## 1 GMRES

1. **Arnoldi step:** For given orthonormal vectors  $q_1, \ldots, q_{r-1} \in \mathbb{R}^n$  considered as a matrix  $Q_{r-1} = [q_1, \ldots, q_{r-1}] \in \mathbb{R}^{n \times (r-1)}$ , and a vector  $v \in \mathbb{R}^n$ , implement a helper function

$$q_r, h_r := Arnoldi\_step(Q_{r-1}, v),$$

which, according to the Arnoldi iteration, appends these vectors (i.e., the matrix  $Q_{r-1}$ ) by an orthonormal vector  $q_r$  through orthogonalizing v against  $q_1, \ldots, q_{r-1}$  and also outputs the numbers  $h_r := (h_{1,r-1}, \ldots, h_{r,r-1})^\top \in \mathbb{R}^r$ , where  $h_{\ell,r-1} := q_\ell^\top v$  for  $\ell \le r$ . You can then call Arnoldi\_step(Q; v) within GMRES(...).

2. GMRES: Implement a function

$$x = GMRES(A, b, x0, tol=1e-6, maxiter=None, N=None),$$

which takes as arguments

- A : a function evaluating the matrix-vector product  $v \mapsto A \cdot v$  for some matrix  $A \in \mathbb{R}^{n \times n}$  (not as an array!)
- ullet b : a vector  $b \in \mathbb{R}^n$
- x0: an arbitrary initial guess  $x^0 \in \mathbb{R}^n$
- $\bullet$  tol : error tolerance as float, which is set to  $10^{-6}$  by default
- maxiter : optional maximum number of iterations, which is set to None by default
- N : optional preconditioner as a function (not as an array), for which  $N(v) \approx A^{-1}v$

and then solves the system Ax=b by applying the GMRES method as presented in the lecture (see pseudocode 1 below). It shall then return

• x : the approximation to the solution.

The iteration shall break if the residual is tolerably small, i.e.,

$$||Ax^k - b||_2 < \text{tol}$$

or the maximum number of iterations maxiter has been reached.

3. Test your solver on a random invertible tridiagonal matrix

$$A = \begin{pmatrix} * & * & 0 & \cdots & 0 \\ * & * & * & \ddots & \vdots \\ 0 & \ddots & \ddots & \ddots & 0 \\ \vdots & \ddots & * & * & * \\ 0 & \cdots & 0 & * & * \end{pmatrix} \in \mathbb{R}^{n \times n}$$

and some right-hand side b and initial guess  $x_0$  of your choice. Choose different n and check how many iterations you need (potentially many!).

Hint: You can generate some random diagonals using numpy.random.rand(n) and then use

to construct a sparse CSR matrix. You can add np.ones(n) to the main diagonal in order to "strengthen" the diagonal of A and thereby to get a better conditioned system. Then implement a function A(x) that outputs the matrix-vector product A.dot(x).

- 4. **Preconditioner:** For the same system, run your GMRES routine with a preconditioner of your choice (for example Jacobi  $N: v \mapsto D^{-1}v$ ). Do you observe any difference in the number of iterations needed? (Don't worry if not!)
- 5. Compare your GMRES solver to scipy.linalg.solve(A\_csr.toarray(),b) for large dimension  $n \ge 10^5$  and measure the time needed in each case. Also, find a SciPy implementation of GMRES and compare to yours.

```
1 INPUT: A \in GL_n(\mathbb{R}), b \in \mathbb{R}^n
 2 OUTPUT: approximation x_r \in K_r(A, b) to the exact solution A^{-1}b
 4 GMRES(A, b, x_0 = 0, \text{ tol = 1e-6, maxiter=None, } N = I):
 5 b := b - Ax_0 //account for initial guess
 6 A := NA, b := Nb //account for preconditioner
 7 //Initialization:
 8 q_1 := \frac{b}{\|b\|_2}, Q_1 := [q_1]
 9 v := Aq_1
10 	ilde{q}_1 := rac{v}{\|v\|_2}, 	ilde{Q}_1 := [	ilde{q}_1], 	ilde{R}_1 = [\|v\|_2]
11 for r = 2, ..., \min(n, \max ter) do
        //STEP 1: use Arnoldi to find column q_r by orthogonalizing v against q_1, \ldots, q_{r-1}
12
        q_r, h_{r-1} := Arnoldi\_step(Q_{r-1}; v) //we don't need h_{r-1} in this variant
13
14
        Q_r := [Q_{r-1}, q_r]
        v := Aq_r
15
16
        //STEP 2: use Arnoldi to find columns \tilde{q}_r, \tilde{r}_r by orthogonalizing v against \tilde{q}_1, \ldots, \tilde{q}_{r-1}
17
        \tilde{q}_r, \tilde{r}_r := \text{Arnoldi\_step}(\tilde{Q}_{r-1}; v)
        Q_r := [Q_{r-1}, \widetilde{q}_r], R_r := [R_{r-1}, \widetilde{r}_r]
18
19
        //STEP 3: solve auxiliary least squares problems to obtain coordinates
20
        c_r := \text{solve\_triangular}(\widetilde{R}_r, \widetilde{Q}_r^\top b)
21
22
        x_r := Q_r c_r
        //Attention: Evaluate the original residual here:
23
        if ||N^{-1}(Ax_r - b)||_2 < \text{tol then}
24
         break
25
        end
26
27 end
28 return x_0 + x_r
```

## Algorithm 1: GMRES

## Solution:

```
Parameters:
   Q : matrix containing the orthonormal vectors as columns
   Returns:
   qr : vector which is orthonormal to all columns in Q
   hr : vector containing the linear coefficients so that
        v = [Q,qr] @ hr
   r = np.shape(Q)[1]
   hr = np.zeros(r+1)
   for j in range(r):
       hr[j] = np.dot(Q[:, j], v)
       v = v - hr[j] * Q[:, j]
   hr[-1] = norm(v)
   return v / hr[-1], hr
def Lanczos_step(Q, v):
   r = np.shape(Q)[1]
   hr = np.zeros(r+1)
   for j in range(max(0, r-2), r):
       hr[j] = np.dot(Q[:, j], v)
       v = v - hr[j] * Q[:, j]
   hr[-1] = norm(v)
   return v / hr[-1], hr
# ----- #
# GMRES
def GMRES(A, b, x0, tol=1e-6, maxiter=None, N=None, sym=False):
   solves a system Ax = b, where A is assumed to be invertible,
   using QR-based GMRES
   Parameters
   A : python function
       for evaluating the matrix-vector product
   b : (n,) numpy.ndarray
        right-hand side
   x0: (n,) numpy.ndarray
        initial guess
   tol : float
         iteration stops if ||Axk - b|| < tol, tol = 1e-6
   maxiter : int (optional)
            maximum number of iterations
   N : python function (optional)
       for evaluation matrix-vector product of preconditioner
   sym : bool (optional)
         indicating whether A is symmetric or not
         if sym=True: Lanczos is used over Arnoldi
   Returns
   x : (n,) numpy.ndarray
       approximate solution to Ax=b, with ||Ax-b||<tol
   info : dict
```

```
# account for initial guess
    boriginal = b
    b = boriginal - A(x0)
    # account for preconditioner A(v) := A(N(v))
    if N:
        Aoriginal = A
        def A(x):
           return N(Aoriginal(x))
        b = N(b)
    # Initializing
    Q = b / norm(b)
    Q = Q[:, np.newaxis]
    v = A(Q[:, 0])
    tQ, tR = v / norm(v), norm(v)
    tQ = tQ[:, np.newaxis]
    n = len(b)
    xr = np.zeros(n)
    # if A is symmetric we (can) use Lanczos instead of Arnoldi
        def ortho(Q, v):
            return Lanczos_step(Q, v)
    else:
        def ortho(Q, v):
            return Arnoldi_step(Q, v)
    if maxiter:
       maxiter = min(maxiter, n+1)
    else:
        maxiter = n+1
    for r in range(1, maxiter):
        # STEP 1: Find next orthonormal basis vector of Krylov subspace
        qr, hr = ortho(Q, v) # we do not use hr in our variant
        Q = np.hstack((Q, qr[:, np.newaxis]))
        v = A(Q[:, -1])
        # STEP 2: Find QR decomposition of AQ_r by appending previous one
        tqr, trr = ortho(tQ, v)
        tQ = np.hstack((tQ, tqr[:, np.newaxis]))
        tR = np.hstack((np.vstack((tR, np.zeros((1, r)))), trr[:, np.newaxis]))
        # STEP 3: Solve least squares problem involving
               AQ_k using its QR-decomposition
        cr = linalg.solve_triangular(tR, tQ.T@b)
        xr = Q @ cr
        # Evaluate current Error and break if its small enough
        lsq_err = norm(boriginal - Aoriginal(x0+xr))
        # collect some infos
        info = dict(iterationCount=r+1,
                    dimension=n,
                    residualNorm=lsq_err)
        if lsq_err < tol:</pre>
           break
    return x0 + xr, info
def main():
```

```
# ----- #
# SETTING
# ----- #
# generate the random system matrix and rhs, as well as initial guess
start_time = time()
n = 10000 # 10000 # 100000
A_{sparseMatrix} = (2 * sparse.eye(n, k=0) -
                   1 * sparse.eye(n , k=1) -
                   1 * sparse.eye(n , k=-1))
# we regularize the matrix to prevent ill-conditioned systems
diagonals = [np.random.rand(n) + 1.5 * np.ones(n),
            np.random.rand(n-1),
            np.random.rand(n-1)
A\_sparseMatrix = scipy.sparse.diags(diagonals, [0, 1, -1]).tocsr()
print("\n Dimension:", n)
print(f" Time to generate the matrix: {time()-start_time:0.2f} [s]")
def A_function(x): # The matrix vector product as a function
   return A_sparseMatrix.dot(x)
b = np.random.rand(n) # np.ones(n) #
x0 = np.random.rand(n) # np.zeros(n) # b#*100 #
# GMRES parameters
tol = 1e-06
maxiter = 1000
sym = False
# runtime parameter
compare_LU_dense = 1
compare_Scipy = 0
for preconditionerChoice in ["none", "Jacobi", "GS"]:
   # ----- #
    # Preconditioner
   # ---
   # Jacobi
    precondJacobiArray = sparse.diags(1. / A_sparseMatrix.diagonal())
   def precondJacobi(b): return precondJacobiArray.dot(b)
   def precondNone(b): return b
   # Gauss-Seidel
    precondGSinvArray = scipy.sparse.tril(A_sparseMatrix).tocsr()
   def precondGS(b): return spla.spsolve(precondGSinvArray, b)
   # choice
    preconditionerFunctionDict = {"none": precondNone,
                                "Jacobi": precondJacobi,
                               "GS": precondGS}
    preconditionerFunction = \
                         preconditionerFunctionDict[preconditionerChoice]
    # ----- #
    # Our GMRES
    # ----- #
    print("-" * 30 + "\n\t Our GMRES \n" + "-" * 30)
    start_time = time()
   xk, info = GMRES(A_function, b, x0=x0, tol=tol,
                   maxiter=maxiter, N=preconditionerFunction, sym=sym)
    print(f" Solving time = {time()-start_time:0.2f} [s]")
    print(" Number of iterations:\t", info["iterationCount"],
         f"\n Residual norm: {info['residualNorm']:0.2e}\n")
    print(" 'Ax = b' is", np.allclose(A_function(xk), b))
```

```
if compare_Scipy:
            # ----- #
            # Scipy's GMRES
            # ----- #
            print("-" * 30 + "\n\t\t SciPy's GMRES \n" + "-" * 30)
            # callback for gmres
            class gmres_counter(object):
                def __init__(self):
                     self.niter = 0
                 def __call__(self, rk=None):
                     self.niter += 1
            counter = gmres_counter()
            start_time = time()
            x, eC = \setminus
            {\tt spla.gmres(A\_sparseMatrix\,,\,b,\,x0=x0\,,\,tol=tol\,,\,maxiter=maxiter\,,}
                        M=spla.LinearOperator((n, n), preconditionerFunction),
                        callback=counter)
            print(" Successful exit: ", eC == 0)
print(f"\n Solving time = {time()-start_time:0.2f} [s]")
            print(" Number of iterations:\t", counter.niter)
print(f" Residual norm: {norm(b - A_sparseMatrix.dot(x)):0.2e}")
             print("\n 'Ax = b' is", np.allclose(A_sparseMatrix.dot(x), b))
    # Compare to Scipy's LU (dense)
    if compare_LU_dense:
        print("\n" + "-" * 30 + "\n\t\t LU dense \n" + "-" * 30)
        start_time = time()
        xk = linalg.solve(A_sparseMatrix.toarray(), b)
        print(f" Solving time = {time()-start_time:0.2f} [s]")
if __name__ == "__main__":
main()
```