is_symmetric(matrix):
         if matrix.shape[0] == matrix.shape[1]:
             if isinstance(matrix[0][0], float):
                 return np.allclose(matrix,matrix.T, atol=1e-06)
             return np.array_equal(matrix,matrix.T)
         return False
```

```
[8]:
    TIMING MODULE:
    https://www.pythoncentral.io/time-a-python-function/

    def wrapper(func, *args, **kwargs):
        def wrapped():
            return func(*args, **kwargs)
        return wrapped
```

```
[9]: num_matrices = 4
     num_tests = 50
     directory = "/Users/AndrewMacbook/Downloads/"
     for i in range(num_matrices):
         file name = "{}mat{}-1.txt".format(directory, i+1)
         matrix = read_matrix(file_name)
         b = generate_ones(matrix)
         truth = np.linalg.solve(matrix, b)
         rel_error_tru = relative_error(truth, truth)
         rel_error_lu = relative_error(truth, p3_solve(matrix, b))
         rel_error_jac = relative_error(truth, jacobi_solve(matrix, b))
         rel_error_gs = relative_error(truth, gauss_seidel_solve(matrix, b))
         linalg_solved
                            = wrapper(np.linalg.solve, matrix, b)
         luCholesky_solved = wrapper(p3_solve, matrix, b)
                            = wrapper(jacobi_solve, matrix, b)
         jacobi solved
         gaussSeidel_solved = wrapper(gauss_seidel_solve, matrix, b)
         print("{} Matrix {} {}".format("#" * 12, i+1, "#" * 12))
         print("{:<8} time: {:.6e}, error: {:.6e}".format("linsol",timeit.</pre>
      -timeit(linalg_solved, number=num_tests) / num_tests, rel_error_tru))
         print("{:<8} time: {:.6e}, error: {:.6e}".format("LU/Chol",timeit.</pre>
      -timeit(luCholesky_solved, number=num_tests) / num_tests, rel_error_lu))
         print("{:<8} time: {:.6e}, error: {:.6e}".format("Jacobi",timeit.</pre>
      -timeit(jacobi_solved, number=num_tests) / num_tests, rel_error_jac))
         print("{:<8} time: {:.6e}, error: {:.6e}".format("GS", timeit.</pre>
      →timeit(gaussSeidel_solved, number=num_tests) / num_tests, rel_error_gs))
```



```
time: 1.087941e-03, error: 0.000000e+00
linsol
LU/Chol time: 2.189109e-02, error: 3.671191e-01
       time: 2.531968e-02, error: 1.704764e-04
Jacobi
       time: 8.308061e-02, error: 1.500389e-06
time: 3.455016e-03, error: 0.000000e+00
linsol
LU/Chol time: 3.668489e-02, error: 2.188609e-02
Jacobi
      time: 9.041695e-02, error: 1.177007e-12
       time: 2.664884e-01, error: 2.635684e-09
time: 9.452872e-04, error: 0.000000e+00
linsol
LU/Chol time: 5.007002e-02, error: 1.002810e-05
       time: 2.076657e-02, error: nan
Jacobi
       time: 1.262019e-03, error: 2.126029e+00
```

```
linsol time: 6.492302e-05, error: 0.000000e+00 LU/Chol time: 3.049418e-03, error: 5.046800e-01 Jacobi time: 8.678296e-04, error: 9.999775e-01 GS time: 3.733116e-03, error: 1.355951e-02
```

0.1.1 EXTRA CREDIT:

1. dwa512: Square Dielectric Waveguide

Not entirely sure what that means, but from the 'structure plot', it looked like a matrix that would be non-singular, and that the iterative methods would shine. Unfortunately the latter proved to be false.

```
[10]: names = ["dwa512"]
      for i in range(len(names)):
          file name = "{}{}.mtx".format(directory, names[i])
          matrix = read_matrix(file_name)
          b = generate_ones(matrix)
          truth = np.linalg.solve(matrix, b)
          rel_error_tru = relative_error(truth, truth)
          rel_error_lu = relative_error(truth, p3_solve(matrix, b))
          rel_error_jac = relative_error(truth, jacobi_solve(matrix, b))
          rel_error_gs = relative_error(truth, gauss_seidel_solve(matrix, b))
          linalg solved
                             = wrapper(np.linalg.solve, matrix, b)
          luCholesky_solved = wrapper(p3_solve, matrix, b)
                             = wrapper(jacobi_solve, matrix, b)
          jacobi solved
          gaussSeidel_solved = wrapper(gauss_seidel_solve, matrix, b)
          print("{} {:^20} {}".format("#" * 12, names[i], "#" * 12))
          print("{:<8} time: {:.6e}, error: {:.6e}".format("linsol", timeit.</pre>
       →timeit(linalg_solved, number=num_tests) / num_tests, rel_error_tru))
          print("{:<8} time: {:.6e}, error: {:.6e}".format("LU/Chol",timeit.</pre>
       →timeit(luCholesky_solved, number=num_tests) / num_tests, rel_error_lu))
          print("{:<8} time: {:.6e}, error: {:.6e}".format("Jacobi",timeit.</pre>
       →timeit(jacobi_solved, number=num_tests) / num_tests, rel_error_jac))
          print("{:<8} time: {:.6e}, error: {:.6e}".format("GS", timeit.</pre>
       →timeit(gaussSeidel_solved, number=num_tests) / num_tests, rel_error_gs))
```

########## dwa512 ######### linsol time: 6.648570e-03, error: 0.000000e+00 LU/Chol time: 6.648564e-01, error: 3.991839e-12 Jacobi time: 1.589301e-01, error: 1.000000e+00 GS time: 3.930483e-01, error: 1.000000e+00

0.1.2 Sources used:

https://courses.grainger.illinois.edu/cs357/sp2020/notes/ref-9-linsys.html

http://www.cim.mcgill.ca/~derek/ecsex43/A1.md.html

https://numpy.org

https://www.pythoncentral.io/time-a-python-function/

https://stackoverflow.com/questions/7370801/how-to-measure-elapsed-time-in-python

https://math.nist.gov/MatrixMarket/data/NEP/dwave/dwa512.html