

# Software Assignment

EE24BTECH11013

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## 1. Introduction to Eigenvalues

In linear algebra, the eigenvalues of a square matrix are a set of scalars that provide critical insights into the matrix's properties. If  $A$  is an  $n \times n$  matrix, a scalar  $\lambda$  is called an eigenvalue of  $A$  if there exists a non-zero vector  $\mathbf{v}$  (called an eigenvector) such that:

$$A\mathbf{v} = \lambda\mathbf{v}.$$

Here,  $\lambda$  represents how the matrix transformation scales the eigenvector  $\mathbf{v}$ .

## 2. Why are Eigenvalues Useful?

Eigenvalues have wide-ranging applications in mathematics, science, and engineering. Their utility arises from the ability to reveal structural and dynamical properties of systems modeled by matrices.

- **Principal Component Analysis (PCA):** Eigenvalues are essential in PCA, a technique used in machine learning and data science to reduce dimensionality while preserving as much information as possible.
- **Mechanical Systems:** In structural engineering, eigenvalues describe the natural frequencies of vibrating systems.
- **Quantum Mechanics:** Eigenvalues correspond to observable quantities such as energy levels of systems.
- **Stability Analysis:** In control theory and differential equations, eigenvalues determine the stability of dynamic systems.
- **Graph Theory:** The eigenvalues of a graph's adjacency or Laplacian matrix provide information about its structure, such as connectivity and clustering.

### 3. Algorithms for Finding Eigenvalues

There are various algorithms to compute eigenvalues, each suited to different types of matrices and accuracy requirements. The most commonly used methods are:

#### 3.1. Characteristic Polynomial Method

This method solves the eigenvalue equation:

$$\det(\lambda I - A) = 0, \quad (1)$$

where  $I$  is the identity matrix. This approach is suitable for small matrices but computationally expensive for larger matrices due to the complexity of determinant calculation.

#### 3.2. Power Iteration

This iterative method finds the largest eigenvalue (in magnitude) of a matrix. It is simple and efficient for sparse or large matrices but cannot find all eigenvalues.

**Steps:**

1. Start with a random vector  $\mathbf{x}_0$ .
2. Compute  $\mathbf{x}_{k+1} = A\mathbf{x}_k$  and normalize it.
3. Repeat until convergence.

#### 3.3. QR Algorithm

The QR Algorithm is an iterative method to find all eigenvalues of a matrix. It repeatedly performs QR decomposition ( $A = QR$ ) and updates  $A$  as  $RQ$  until it converges to an upper triangular matrix, where the eigenvalues are the diagonal elements.

#### 3.4. Jacobi Method

The Jacobi method is specifically used for symmetric matrices. It applies a sequence of orthogonal transformations to make the matrix diagonal, where the diagonal entries are the eigenvalues.

#### 3.5. Divide-and-Conquer Method

This approach is efficient for dense symmetric matrices. It recursively partitions the matrix into smaller submatrices, solves for their eigenvalues, and combines the results.

### 3.6. Arnoldi and Lanczos Algorithms

These iterative methods are used to compute a few eigenvalues and eigenvectors of large sparse matrices. They are widely used in scientific computing.

## 4. Comparison of Algorithms

Method	Advantages
Characteristic Polynomial	Straightforward for small matrices
Power Iteration	Simple and efficient for largest eigenvalue
QR Algorithm	Finds all eigenvalues
Jacobi Method	Suitable for symmetric matrices
Arnoldi/Lanczos	Efficient for large sparse matrices

Table 1: Comparison of Eigenvalue Algorithms (Method and Advantages)

## 5. Conclusion

Eigenvalues are fundamental in understanding the behavior of linear systems, from stability analysis to data science. The choice of an algorithm for finding eigenvalues depends on the matrix's properties, such as size, sparsity, and symmetry. For small matrices, direct methods like the characteristic polynomial method are effective. For large matrices, iterative methods like the QR algorithm or Arnoldi method are preferred due to their efficiency and scalability.

## Why Choose the QR Algorithm?

The QR Algorithm was chosen for computing eigenvalues due to its versatility and efficiency in handling a wide range of matrices. While other methods have their merits. The other algorithms like Arnoldi or Lanczos are better suited for very large sparse matrices, the QR Algorithm is ideal for the matrix sizes typically encountered in many practical applications.

## Introduction to QR Algorithm

The QR Algorithm is a numerical method for computing the eigenvalues of a square matrix. It is based on the iterative decomposition of a matrix  $A$  into the product of an orthogonal matrix  $Q$  and an upper triangular matrix  $R$ , followed by recomputing  $A = RQ$ . This process continues until the matrix converges to a form where the eigenvalues can be extracted from the diagonal.

## Why is the QR Algorithm Useful?

The QR Algorithm is widely used in numerical linear algebra for several reasons:

- **Efficiency:** It is computationally efficient for large matrices.
- **Stability:** The orthogonal nature of  $Q$  ensures numerical stability.
- **Generality:** The algorithm is applicable to both symmetric and non-symmetric matrices.
- **Practicality:** It forms the basis for eigenvalue solvers in software libraries such as LAPACK and MATLAB.

## Steps in the QR Algorithm

**Hessenberg Reduction and its Role in Eigenvalue Computation Using the QR Algorithm**

## What is Hessenberg Reduction?

A matrix  $A$  is said to be in **Hessenberg form** if:

$$A = \begin{bmatrix} a_{11} & a_{12} & a_{13} & \cdots & a_{1n} \\ a_{21} & a_{22} & a_{23} & \cdots & a_{2n} \\ 0 & a_{32} & a_{33} & \cdots & a_{3n} \\ \vdots & \vdots & \ddots & \ddots & \vdots \\ 0 & 0 & \cdots & a_{n-1,n-1} & a_{n-1,n} \\ 0 & 0 & \cdots & 0 & a_{nn} \end{bmatrix},$$

where all elements below the first sub-diagonal are zero. Hessenberg reduction transforms a general square matrix  $A$  into this form via unitary (orthogonal for real matrices) similarity transformations:

$$H = Q^\dagger A Q,$$

where  $Q$  is a unitary matrix,  $Q^\dagger Q = I$ , and  $H$  is the resulting Hessenberg matrix.

## Why Hessenberg Reduction?

- The QR decomposition of a Hessenberg matrix is much faster than that of a general matrix because of its sparse structure.
- The eigenvalues of the matrix are preserved during Hessenberg reduction.
- It reduces the computational complexity of the QR algorithm from  $O(n^3)$  per iteration for a general matrix to  $O(n^2)$  for a Hessenberg matrix.

## Hessenberg Reduction and QR Algorithm

The QR algorithm iteratively finds the eigenvalues of a matrix by decomposing it into the product of an orthogonal matrix  $Q_k$  and an upper triangular matrix  $R_k$ :

$$A_k = Q_k R_k,$$

and updating  $A_{k+1}$  as:

$$A_{k+1} = R_k Q_k.$$

When the input matrix  $A$  is reduced to Hessenberg form  $H$  before the QR algorithm:

1. Hessenberg reduction simplifies the matrix to have a sparse structure with zeros below the first sub-diagonal.
2. The QR decomposition of  $H$  can be performed more efficiently, as  $R_k$  inherits the upper triangular form of  $H$ , and  $Q_k$  can be computed with reduced operations.
3. The iterative process eventually converges to a quasi-upper triangular matrix, where:
  - Diagonal elements represent real eigenvalues.
  - $2 \times 2$  blocks represent complex conjugate eigenvalue pairs.

## Steps in Eigenvalue Computation

1. **Input Matrix  $A$ :** The algorithm starts with the given square matrix  $A$ .
2. **Hessenberg Reduction:** Transform  $A$  to Hessenberg form  $H$  using Givens rotations or Householder reflections.
3. **QR Algorithm:**
  - (a) Compute  $Q_k$  and  $R_k$  such that  $H_k = Q_k R_k$ .
  - (b) Update  $H_{k+1} = R_k Q_k$ .
  - (c) Check convergence: When all sub-diagonal elements are below a pre-defined tolerance, the iteration stops.
4. **Eigenvalue Extraction:**
  - Diagonal elements of the final matrix represent real eigenvalues.
  - For  $2 \times 2$  blocks, solve the characteristic equation:

$$\lambda^2 - (\text{trace})\lambda + (\text{determinant}) = 0.$$

## Advantages of Hessenberg Reduction in QR Algorithm

- **Computational Efficiency:** Reduces the complexity of each QR iteration, especially for large matrices.
- **Numerical Stability:** The orthogonal transformations preserve the eigenvalues and improve stability.
- **Flexibility:** Works seamlessly with both real and complex matrices.

## Expected Output

For the input matrix:

$$A = \begin{bmatrix} 4 & 2 & 1 \\ 0 & 3 & -1 \\ 2 & 0 & 2 \end{bmatrix}$$

The eigenvalues computed using the QR algorithm are:

$$\lambda_1 = 0.80417665+0.j, \quad \lambda_2 = 4.09791167+0.78500326j, \quad \lambda_3 = 4.09791167-0.78500326j$$

## Conclusion

The QR Algorithm is a robust and efficient method for computing the eigenvalues of square matrices. Its iterative approach and reliance on orthogonal transformations ensure both accuracy and stability, making it a cornerstone of numerical linear algebra.