CE 601: Numerical Methods

Lecture 41

Eigen Values & Eigen Vectors

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We have studied earlier about system of linear equations

i.e.
$$[A]{x} = {b}$$

- \rightarrow A linear system will be homogeneous, say if $[B]{x} = 0$
- → This homogeneous sytem, if we can represent

as
$$[A - \lambda I] \{x\} = 0 \rightarrow (1A)$$

where $I \rightarrow$ unit matrix

 $A \rightarrow$ another square matrix

 $\lambda \rightarrow$ some scalar values

Then that means that $[B] = [A - \lambda I]$

or,
$$[A]{x} = \lambda {x} \rightarrow (1B)$$

i.e. the matrix [A] transforms the vector $\{x\}$ by simply multiplying with a scalar quantity.

Eq.(1) has trivial solution $\{x\} = 0$.

But this is of not interest to us.

Therefore, we need to make $[A - \lambda I]$ as singular

i.e.
$$\det(A - \lambda I) = 0$$
.

The parameters $'\lambda'$ that can make the matrix $[A - \lambda I]$ as singular are called eigen values.

Again in
$$[A]{x} = \lambda {x}$$

- \rightarrow For different values of λ , we get different vectors $\{x\}$ i.e. for $\lambda_1 \rightarrow \{x\}_1, \lambda_2 \rightarrow \{x\}_2, \dots$
- \rightarrow These vectors that are solutions of $[A]\{x\} = \lambda \{x\}$ are called eigen vactors.

 \rightarrow To find eigen values:

$$\det(A - \lambda I) = 0$$

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i.e.
$$\begin{bmatrix} (a_{11} - \lambda) & a_{12} & \cdots & a_{1n} \\ a_{21} & (a_{22} - \lambda) & \cdots & a_{2n} \\ \vdots & \vdots & \ddots & \vdots \\ a_{n1} & a_{n2} & \cdots & (a_{nn} - \lambda) \end{bmatrix} \begin{bmatrix} x_1 \\ x_2 \\ \vdots \\ x_n \end{bmatrix} = 0$$

determinant is some polynomial in λ

$$p(\lambda) = (-1)^n \left[\lambda^n + c_1 \lambda^{n-1} + \dots + c_{n-1} \lambda + c_n \right] = 0$$

Solving this non-linear polynomial will give you \rightarrow The Eigen values.

Numerical Methods to Determine Eigen Values & Eigen Vectors

→ The traditional method to solve non-linear equations as discussed in earlier lectures can be employed to solve non-linear polynomial

$$(-1)^{n} \left[\lambda^{n} + c_{1} \lambda^{n-1} + c_{2} \lambda^{n-2} + \cdots + c_{n-1} \lambda + c_{n} \right] = 0$$

These are other methods as well.

I. Power Method

- * This is an iterative procedure.
- * It is based on repetitive multiplication of a trial eigen vector $\{x\}^{(0)}$ with a scaling factor.
- * We assume a trial eigen vector.
- * The normalisation is done on contents of $\{x\}$.
- * The procedure will determine the dominant eigen pair $\left[\lambda_{\max}, \{x\}\right]$.

 \rightarrow If for an $n \times n$ non-singular matrix [A] if the eigen values $\lambda_1 > \lambda_2 > \lambda_3 > \cdots > \lambda_n$, then $\lambda_1 \to$ dominant eigen value and $\{x\}_1 \to$ eigen vector corresponding to domiant eigen value. The procedure is

- (i) Assume $\{x\}_{1}^{(0)}$
- (ii) Assign $\{y\}^{(1)} = [A]\{x\}^{(0)}$
- (iii) Identity $\lambda_1^{(1)} = \max \left| y_i^{(1)} \right|$
- (iv) Evaluate $\{x\}^{(1)} = \frac{1}{\lambda_1^{(1)}} \{y\}^{(1)}$

(v) Again evaluate
$$\{y\}^{(2)} = [A]\{x\}^{(1)}$$

Determine as above $\lambda_1^{(2)} = \max |y_i^{(2)}|$

$$\{x\}^{(2)} = \frac{1}{\lambda_1^{(2)}} \{y\}^{(2)}$$

(vi) Continue the process till convergence

i.e.
$$\frac{\left|\lambda_{1}^{(k+1)} - \lambda_{1}^{(k)}\right|}{\left|\lambda_{1}^{(k)}\right|} \le \varepsilon$$
 or as per requirement.

Example:

Compute the largest eigen value of
$$\begin{bmatrix} A \end{bmatrix} = \begin{bmatrix} 2 & -1 & 0 & 0 \\ -1 & 4 & -1 & 0 \\ 0 & -1 & 4 & -1 \\ 0 & 0 & -1 & 2 \end{bmatrix}$$

Answer. The matrix A is 4×4 .

... There are four eigen values for [A]. The homogeneous system is $[A - \lambda I]\{x\} = 0$. Let $|\lambda_1| \ge |\lambda_2| \ge |\lambda_3| \ge |\lambda_4|$. To find λ_1

Assign
$$\left\{x\right\}_{1}^{(0)} = \begin{cases} 1\\0\\0\\0 \end{cases}$$

$$\{y\}^{(1)} = [A]\{x\}^{(0)} = \begin{bmatrix} 2 & -1 & 0 & 0 \\ -1 & 4 & -1 & 0 \\ 0 & -1 & 4 & -1 \\ 0 & 0 & -1 & 2 \end{bmatrix} \begin{bmatrix} 1 \\ 0 \\ 0 \\ 0 \end{bmatrix} = \begin{bmatrix} 2 \\ -1 \\ 0 \\ 0 \end{bmatrix}$$

i.e.
$$\{y\}^{(1)} = \begin{cases} 2 \\ -1 \\ 0 \\ 0 \end{cases}, \therefore \lambda_1^{(1)} = \max |y_i^{(i)}| = 2$$

$$\{x\}^{(1)} = \frac{1}{\lambda_1^{(1)}} \{y\}^{(1)} = \frac{1}{2} \begin{cases} 2 \\ -1 \\ 0 \\ 0 \end{cases} = \begin{cases} 1 \\ -0.5 \\ 0 \\ 0 \end{cases}$$

$$\{y\}^{(2)} = [A]\{x\}^{(1)} = \begin{bmatrix} 2 & -1 & 0 & 0 \\ -1 & 4 & -1 & 0 \\ 0 & -1 & 4 & -1 \\ 0 & 0 & -1 & 2 \end{bmatrix} \begin{bmatrix} 1 \\ -0.5 \\ 0 \\ 0 \end{bmatrix} = \begin{bmatrix} 2.5 \\ -3.0 \\ 0.5 \\ 0 \end{bmatrix}$$

$$\therefore \lambda_1^{(2)} = 3.0$$

and
$$\{x\}^{(2)} = \frac{1}{3.0} \begin{cases} 2.5 \\ -3.0 \\ 0.5 \\ 0 \end{cases} = \begin{cases} 0.83373 \\ -1.0000 \\ 0.16667 \\ 0.00000 \end{cases}$$

The process goes on till there is no change in value of λ for successive iteration.

After many iterations,

$$\lambda_1 = 5.30269$$

$$\left\{x\right\}_{1} = \begin{cases} 0.30279 \\ -1.0000 \\ 0.99993 \\ -0.30274 \end{cases}$$

II. Inverse Power Method

For an $n \times n$ matrix [A] with unique eigen value of minimum absolute magnitude, an be obtained by power method

 $[A] \rightarrow \lambda_i$ (Eigen values for matrix A)

for $[A]^{-1} \to \frac{1}{\lambda_i}$ (Eigen values of matrix A^{-1})

... Applying power method on $[A]^{-1}$ will give you the dominant eigen value of $[A]^{-1}$, which happens to be the minimum eigen value of [A]. $[A]\{x\} = \lambda \{x\}$

$$[A]^{-1}[A]\{x\} = \lambda [A]^{-1}\{x\}$$

$$\therefore [A]^{-1} \{x\} = \frac{1}{\lambda} [I] \{x\} = \frac{1}{\lambda} \{x\}$$

III. Shifted Power Method

$$[A]\{x\} = \lambda \{x\}$$

If you have some scalar value 's' and you are interested to find the closest eigen value of A corresponding to 's'.

Then let,
$$s[I]{x} = s{x}$$

$$[A]{x} - s{x} = \lambda {x} - s{x}$$

$$\Rightarrow [A - sI]\{x\} = (\lambda - s)\{x\}$$

 $[A-sI] \rightarrow \text{shifted matrix}.$

$$[A-sI]_{shifted} \{x\} = \lambda_{shifted} \{x\},\,$$

where, $\lambda_{shifted}$ is the shifted eigen value.

Apply power method to detrmine the dominant eigen value of the shifted system.