

## Unit 4. Eigenvalues

Numerical Analysis

EE/NTHU

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### Eigenvalues and Eigenvectors

#### Definition 4.1.1.

Given a real  $n \times n$  matrix  $\mathbf{A}$ , the number  $\lambda \in \mathbb{C}$  is an **eigenvalue** of  $\mathbf{A}$  if there is a nonnull vector  $\mathbf{x} \in \mathbb{R}$  such that

$$\mathbf{A}\mathbf{x} = \lambda\mathbf{x}. \quad (4.1.1)$$

The vector  $\mathbf{x}$  is the **eigenvector** associated with the eigenvalue  $\lambda$  and the set of eigenvalues of  $\mathbf{A}$  is the **spectrum** of  $\mathbf{A}$ , denoted by  $\sigma(\mathbf{A})$ .

#### Theorem 4.1.2.

The eigenvalue  $\lambda$  corresponding to the eigenvector  $\mathbf{x}$  can be determined by computing the **Rayleigh quotient**

$$\lambda = \mathbf{x}^T \mathbf{A} \mathbf{x} / \mathbf{x}^T \mathbf{x}. \quad (4.1.2)$$

#### Definition 4.1.3.

The **characteristic polynomial** of the real  $n \times n$  matrix  $\mathbf{A}$  is

$$p_A(\lambda) = \det(\mathbf{A} - \lambda \mathbf{I}). \quad (4.1.3)$$

## Eigenvalues and Eigenvectors, II

### Theorem 4.1.4.

The eigenvalue  $\lambda$  is the solution of the **characteristic equation**

$$p_A(\lambda) = 0. \quad (4.1.4)$$

- Since the characteristic polynomial is of degree  $n$  with respect to  $\lambda$ , there exist  $n$  eigenvalues of  $\mathbf{A}$ . But these eigenvalues may not be distinct from each other.

### Theorem 4.1.5.

Given real  $n \times n$  matrix  $\mathbf{A}$  with the eigenvalues  $\lambda_i$ ,  $i = 1, \dots, n$  then

$$\det(\mathbf{A}) = \prod_{i=1}^n \lambda_i, \quad (4.1.5)$$

$$\text{tr}(\mathbf{A}) = \sum_{i=1}^n \lambda_i. \quad (4.1.6)$$

## Eigenvalues and Eigenvectors, III

### Theorem 4.1.6.

Given a real  $n \times n$  matrix  $\mathbf{A}$  with the eigenvalues  $\lambda_i$ ,  $i = 1, \dots, n$ , then

1.  $\mathbf{A}$  is singular if and only if there is a  $\lambda_i = 0$ ,  $1 \leq i \leq n$ .
2. Complex eigenvalues of  $\mathbf{A}$  occur in conjugate pairs.
3. The matrix polynomial  $p_A(\mathbf{A})$  satisfies

$$p_A(\mathbf{A}) = \mathbf{0}. \quad (4.1.7)$$

### Theorem 4.1.7.

The **spectral radius** of a real matrix  $\mathbf{A}$  is defined as

$$\rho(\mathbf{A}) = \max_{\lambda \in \sigma(\mathbf{A})} |\lambda|. \quad (4.1.8)$$

# Eigenvalues and Eigenvectors, IV

## Theorem 4.1.8.

If the real matrix  $\mathbf{A}$  has the following block triangular form

$$\mathbf{A} = \begin{bmatrix} \mathbf{A}_{11} & \mathbf{A}_{12} & \cdots & \mathbf{A}_{1k} \\ 0 & \mathbf{A}_{22} & \cdots & \mathbf{A}_{2k} \\ \vdots & \vdots & \ddots & \vdots \\ 0 & \cdots & 0 & \mathbf{A}_{kk} \end{bmatrix} \quad (4.1.9)$$

Then

$$p_A(\lambda) = p_{A_{11}}(\lambda) \cdot p_{A_{22}}(\lambda) \cdots p_{A_{kk}}(\lambda), \quad (4.1.10)$$

$$\sigma(\mathbf{A}) = \bigcup_{j=1}^k \sigma(\mathbf{A}_{jj}). \quad (4.1.11)$$

## Theorem 4.1.9.

If the real matrix  $\mathbf{A}$  is triangular then

$$\sigma(\mathbf{A}) = \{a_{ii} | i = 1, \dots, n\}.$$

# Eigenvalues and Eigenvectors, V

## Definition 4.1.10.

Given a real  $n \times n$  matrix  $\mathbf{A}$  with the characteristic polynomial

$$p_A(\lambda) = \prod_{i=1}^k (\lambda - \lambda_i)^{\delta_i} \quad (4.1.12)$$

with  $\lambda_i \neq \lambda_j$ ,  $1 \leq i, j \leq k$ , and  $\sum_{i=1}^k \delta_i = n$ . The **algebraic multiplicity** of  $\lambda_i$  is  $\delta_i$ ,  $1 \leq i \leq k$ .

## Theorem 4.1.11.

The set of eigenvectors associated with a eigenvalue,  $\lambda$ , of a real  $n \times n$  matrix  $\mathbf{A}$  forms a subspace. The dimension of this subspace is called **geometric multiplicity** of the eigenvalue  $\lambda$ . For any  $\lambda$  the geometric multiplicity is less than or equal to the algebraic multiplicity.

- The power method can be used to estimate the largest and the smallest eigenvalues.
- These eigenvalues are needed for the condition number of a linear system solution, and other applications.
- In this section, we assume the real  $n \times n$  matrix  $\mathbf{A}$  has  $n$  distinct real eigenvalues and  $\mathbf{x}_i$  is the eigenvector associated with eigenvalue  $\lambda_i$ ,  $i = 1, \dots, n$ .
- It is further assumed that the eigenvalues are ordered as

$$|\lambda_1| > |\lambda_2| \geq |\lambda_3| \cdots \geq |\lambda_n|, \quad (4.1.13)$$

where  $\lambda_1$  has the algebraic multiplicity of 1.

- Under these assumptions,  $\lambda_1$  is called the **dominant** eigenvalue of matrix  $\mathbf{A}$ .

## The Power Method, II

### Algorithm 4.1.12. Power Method

Given a diagnosable matrix  $\mathbf{A}$  and an arbitrary initial vector  $\mathbf{q}^{(0)}$ , then

$$\mathbf{q}^{(k)} = \frac{\mathbf{A}^k \mathbf{q}^{(0)}}{\|\mathbf{A}^k \mathbf{q}^{(0)}\|_2}, \quad (4.1.14)$$

$$\nu^{(k)} = (\mathbf{q}^{(k)})^T \mathbf{A} \mathbf{q}^{(k)}. \quad (4.1.15)$$

- Since  $\mathbf{A}$  is diagnosable, the eigenvectors form a basis of  $\mathbb{R}^n$ . Thus,  $\mathbf{q}^{(0)}$  can be expressed as

$$\mathbf{q}^{(0)} = \sum_{i=1}^n \alpha_i \mathbf{x}_i = \alpha_1 \mathbf{x}_1 + \alpha_2 \mathbf{x}_2 + \cdots + \alpha_n \mathbf{x}_n$$

Then

$$\mathbf{A} \mathbf{q}^{(0)} = \mathbf{A} \sum_{i=1}^n \alpha_i \mathbf{x}_i = \sum_{i=1}^n \alpha_i \lambda_i \mathbf{x}_i$$

# The Power Method, III

- And

$$\begin{aligned}\mathbf{A}^k \mathbf{q}^{(0)} &= \mathbf{A}^k \sum_{i=1}^n \alpha_i \mathbf{x}_i = \sum_{i=1}^n \alpha_i \mathbf{A}^k \mathbf{x}_i = \sum_{i=1}^n \alpha_i \lambda_i^k \mathbf{x}_i \\ &= \alpha_1 \lambda_1^k \left( \mathbf{x}_1 + \sum_{i=2}^n \frac{\alpha_i}{\alpha_1} \left( \frac{\lambda_i}{\lambda_1} \right)^k \mathbf{x}_i \right)\end{aligned}$$

When  $k \rightarrow \infty$ ,

$$\mathbf{A}^k \mathbf{q}^{(0)} \rightarrow \alpha_1 \lambda_1^k \mathbf{x}_1, \text{ if } \alpha_1 \neq 0.$$

Or, we can write

$$\mathbf{q}^{(k)} = \frac{\alpha_1 \lambda_1^k (\mathbf{x}_1 + \mathbf{y}^{(k)})}{\|\alpha_1 \lambda_1^k (\mathbf{x}_1 + \mathbf{y}^{(k)})\|_2}$$

where  $\mathbf{y}^{(k)} = \sum_{i=2}^n \frac{\alpha_i}{\alpha_1} \left( \frac{\lambda_i}{\lambda_1} \right)^k \mathbf{x}_i$ , and  $\mathbf{y}^{(k)} \rightarrow 0$  when  $k \rightarrow \infty$ .

Also, if  $\alpha_1 \neq 0$ , as  $k \rightarrow \infty$

$$\mathbf{q}^{(k)} \rightarrow \mathbf{x}_1 \quad (4.1.16)$$

$$\nu^{(k)} \rightarrow \lambda_1 \quad (4.1.17)$$

# The Power Method, IV

## Theorem 4.1.13.

Let  $\mathbf{A} \in \mathbb{C}^{n \times n}$  be a diagonizable matrix whose eigenvalues satisfy equation (4.1.13). Assuming  $\alpha_1 \neq 0$ , there is a constant  $C > 0$  such that

$$\|\tilde{\mathbf{q}}^{(k)} - \mathbf{x}_1\|_2 \leq C \left| \frac{\lambda_2}{\lambda_1} \right|^k, \quad k \geq 1, \quad (4.1.18)$$

where

$$\tilde{\mathbf{q}}^{(k)} = \frac{\mathbf{q}^{(k)} \|\mathbf{A}^k \mathbf{q}^{(0)}\|_2}{\alpha_1 \lambda_1^k} = \mathbf{x}_1 + \sum_{i=2}^n \frac{\alpha_i}{\alpha_1} \left( \frac{\lambda_i}{\lambda_1} \right)^k \mathbf{x}_i, \quad k = 1, 2, \dots \quad (4.1.19)$$

$$\begin{aligned}\|\tilde{\mathbf{q}}^{(k)} - \mathbf{x}_1\|_2 &= \left\| \sum_{i=2}^n \frac{\alpha_i}{\alpha_1} \left( \frac{\lambda_i}{\lambda_1} \right)^k \mathbf{x}_i \right\|_2 \leq \left( \sum_{i=2}^n \left[ \frac{\alpha_i}{\alpha_1} \left( \frac{\lambda_i}{\lambda_1} \right)^k \right]^2 \right)^{1/2} \\ &\leq \left( \sum_{i=2}^n \left[ \frac{\alpha_i}{\alpha_1} \left( \frac{\lambda_2}{\lambda_1} \right)^k \right]^2 \right)^{1/2} = \left| \frac{\lambda_2}{\lambda_1} \right|^k \left[ \sum_{i=2}^n \left( \frac{\alpha_i}{\alpha_1} \right)^2 \right]^{1/2}\end{aligned}$$

Thus,  $C = \left[ \sum_{i=2}^n \left( \frac{\alpha_i}{\alpha_1} \right)^2 \right]^{1/2}$ , and  $C$  is independent of  $k$ .

- The preceding theorem states that the power method converges with the rate  $\left| \frac{\lambda_2}{\lambda_1} \right|$ .
  - If  $|\lambda_2| \ll |\lambda_1|$  then it converges quickly,
  - On the other hand, if  $|\lambda_2| \approx |\lambda_1|$  then it converges slowly.
- $\tilde{\mathbf{q}}^{(k)}$  converges to  $\mathbf{x}_1$
- Since  $\tilde{\mathbf{q}}^{(k)} = \frac{\mathbf{q}^{(k)} \|\mathbf{A}^k \mathbf{q}^{(0)}\|_2}{\alpha_1 \lambda_1^k}$ ,  $\mathbf{q}^{(k)}$  converges, too.
  - See textbook.
- And  $\nu^{(k)} = (\mathbf{q}^{(k)})^T \mathbf{A} \mathbf{q}^{(k)}$  converges to  $\lambda_1$ .
  - By the rate  $\left| \frac{\lambda_2}{\lambda_1} \right|$ .
- If  $\mathbf{A}$  is real and symmetric and  $\alpha_1 \neq 0$ , then it can be shown that

$$|\lambda_1 - \nu^{(k)}| \leq |\lambda_1 - \lambda_n| \cdot \tan^2(\theta_0) \cdot \left| \frac{\lambda_2}{\lambda_1} \right|^{2k}. \quad (4.1.20)$$

where  $\cos(\theta_0) = |\mathbf{x}_1^T \mathbf{q}^{(0)}| \neq 0$ .

- In this case, the convergence rate is quadratic,  $\left| \frac{\lambda_2}{\lambda_1} \right|^2$ .

## Stopping Criteria

- So far, we know that  $\lim_{k \rightarrow \infty} \mathbf{q}^{(k)} = \mathbf{x}_1$ , the eigenvector associated with  $\lambda_1$  of matrix  $\mathbf{A}$ , and  $\lim_{k \rightarrow \infty} \nu^{(k)} = \lambda_1$ , which is the eigenvalue with the largest module.
- Define the residue at iteration  $k$  as

$$\mathbf{r}^{(k)} = \mathbf{A} \mathbf{q}^{(k)} - \nu^{(k)} \mathbf{q}^{(k)}, \quad k \geq 1. \quad (4.1.21)$$

Then, as  $k \rightarrow \infty$ ,  $\mathbf{r}^{(k)} \rightarrow \mathbf{0}$ . One can use  $\|\mathbf{r}^{(k)}\|_2$  as a stopping criterion.

- In fact, it has been shown that

$$|\lambda_1 - \nu^{(k)}| \simeq \frac{\|\mathbf{r}^{(k)}\|_2}{|(\mathbf{w}^{(k)})^T \mathbf{q}^{(k)}|}, \quad k \geq 1, \quad (4.1.22)$$

where  $\mathbf{w}^{(k)}$  satisfies  $(\mathbf{w}^{(k)})^T \mathbf{A} = \nu^{(k)} (\mathbf{w}^{(k)})^T$ , and as  $k \rightarrow \infty$ ,  $\mathbf{w}^{(k)} \rightarrow \mathbf{w}$  and  $\mathbf{w}^T \mathbf{A} = \lambda_1 \mathbf{w}^T$  is the left eigenvector associated with  $\lambda_1$ .

If  $\mathbf{A}$  is symmetric the  $\mathbf{w} = \mathbf{q}$ .

- One approach is to use Eq. (4.1.22) as the stopping criterion.



## Algorithm 4.1.14. The Power Method

Given a diagonizable matrix  $\mathbf{A}$ , an initial vector  $\mathbf{q}^{(0)}$ , a small number  $\epsilon$  and a large integer *maxiter*, let

$$tol = 1 + \epsilon, \mathbf{q}^{(0)} = \frac{\mathbf{q}^{(0)}}{\|\mathbf{q}^{(0)}\|_2}, \text{ and } k = 0,$$

while ( $tol \geq \epsilon$  and  $k \leq \text{maxiter}$ ) {

$$\mathbf{z} = \mathbf{A}\mathbf{q}^{(k)},$$

$$k = k + 1,$$

$$\mathbf{q}^{(k)} = \frac{\mathbf{z}}{\|\mathbf{z}\|_2},$$

$$\nu^{(k)} = (\mathbf{q}^{(k)})^T \mathbf{A} \mathbf{q}^{(k)},$$

$$\mathbf{r}^{(k)} = \mathbf{A} \mathbf{q}^{(k)} - \nu^{(k)} \mathbf{q}^{(k)},$$

$$(\mathbf{u}^{(k)})^T = (\mathbf{q}^{(k)})^T \mathbf{A},$$

$$\mathbf{w}^{(k)} = \frac{\mathbf{u}}{\|\mathbf{u}\|_2},$$

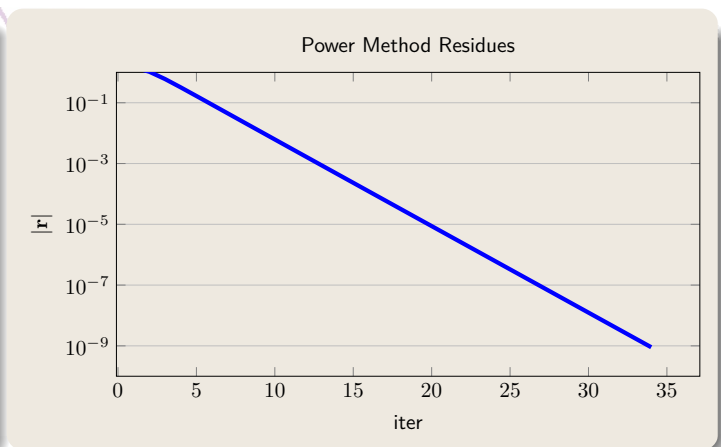
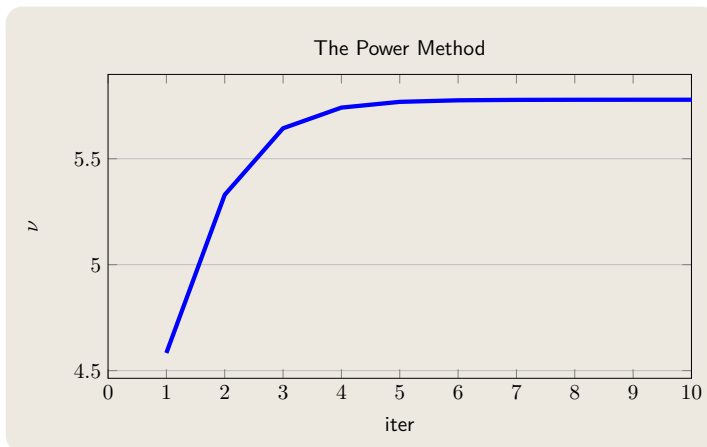
$$tol = \frac{\|\mathbf{r}^{(k)}\|_2}{|(\mathbf{w}^{(k)})^T \mathbf{q}^{(k)}|},$$

}

## The Algorithm, II

- Note that it can also check for  $|(\mathbf{w}^{(k)})^T \mathbf{q}^{(k)}|$ . If this number is 0, then  $\lambda_1$  does not have the algebraic multiplicity of 1.
- If the algebraic multiplicity of  $\lambda_1$  is greater than 1, then  $\mathbf{q}^{(k)}$  may not converge though  $\nu^{(k)}$  is convergent.
  - $\mathbf{r}^{(k)}$  may not be convergent either.
  - Thus, using  $\|\mathbf{r}^{(k)}\|_2$  as the stopping criterion may not work.
- If  $\mathbf{A}$  is symmetric then the left eigenvector  $\mathbf{q}_i$  of a eigenvalue  $\lambda_i$  is always the same as the right eigenvector  $\mathbf{w}_i$  associated with  $\lambda_i$ , and  $|(\mathbf{w}^{(k)})^T \mathbf{q}^{(k)}| = 1$ .
- For asymmetric  $\mathbf{A}$ , this property does not hold but  $|(\mathbf{w}^{(k)})^T \mathbf{q}^{(k)}|$  is convergent to a single number as both eigenvectors converges.

# The Power Method, Example



- Using the resistor network example of Unit 3
- The power method is shown to be convergent with a constant rate
  - $\lambda_1 = 5.77846$ .

## The Power Method, Complexity

- The first power method algorithm, Algorithm (4.1.12), is usually formulated as

$$\mathbf{q}^{(k+1)} = \frac{\mathbf{A}\mathbf{q}^{(k)}}{\|\mathbf{A}\mathbf{q}^{(k)}\|_2}, \quad (4.1.23)$$

$$\nu^{(k+1)} = (\mathbf{q}^{(k+1)})^T \mathbf{A} \mathbf{q}^{(k+1)}. \quad (4.1.24)$$

- The computation is dominated by  $\mathbf{A}\mathbf{q}^{(k)}$ .
  - Matrix-vector multiplication.
  - Computational complexity is  $\mathcal{O}(n^2)$  per iteration.
  - Overall computational complexity is  $\mathcal{O}(N_{iter} \times n^2)$ .
    - $N_{iter}$  is the number of iterations needed to reach to a converged solution.
    - A function of  $\left| \frac{\lambda_2}{\lambda_1} \right|$ .
- The computation of the second form of power method, Algorithm (4.1.14), is also dominated by  $\mathbf{A}\mathbf{q}^{(k)}$  and  $(\mathbf{q}^{(k)})^T \mathbf{A}$ .
  - Matrix-vector multiplications.
  - Overall computational complexity remains  $\mathcal{O}(N_{iter} \times n^2)$ .
  - But with a larger coefficient.



# Inverse Power Method

- The power method can be modified to find the eigenvalue with the smallest module, and the eigenvector associated.

## Algorithm 4.1.15. Inverse Power Method

Given a diagonable matrix  $\mathbf{A}$  and an arbitrary initial guess  $\mathbf{q}^{(0)}$  with unit Euclidean norm, iterate for  $k = 1, \dots$

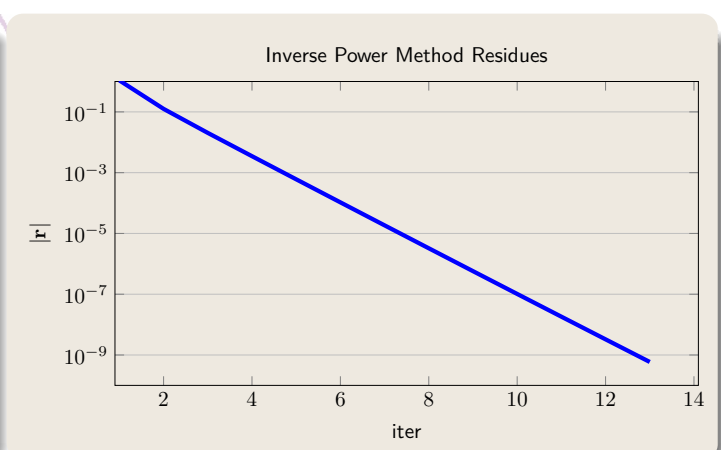
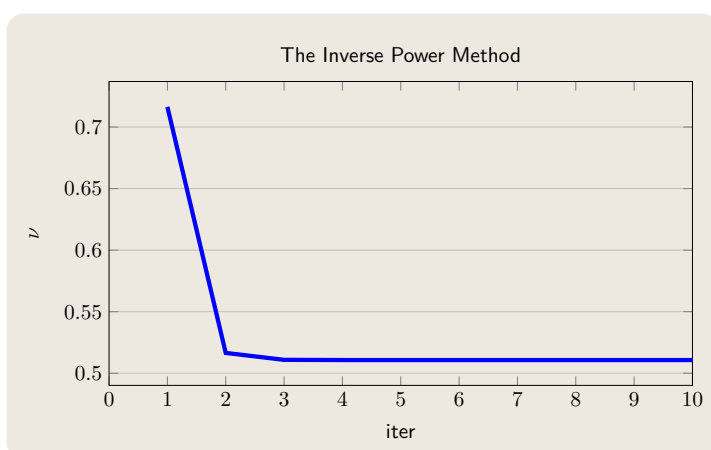
$$\mathbf{A}\mathbf{z}^{(k)} = \mathbf{q}^{(k-1)}, \quad (4.1.25)$$

$$\mathbf{q}^{(k)} = \frac{\mathbf{z}^{(k)}}{\|\mathbf{z}^{(k)}\|_2}, \quad (4.1.26)$$

$$\mu^{(k)} = (\mathbf{q}^{(k)})^T \mathbf{A} \mathbf{q}^{(k)}. \quad (4.1.27)$$

- Note that the vector  $\mathbf{z}^{(k)}$  can be found using LU decomposition or any linear system solution method.
- Other than that this is the power method and its convergence rate is determined by  $\left| \frac{\lambda_n}{\lambda_{n-1}} \right|$ .

## Inverse Iteration Example



- Using the resistor network example of Unit 3
- The power method is shown to be convergent with a constant rate
  - $\lambda_7 = 0.510711$ .

# The Inverse Power Method with Shifting

- The inverse power method can be generalized to find eigenvalue that is closest to a specific number,  $\omega$ , and the eigenvector associated.

## Algorithm 4.1.16. Inverse Power Method with Shifting

Given a diagonosable matrix  $\mathbf{A}$  and an arbitrary initial guess  $\mathbf{q}^{(0)}$  with unit Euclidean norm and  $\omega \in \mathbb{R}$ , iterate for  $k = 1, \dots$

$$(\mathbf{A} - \omega \mathbf{I})\mathbf{z}^{(k)} = \mathbf{q}^{(k-1)}, \quad (4.1.28)$$

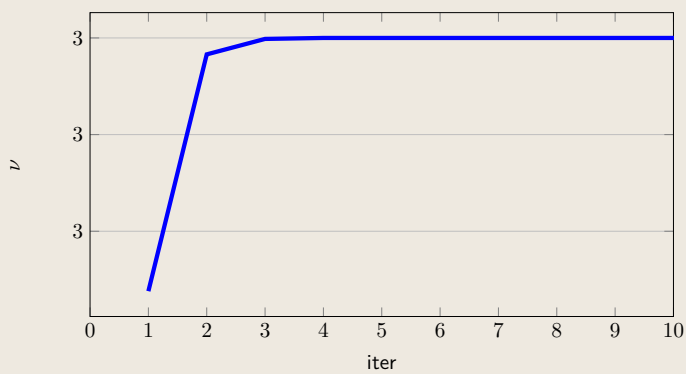
$$\mathbf{q}^{(k)} = \frac{\mathbf{z}^{(k)}}{\|\mathbf{z}^{(k)}\|_2}, \quad (4.1.29)$$

$$\mu^{(k)} = (\mathbf{q}^{(k)})^T \mathbf{A} \mathbf{q}^{(k)}. \quad (4.1.30)$$

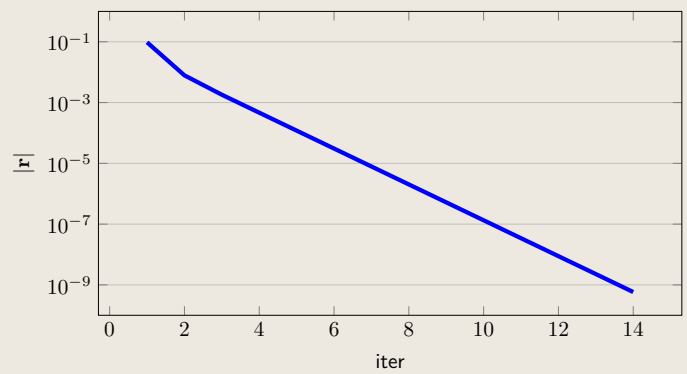
- Except Eq. (4.1.28), this algorithm is identical to the inverse power method.
- And the convergence rate is determined by the ratio of the two eigenvalues that are closest to  $\omega$ .

## Inverse Power Method with Shifting Examples

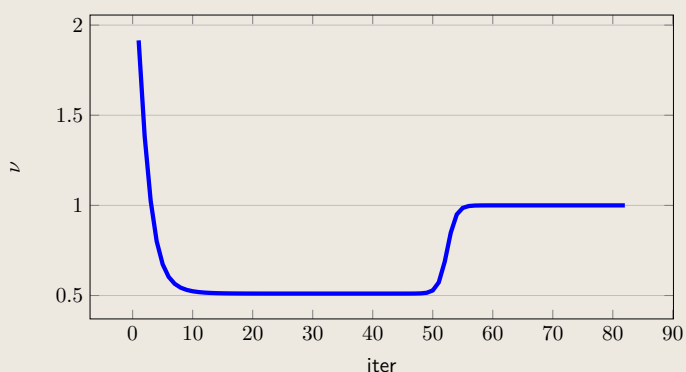
The Inverse Power Method with Shifting,  $\omega = 3.1$



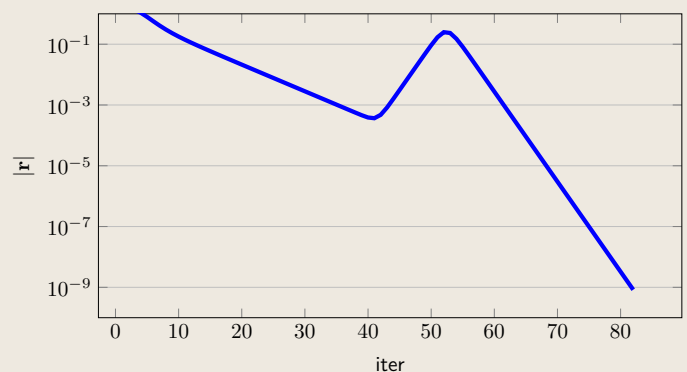
Inverse Power Method with Shifting ( $\omega = 3.1$ ), Residues



The Inverse Power Method with Shifting,  $\omega = 1.5$



Inverse Power Method with Shifting ( $\omega = 1.5$ ), Residues



- In our development we have assumed that  $\lambda_1 > \lambda_2$ , or the algebraic multiplicity of  $\lambda_1$  is 1. But, the power method is convergent when  $\lambda_1 = \lambda_2$ , since the vectors generated by Eq. (4.1.14) converge to the subspace spanned by  $\mathbf{x}_1$  and  $\mathbf{x}_2$ , and  $\nu$  converges to  $\lambda_1$ . (The original form of power method, Algorithm (4.1.12).)
- When  $\lambda_1 = -\lambda_2$  then the power method would oscillate and not converge.
- If  $\lambda_1 = \overline{\lambda_2}$ , it would also oscillate and not converge.
- The inverse power method with shifting is very effective in find eigenvalues and eigenvectors for diagnosable matrices.
- With  $\omega$  is properly positioned, the eigenvalue closest to  $\omega$  can be quickly found. The convergence rate can be improved greatly if  $\omega$  is close to  $\lambda_i$ .
- Initial guess  $\mathbf{q}^{(0)}$  can affect the convergence rate as shown in the last example. But, due to the computer round off, the eigenvalue closest to  $\omega$  is always found, even  $\mathbf{q}^{(0)}$  is an eigenvector of a different eigenvalue.
- The inverse power method with shifting is more effective even though it needs more operations for each iteration: matrix addition and forward and backward substitutions. (note that LU decomposition needs to be done only once.)

## Summary

- The power method.
  - Simple form.
  - More elaborated form.
- The inverse power method.
- Power method with shifting.