

(1) Show that the Hilbert matrix is positive definite.

*Proof:*

Suppose we have the  $n \times n$  Hilbert matrix

$$H_{ij} = \frac{1}{i+j-1}, \quad i, j = 1, \dots, n.$$

Then, to show that  $H$  is positive definite, we need to show that  $x^* H x > 0$  for all  $x \neq 0$  in  $\mathbb{C}^n$ . So, suppose we have any nonzero  $x \in \mathbb{C}^n$ . Then,

$$\begin{aligned} x^* H x &= \sum_{i=1}^n \left( \bar{x}_i \sum_{j=1}^n x_j H_{ij} \right) = \sum_{i=1}^n \sum_{j=1}^n \bar{x}_i x_j \frac{1}{i+j-1} \\ &= \sum_{i=1}^n \sum_{j=1}^n \bar{x}_i x_j \int_0^1 t^{i+j-2} dt \\ &= \sum_{i=1}^n \sum_{j=1}^n \int_0^1 (\bar{x}_i t^{i-1}) (x_j t^{j-1}) dt \\ &= \int_0^1 \sum_{i=1}^n \sum_{j=1}^n (\bar{x}_i t^{i-1}) (x_j t^{j-1}) dt \\ &= \int_0^1 \left( \sum_{i=1}^n \bar{x}_i t^{i-1} \right) \left( \sum_{j=1}^n x_j t^{j-1} \right) dt. \end{aligned}$$

Now if we let  $\alpha = \sum_{j=1}^n x_j t^{j-1}$ , our equation becomes

$$\begin{aligned} x^* H x &= \int_0^1 \left( \sum_{i=1}^n \bar{x}_i t^{i-1} \right) \left( \sum_{j=1}^n x_j t^{j-1} \right) dt \\ &= \int_0^1 \bar{\alpha} \alpha dt \\ &= \int_0^1 |\alpha|^2 dt \\ &> 0. \end{aligned}$$

So, our Hilbert matrix must be positive definite. □

- (2) Using my power iteration code (attached in section: Code Used), I generated the table below of the largest eigenvalues for the first 10 square Hilbert matrices.

$n$	Eigenvalue	Eigenvector
1	1.00	$[1.0]^T$
2	1.27	$[0.88, 0.47]^T$
3	1.41	$[0.83, 0.46, 0.32]^T$
4	1.50	$[0.79, 0.45, 0.32, 0.25]^T$
5	1.57	$[0.77, 0.45, 0.32, 0.25, 0.21]^T$
6	1.62	$[0.75, 0.44, 0.32, 0.25, 0.21, 0.18]$
7	1.66	$[0.73, 0.44, 0.32, 0.25, 0.21, 0.18, 0.16]^T$
8	1.70	$[0.72, 0.43, 0.32, 0.26, 0.21, 0.18, 0.16, 0.15]^T$
9	1.73	$[0.71, 0.43, 0.32, 0.26, 0.21, 0.19, 0.16, 0.15, 0.13]^T$
10	1.75	$[0.70, 0.43, 0.32, 0.26, 0.22, 0.19, 0.16, 0.15, 0.13, 0.12]^T$

- (3) To find the smallest eigenvalues using a power iteration, I just change the matrix-vector multiplication to a backslash as to solve  $Ax = b$  instead of computing  $x = Ab$ . This change has the effect of making the smallest eigenvalue of  $A$  dominate the power iteration. For  $n = 16$ , I obtained the eigenvalue:

$$\lambda_{\min} = -2.765686556840889 \cdot 10^{-18}$$

which is not even positive even though we know Hilbert matrices are symmetric positive definite. So my naive Julia code is definitely not converging properly under double precision. Using Mathematica, I found that the actual smallest eigenvalue was

$$\lambda_{\min} = 9.1974198206514524197318 \cdot 10^{-23}$$

- (4) Assume that a real matrix  $A$  has eigenvalues  $\lambda_1 = -\lambda_2$  and  $|\lambda_1| = |\lambda_2| > |\lambda_3| \geq |\lambda_n|$ .

Without loss of generality, assume  $\lambda_1 > 0$  which makes  $\lambda_2 < 0$ . Now, note that because  $A$ , eigenvectors corresponding to different eigenvalues of  $A$  are orthogonal. Then, to find the eigenvectors corresponding to  $\lambda_1$  and  $\lambda_2$ , run the standard power iteration (until convergence) to find some normalized vector,  $\vec{v}_0$ , in the span of  $\vec{\lambda}_1$  and  $\vec{\lambda}_2$  where  $\vec{\lambda}_1$  and  $\vec{\lambda}_2$  are eigenvectors corresponding to  $\lambda_1$  and  $\lambda_2$  respectively. Then, because  $\vec{\lambda}_1$  and  $\vec{\lambda}_2$  are orthogonal, we can uniquely decompose  $\vec{v}_0$  as

$$\vec{v}_0 = a\vec{\lambda}_1 + b\vec{\lambda}_2$$

for some constants  $a$  and  $b$ . Then, one more power iteration on  $\vec{v}_0$  will yield

$$\vec{v}_1 = a\vec{\lambda}_1 - b\vec{\lambda}_2$$

because  $\lambda_2 < 0$  and the eigenvectors are orthogonal. Then, we can simply find an eigenvector corresponding to  $\lambda_1$  as

$$\vec{v}_0 + \vec{v}_1 = 2a\vec{\lambda}_1$$

and an eigenvector corresponding to  $\lambda_2$  as

$$\vec{v}_0 - \vec{v}_1 = 2b\vec{\lambda}_2.$$

- (5) A real symmetric matrix  $A$  has an eigenvalue 1 of multiplicity 8; the rest of the eigenvalues are  $\leq 0.1$  in absolute value.

We can find an orthogonal basis for the 8-dimensional eigenspace corresponding to the dominant eigenvalue by combining the standard power iteration with the Gram-Schmidt process. To do so:

- (i) Generate 8, random vectors.
- (ii) Apply power iteration to each vector until convergence. The resulting vectors will be in the span of the dominant eigenvectors.
- (iii) Apply Gram-Schmidt to the set of generated vectors.
- (iv) You're all done!

To estimate how long it will take to find this basis for an  $n \times n$  matrix, we need to figure out how long it will take each power iteration to converge to double-precision. If each subdominant eigenvalue has a magnitude less than or equal to 0.1, we should expect our iteration to converge in

$$(n - 8)(0.1)^i \leq 10^{-16}$$

or

$$i = \log_{0.1} \frac{10^{-16}}{n - 8}$$

iterations. This calculation comes from the fact that we need  $n - 8$  vectors with  $|\lambda| < 0.1$  to vanish from power iteration to have convergence to double-precision. Then, we need to do 8 power iterations for a total of

$$8i = 8 \log_{0.1} \frac{10^{-16}}{n - 8}$$

iterations minimum for convergence. For  $n \leq 1000$ ,  $i \leq 19$  so for any matrix smaller than  $1000 \times 1000$ , we should expect convergence to our basis in a minimum of 152 iterations.

## Code Used