



- 1 a) Since A is normal, it can be orthogonally diagonalized such that $A = Q\Lambda Q^T$, where the columns of Q are the eigenvectors q_i , and the entries of Λ are the eigenvalues λ_i of A .

First, assume that $x^T Ax > 0$ for all $x \neq 0$. This must be true for $x = q_i$, so $q_i^T A q_i = \lambda_i q_i^T q_i = \lambda_i \|q_i\|_2^2 > 0$. We have proved that if A is positive definite, then all the eigenvalues are positive.

Next, assume that the eigenvalues of A are positive. Since the eigenvectors q_i form a basis for \mathbb{R}^n , we can represent any vector as a linear combination of the q_i . Let $u = Qc$ be any nonzero vector in \mathbb{R}^n . Then $u^T Au = c^T Q^T Q \Lambda Q^T Q c = c^T \Lambda c = \sum_i \lambda_i c_i^2$. All the λ_i are positive, and at least one of the c_i are nonzero, so $u^T Au > 0$. Thus, we have proved the other direction.

- b) The matrix A is normal, so it can be unitarily diagonalized, $A = Q\Lambda Q^H$. Thus, $AA^H = Q\Lambda\Lambda^H Q^H = Q\Lambda^H\Lambda Q^H = A^H A$. We can then calculate the condition number:

$$\kappa(A) = \|A\|_2 \|A^{-1}\|_2 = \sqrt{\rho(AA^H)} \sqrt{\rho(A^{-1}A^{-H})} = \max_i |\lambda_i| \max_i \frac{1}{|\lambda_i|} = \frac{\max_i |\lambda_i|}{\min_i |\lambda_i|}.$$

- c) Assume first that A is positive definite. Then A is non-singular, so given a vector $y \neq 0$, we can find a unique $x \neq 0$ so that $y = Ax$. Thus, $y^T A^{-1} y = x^T A^T x = x^T Ax > 0$. To prove the converse, just substitute $A \leftrightarrow A^{-1}$.
- d) Since A is normal, we have that $A = Q\Lambda Q^T$ with $Q^T Q = Q Q^T = I$. Hence, we can express the Rayleigh quotient as

$$R(x) = \frac{x^T Ax}{x^T x} = \frac{x^T Q \Lambda Q^T x}{x^T Q Q^T x}.$$

If we now define $y = Q^T x$, we have

$$R = \frac{y^T \Lambda y}{y^T y} = \frac{\sum_{i=1}^n \lambda_i y_i^2}{\sum_{i=1}^n y_i^2},$$

and using the fact that all the eigenvalues of A are positive, we obtain the bounds

$$R = \frac{\sum_{i=1}^n \lambda_i y_i^2}{\sum_{i=1}^n y_i^2} \leq \frac{\lambda_n \sum_{i=1}^n y_i^2}{\sum_{i=1}^n y_i^2} = \lambda_n, \quad R = \frac{\sum_{i=1}^n \lambda_i y_i^2}{\sum_{i=1}^n y_i^2} \geq \frac{\lambda_1 \sum_{i=1}^n y_i^2}{\sum_{i=1}^n y_i^2} = \lambda_1.$$

Here, λ_1 is the smallest and λ_n is the largest eigenvalue of A .

- 2 a) A is normal if $AA^H = A^H A$. Here,

$$A = \begin{bmatrix} 2 & 0 \\ -1 & 1 \end{bmatrix}, \quad A^H = \begin{bmatrix} 2 & -1 \\ 0 & 1 \end{bmatrix},$$

which leads to

$$AA^H = \begin{bmatrix} 4 & -2 \\ -2 & 2 \end{bmatrix} \quad \text{and} \quad A^H A = \begin{bmatrix} 5 & -1 \\ -1 & 0 \end{bmatrix}.$$

Hence, A is not normal.

- b)** Here, A is a lower triangular matrix, and we have the eigenvalues $\lambda_1 = 2$ and $\lambda_2 = 1$ on the diagonal. For the eigenvectors, we have that

$$Av_1 = 2v_1 \implies v_1 = \begin{bmatrix} 1 \\ -1 \end{bmatrix}, \quad \text{and} \quad Av_2 = v_2 \implies v_2 = \begin{bmatrix} 0 \\ 1 \end{bmatrix}.$$

- c)** We see that the eigenvectors are linearly independent, but they are not orthogonal since $v_1^H v_2 \neq 0$.

- d)** Yes. All the eigenvalues are nondefective, so the eigenvectors form a full set of linearly independent vectors for \mathbb{R}^2 . This implies that A can be diagonalized, i.e. we can write $A = V\Lambda V^{-1}$, where the columns of V are the eigenvectors of A and Λ is a diagonal matrix containing the eigenvalues of A .

- e)** We rewrite A as

$$A = \underbrace{\frac{1}{2}(A + A^T)}_H + \underbrace{\frac{1}{2}(A - A^T)}_S,$$

where

$$\begin{aligned} H^T &= H && \text{i.e. symmetric,} \\ S^T &= -S && \text{i.e. skew-symmetric.} \end{aligned}$$

Since H is symmetric, we can orthogonally diagonalize it, $H = Q\Lambda Q^T$. Hence, for all $u \in \mathbb{R}^2$, letting $c = Q^T u$,

$$u^T A u = u^T H u + \underbrace{u^T S u}_0 = u^T H u = c^T \Lambda c = \sum_i \lambda_i c_i^2.$$

We find that the matrix H is

$$H = \begin{bmatrix} 2 & -1/2 \\ -1/2 & 1 \end{bmatrix},$$

which has eigenvalues $\lambda = (3 \pm \sqrt{2})/2$, which are both positive. Thus, we choose $\alpha = \lambda_{\min}(H) = (3 - \sqrt{2})/2$, which makes

$$u^T A u = \sum_i \lambda_i c_i^2 \geq \alpha \|c\|_2^2 = \alpha \|u\|_2^2.$$

- f)** Yes, this follows from **e)** since $\alpha > 0$.
- g)** A Schur factorization is a product $A = QRQ^H$ for Q unitary and R upper triangular. From **d)**, we know that A can be diagonalized, $A = V\Lambda V^{-1}$, however, V is not unitary, so this is not a Schur factorization. We orthonormalize V using, say, the Gram–Schmidt process, to obtain $V = Q\tilde{R}$, where Q is unitary and \tilde{R} is upper triangular. Substituting this factorization into the diagonalization leads to a Schur decomposition

$$A = QRQ^H, \quad \text{where} \quad R = \tilde{R}\Lambda\tilde{R}^{-1}.$$

n	Time (s)	Time/ n^3
500	$1.30 \cdot 10^{-2}$	$1.04 \cdot 10^{-10}$
600	$2.29 \cdot 10^{-2}$	$1.06 \cdot 10^{-10}$
700	$3.52 \cdot 10^{-2}$	$1.03 \cdot 10^{-10}$
800	$5.27 \cdot 10^{-2}$	$1.03 \cdot 10^{-10}$
900	$7.27 \cdot 10^{-2}$	$1.00 \cdot 10^{-10}$

Table 1: Two-dimensional Poisson problem with diagonalization method

The Gram–Schmidt process yields

$$w_1 = v_1, \quad q_1 = \frac{w_1}{\|w_1\|_2} = \frac{1}{\sqrt{2}} \begin{bmatrix} 1 \\ -1 \end{bmatrix}$$

$$w_2 = v_2 - (v_2^H q_1) q_1 = \begin{bmatrix} 0 \\ 1 \end{bmatrix} - \frac{1}{2} \begin{bmatrix} -1 \\ 1 \end{bmatrix} = \frac{1}{2} \begin{bmatrix} 1 \\ 1 \end{bmatrix}, \quad q_2 = \frac{w_2}{\|w_2\|_2} = \frac{1}{\sqrt{2}} \begin{bmatrix} 1 \\ 1 \end{bmatrix}.$$

Thus

$$Q = \frac{1}{\sqrt{2}} \begin{bmatrix} 1 & 1 \\ -1 & 1 \end{bmatrix},$$

giving

$$\begin{aligned} R &= Q^H A Q \\ &= \frac{1}{2} \begin{bmatrix} 1 & -1 \\ 1 & 1 \end{bmatrix} \begin{bmatrix} 2 & 0 \\ -1 & 1 \end{bmatrix} \begin{bmatrix} 1 & 1 \\ -1 & 1 \end{bmatrix} \\ &= \begin{bmatrix} 2 & 1 \\ 0 & 1 \end{bmatrix}. \end{aligned}$$

3 We consider again the Poisson problem

$$\begin{aligned} -\Delta u &= f, & \text{in } \Omega &= (0, 1) \times (0, 1), \\ u &= 0, & \text{on } \partial\Omega. \end{aligned}$$

and discretize this with the five point formula using a uniform grid spacing, $h = 1/n$, in each spatial direction. The number of unknowns is therefore $(n-1)^2 \approx n^2$.

- a)
 - i) First we solve the Poisson problem using the diagonalization method. See Table 1.
 - ii) Next, we solve the same problem using LU factorization exploiting sparsity. See Table 2
 - iii) Finally, we solve the same problem using full LU factorization without exploiting sparsity. See Table 3.
- b) All the methods perform as expected. Note the difference in absolute solution times. The diagonalization method is clearly much better than either of the other methods.
- c) For the one-dimensional Laplace operator the continuous eigenfunctions and eigenvalues are (see the note by Rønquist)

$$\begin{aligned} u_j^*(x) &= \sin(j\pi x), \\ \lambda_j^* &= j^2\pi^2, \quad j = 1, \dots, \infty. \end{aligned}$$

n	Time (s)	Time/ n^4
50	$1.55 \cdot 10^{-2}$	$2.47 \cdot 10^{-9}$
60	$2.97 \cdot 10^{-2}$	$2.29 \cdot 10^{-9}$
70	$5.12 \cdot 10^{-2}$	$2.13 \cdot 10^{-9}$
80	$8.34 \cdot 10^{-2}$	$2.04 \cdot 10^{-9}$
90	$1.27 \cdot 10^{-1}$	$1.93 \cdot 10^{-9}$

Table 2: Two-dimensional Poisson problem with LU factorization exploiting sparsity

n	Time (s)	Time/ n^6
50	0.153	$9.78 \cdot 10^{-12}$
60	0.404	$8.67 \cdot 10^{-12}$
70	0.945	$8.03 \cdot 10^{-12}$
80	2.031	$7.75 \cdot 10^{-12}$
90	4.026	$7.58 \cdot 10^{-12}$

Table 3: Two-dimensional Poisson problem with full LU factorization

From the note, we can also see that if we discretize the Laplace operator using a central difference scheme with uniform grid spacing, $h = 1/n$, we get the matrix A^{1D} with eigenvalues

$$\lambda_j = \frac{2}{h^2} \left(1 - \cos\left(\frac{j\pi}{n}\right) \right), \quad j = 1, \dots, n-1.$$

Hence, we expect

$$\begin{aligned} \lambda_{\min} &= \frac{2}{h^2} \left(1 - \cos\left(\frac{\pi}{n}\right) \right) \approx \frac{1}{h^2} \frac{\pi^2}{n^2}, \\ \lambda_{\max} &= \frac{2}{h^2} \left(1 - \cos\left(\frac{(n-1)\pi}{n}\right) \right) \approx \frac{4}{h^2}, \\ &\Downarrow \\ \kappa(A^{1D}) &= \frac{\lambda_{\max}}{\lambda_{\min}} \approx \frac{4}{\pi^2} n^2. \end{aligned}$$

The condition number should scale as $O(n^2)$ in 1D.

d) Consider now the two-dimensional Poisson problem

$$\begin{aligned} -\Delta u &= f, & \text{in } \Omega &= (0, 1) \times (0, 1), \\ u &= 0, & \text{on } \partial\Omega. \end{aligned}$$

In the continuous case, the eigenfunctions and eigenvalues of the Laplace operator are given as

$$\begin{aligned} u_{j,k}^*(x, y) &= \sin(j\pi x) \sin(k\pi y), \\ \lambda_{j,k}^* &= j^2\pi^2 + k^2\pi^2 = (j^2 + k^2)\pi^2, \\ j &= 1, \dots, \infty, \quad k = 1, \dots, \infty. \end{aligned}$$

These results are obtained simply by using separation of variables.

n	$\lambda_{\min}(\hat{A}^{1D})$	$\lambda_{\min}(\hat{A}^{1D})$	$\kappa(A^{1D})$
10	$8.10 \cdot 10^{-2}$	3.92	48.37
20	$2.23 \cdot 10^{-2}$	3.98	178.06
30	$1.03 \cdot 10^{-2}$	3.99	388.81
40	$5.87 \cdot 10^{-3}$	3.99	680.62

(a) One dimension

n	$\lambda_{\min}(\hat{A}^{2D})$	$\lambda_{\min}(\hat{A}^{2D})$	$\kappa(A^{2D})$
10	$1.62 \cdot 10^{-1}$	7.84	48.37
20	$4.47 \cdot 10^{-2}$	7.96	178.06
30	$2.05 \cdot 10^{-2}$	7.98	388.81
40	$1.17 \cdot 10^{-2}$	7.99	680.62

(b) Two dimensions

Table 4: Condition numbers

Again, we discretize the Poisson problem using the five-point formula, and get the discrete Laplace operator (or matrix), A^{2D} . The eigenvalues of A^{2D} are

$$\lambda_{j,k} = \frac{2}{h^2} \left(1 - \cos\left(\frac{j\pi}{n}\right) \right) + \frac{2}{h^2} \left(1 - \cos\left(\frac{k\pi}{n}\right) \right),$$

so we expect that

$$\lambda_{\min} = \frac{2}{h^2} \left(1 - \cos\left(\frac{\pi}{n}\right) \right) + \frac{2}{h^2} \left(1 - \cos\left(\frac{\pi}{n}\right) \right) \approx \frac{2}{h^2} \frac{\pi^2}{n^2},$$

$$\lambda_{\min} = \lambda_{j,k} = \frac{2}{h^2} \left(1 - \cos\left(\frac{(n-1)\pi}{n}\right) \right) + \frac{2}{h^2} \left(1 - \cos\left(\frac{(n-1)\pi}{n}\right) \right) \approx \frac{8}{h^2},$$

↓

$$\kappa(A^{2D}) = \frac{\lambda_{\max}}{\lambda_{\min}} \approx \frac{4}{\pi^2} n^2.$$

We compute the maximum and minimum eigenvalues for the matrices $\hat{A}^{1D} = h^2 A^{1D}$ and $\hat{A}^{2D} = h^2 A^{2D}$ by using the command `eig` in MATLAB. The condition number is found by the use of `cond`. See Table 4.

We observe that the condition number κ is the same for one and two dimensions. We also get the expected behavior for the minimum and maximum eigenvalues.