



MATHEMATICS-II

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UNIT-1

Vector Calculus

INTRODUCTION

- In this chapter, vector differential calculus is considered, which extends the basic concepts of differential calculus, such as, continuity and differentiability to vector functions in a simple and natural way. Also, the new concepts of gradient, divergence and curl are introduced.
- *Example:* $\mathbf{i}, \mathbf{j}, \mathbf{k}$ are unit vectors.

VECTOR DIFFERENTIAL OPERATOR

- The vector differential operator Δ is defined as $\Delta = i \partial/\partial x + j \partial/\partial y + k \partial/\partial z$. This operator possesses properties analogous to those of ordinary vectors as well as differentiation operator. We will define now some quantities known as gradient, divergence and curl involving this operator.

GRADIENT

- Let $f(x,y,z)$ be a scalar point function of position defined in some region of space. Then gradient of f is denoted by $\text{grad } f$ or Δf and is defined as

$$\text{grad } f = i \frac{\partial f}{\partial x} + j \frac{\partial f}{\partial y} + k \frac{\partial f}{\partial z}$$

- *Example:* If $f=2x+3y+5z$ then $\text{grad } f = 2i+3j+5k$

DIRECTIONAL DERIVATIVE

- The directional derivative of a scalar point function f at a point $P(x,y,z)$ in the direction of g at P and is defined as $\text{grad } g / |\text{grad } g| \cdot \text{grad } f$
- *Example:* The directional derivative of $f=xy+yz+zx$ in the direction of the vector $i+2j+2k$ at the point $(1,2,0)$ is $10/3$

DIVERGENCE OF A VECTOR

- Let f be any continuously differentiable vector point function. Then divergence of f and is written as $\text{div } f$ and is defined as

$$\text{Div } f = \partial f_1 / \partial x + j \partial f_2 / \partial y + k \partial f_3 / \partial z$$

- *Example 1:* The divergence of a vector $2xi+3yj+5zk$ is 10
- *Example 2:* The divergence of a vector $f=xy^2i+2x^2yzj-3yz^2k$ at $(1,-1,1)$ is 9

SOLENOIDAL VECTOR

- A vector point function f is said to be solenoidal vector if its divergent is equal to zero i.e., $\text{div } f=0$
- *Example 1:* The vector $f=(x+3y)i+(y-2z)j+(x-2z)k$ is solenoidal vector.
- *Example 2:* The vector $f=3y^4z^2i+z^3x^2j-3x^2y^2k$ is solenoidal vector.

CURL OF A VECTOR

- Let f be any continuously differentiable vector point function. Then the vector function curl of f is denoted by $\text{curl } f$ and is defined as $\text{curl } f = i \frac{\partial f}{\partial x} + j \frac{\partial f}{\partial y} + k \frac{\partial f}{\partial z}$
- *Example 1:* If $f = xy^2i + 2x^2yzj - 3yz^2k$ then $\text{curl } f$ at $(1, -1, 1)$ is $-i - 2k$
- *Example 2:* If $r = xi + yj + zk$ then $\text{curl } r$ is 0

IRROTATIONAL VECTOR

- Any motion in which curl of the velocity vector is a null vector i.e., $\text{curl } \mathbf{v} = 0$ is said to be irrotational. If \mathbf{f} is irrotational, there will always exist a scalar function $f(x,y,z)$ such that $\mathbf{f} = \text{grad } g$. This g is called scalar potential of \mathbf{f} .
- *Example:* The vector $\mathbf{f} = (2x+3y+2z)\mathbf{i} + (3x+2y+3z)\mathbf{j} + (2x+3y+3z)\mathbf{k}$ is irrotational vector.

VECTOR INTEGRATION

- INTRODUCTION: In this chapter we shall define line, surface and volume integrals which occur frequently in connection with physical and engineering problems. The concept of a line integral is a natural generalization of the concept of a definite integral of $f(x)$ exists for all x in the interval $[a,b]$

WORK DONE BY A FORCE

- If F represents the force vector acting on a particle moving along an arc AB , then the work done during a small displacement $F \cdot dr$. Hence the total work done by F during displacement from A to B is given by the line integral $\int F \cdot dr$
- *Example:* If $f = (3x^2 + 6y)i - 14yzj + 20xz^2k$ along the lines from $(0,0,0)$ to $(1,0,0)$ then to $(1,1,0)$ and then to $(1,1,1)$ is $23/3$

SURFACE INTEGRALS

- The surface integral of a vector point function F expresses the normal flux through a surface. If F represents the velocity vector of a fluid then the surface integral $\int F \cdot n \, dS$ over a closed surface S represents the rate of flow of fluid through the surface.
- *Example:* The value of $\int F \cdot n \, dS$ where $F = 18zi - 12j + 3yk$ and S is the part of the surface of the plane $2x + 3y + 6z = 12$ located in the first octant is 24.

VOLUME INTEGRAL

- Let $\mathbf{f}(\mathbf{r}) = f_1\mathbf{i} + f_2\mathbf{j} + f_3\mathbf{k}$ where f_1, f_2, f_3 are functions of x, y, z .
We know that $dv = dx dy dz$. The volume integral is given by
$$\int \mathbf{f} dv = \iiint (f_1\mathbf{i} + f_2\mathbf{j} + f_3\mathbf{k}) dx dy dz$$
- *Example:* If $\mathbf{F} = 2xzi\mathbf{i} - xj\mathbf{j} + y^2k\mathbf{k}$ then the value of $\int \mathbf{f} dv$ where v is the region bounded by the surfaces $x=0, x=2, y=0, y=6, z=x^2, z=4$ is $128\mathbf{i} - 24\mathbf{j} - 384\mathbf{k}$

VECTOR INTEGRAL THEOREMS

- In this chapter we discuss three important vector integral theorems.
- 1) Gauss divergence theorem
- 2) Green's theorem
- 3) Stokes theorem

GAUSS DIVERGENCE THEOREM

- This theorem is the transformation between surface integral and volume integral. Let S be a closed surface enclosing a volume v . If f is a continuously differentiable vector point function, then
- $\int \text{div } f \, dv = \int f \cdot n \, dS$
- Where n is the outward drawn normal vector at any point of S .

GREEN'S THEOREM

- This theorem is transformation between line integral and double integral. If S is a closed region in xy plane bounded by a simple closed curve C and in M and N are continuous functions of x and y having continuous derivatives in R , then
- $\int Mdx + Ndy = \iint (\partial N / \partial x - \partial M / \partial y) dx dy$

STOKES THEOREM

- This theorem is the transformation between line integral and surface integral. Let S be a open surface bounded by a closed, non-intersecting curve C . If F is any differentiable vector point function then
- $\int F \cdot dr = \int \text{Curl } F \cdot n \, ds$

UNIT-2

Fourier Series and Fourier Transform

INTRODUCTION

- Suppose that a given function $f(x)$ defined in $(-\pi, \pi)$ or $(0, 2\pi)$ or in any other interval can be expressed as a trigonometric series as

$$f(x) = a_0/2 + (a_1 \cos x + a_2 \cos 2x + a_3 \cos 3x + \dots + a_n \cos nx) + (b_1 \sin x +$$

$$b_2 \sin 2x + \dots + b_n \sin nx) + \dots$$

$$f(x) = a_0/2 + \sum (a_n \cos nx + b_n \sin nx)$$

- Where a and b are constants with in a desired range of values of the variable such series is known as the fourier series for $f(x)$ and the constants a_0, a_n, b_n are called fourier coefficients of $f(x)$
- It has period 2π and hence any function represented by a series of the above form will also be periodic with period 2π

POINTS OF DISCONTINUITY

- In deriving the Euler's formulae for a_0, a_n, b_n it was assumed that $f(x)$ is continuous. Instead a function may have a finite number of discontinuities. Even then such a function is expressible as a fourier series.

DISCONTINUITY FUNCTION

- For instance, let the function $f(x)$ be defined by
- $f(x) = \phi(x), c < x < x_0$
- $= \Psi(x), x_0 < x < c + 2\pi$
- where x_0 is the point of discontinuity in $(c, c + 2\pi)$.

DISCONTINUITY FUNCTION

In such cases also we obtain the fourier series for $f(x)$ in the usual way. The values of a_0, a_n, b_n are given by

$$a_0 = 1/\pi \left[\int \phi(x) dx + \int \Psi(x) dx \right]$$

$$a_n = 1/\pi \left[\int \phi(x) \cos nx dx + \int \Psi(x) \cos nx dx \right]$$

$$b_n = 1/\pi \left[\int \phi(x) \sin nx dx + \int \Psi(x) \sin nx dx \right]$$

EULER'S FORMULAE

- The fourier series for the function $f(x)$ in the interval $C \leq x \leq C+2\pi$ is given by

$$f(x) = a_0/2 + \sum (a_n \cos nx + b_n \sin nx)$$

where $a_0 = 1/\pi \int f(x) dx$

$$a_n = 1/\pi \int f(x) \cos nx dx$$

$$b_n = 1/\pi \int f(x) \sin nx dx$$

These values of a_0 , a_n , b_n are known as Euler's formulae

EVEN AND ODD FUNCTIONS

- A function $f(x)$ is said to be even if $f(-x)=f(x)$ and odd if $f(-x) = -f(x)$.
- If a function $f(x)$ is even in $(-\pi, \pi)$, its fourier series expansion contains only cosine terms, and their coefficients are a_0
- and a_n .
- $f(x) = a_0/2 + \sum a_n \cos nx$
- where $a_0 = 2/\pi \int f(x) dx$
- $a_n = 2/\pi \int f(x) \cos nx dx$

ODD FUNCTION

- When $f(x)$ is an odd function in $(-\pi, \pi)$ its fourier expansion contains only sine terms.
- And their coefficient is b_n
- $f(x) = \sum b_n \sin nx$
- where $b_n = 2/\pi \int f(x) \sin nx \, dx$

HALF RANGE FOURIER SERIES

- THE SINE SERIES: If it be required to express $f(x)$ as a sine series in $(0, \pi)$, we define an odd function $f(x)$ in $(-\pi, \pi)$, identical with $f(x)$ in $(0, \pi)$.
- Hence the half range sine series $(0, \pi)$ is given by
- $$f(x) = \sum b_n \sin nx$$
- Where $b_n = \frac{2}{\pi} \int_0^\pi f(x) \sin nx \, dx$

HALF RANGE SERIES

- The cosine series: If it be required to express $f(x)$ as a cosine series, we define an even function $f(x)$ in $(-\pi, \pi)$, identical with $f(x)$ in $(0, \pi)$, i.e we extend the function reflecting it with respect to the y-axis, so that $f(-x)=f(x)$.

HALF RANGE COSINE SERIES

- Hence the half range series in $(0, \pi)$ is given by
- $f(x) = a_0/2 + \sum a_n \cos nx$
- where $a_0 = 2/\pi \int f(x) dx$
- $a_n = 2/\pi \int f(x) \cos nx dx$

CHANGE OF INTERVAL

- So far we have expanded a given function in a Fourier series over the interval $(-\pi, \pi)$ and $(0, 2\pi)$ of length 2π . In most engineering problems the period of the function to be expanded is not 2π but some other quantity say $2l$. In order to apply earlier discussions to functions of period $2l$, this interval must be converted to the length 2π .

PERIODIC FUNCTION

- Let $f(x)$ be a periodic function with period $2l$ defined in the interval $c < x < c + 2l$. We must introduce a new variable z such that the period becomes 2π .

CHANGE OF INTERVAL

- The fourier expansion in the change of interval is given by
- $f(x) = a_0/2 + \sum a_n \cos n\pi x/l + \sum b_n \sin n\pi x/l$
- Where $a_0 = 1/l \int f(x) dx$
- $a_n = 1/l \int f(x) \cos n\pi x/l dx$
- $b_n = 1/l \int f(x) \sin n\pi x/l dx$

EVEN AND ODD FUNCTION

- Fourier cosine series : Let $f(x)$ be even function in $(-l, l)$ then
- $f(x) = a_0/2 + \sum a_n \cos n\pi x/l$
- where $a_0 = 2/l \int f(x) dx$
- $a_n = 2/l \int f(x) \cos n\pi x/l dx$

FOURIER SINE SERIES

- Fourier sine series : Let $f(x)$ be an odd function in $(-l,l)$ then
- $f(x) = \sum b_n \sin n\pi x/l$
- where $b_n = 2/l \int f(x) \sin n\pi x/l \, dx$
- Once ,again here we remarks that the even nature or odd nature of the function is to be considered only when we deal with the interval $(-l,l)$.

HALF-RANGE EXPANSION


- Cosine series: If it is required to expand $f(x)$ in the interval $(0, l)$ then we extend the function reflecting in the y-axis, so that $f(-x) = f(x)$. We can define a new function $g(x)$ such that $f(x) = a_0/2 + \sum a_n \cos n\pi x/l$
- where $a_0 = 2/l \int f(x) dx$
- $a_n = 2/l \int f(x) \cos n\pi x/l dx$

HALF RANGE SINE SERIES

- Sine series : If it be required to expand $f(x)$ as a sine series in $(0,l)$, we extend the function reflecting it in the origin so that $f(-x) = f(x)$. we can define the fourier series in $(-l,l)$ then,
- $$f(x) = a_0/2 + \sum b_n \sin n\pi x/l$$
- where $b_n = 2/l \int f(x) \sin n\pi x/l \, dx$

FOURIER INTEGRAL TRANSFORMS

- INTRODUCTION: A transformation is a mathematical device which converts or changes one function into another function. For example, differentiation and integration are transformations.
- In this we discuss the application of finite and infinite fourier integral transforms which are mathematical devices from which we obtain the solutions of boundary value.

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- We obtain the solutions of boundary value problems related to engineering. For example conduction of heat, free and forced vibrations of a membrane, transverse vibrations of a string, transverse oscillations of an elastic beam etc.

- DEFINITION: The integral transforms of a function $f(t)$ is defined by
- $F(p) = I[f(t)] = \int f(t) k(p, t) dt$
- Where $k(p, t)$ is called the kernel of the integral transform and is a function of p and t .

FOURIER COSINE AND SINE INTEGRAL

- When $f(t)$ is an odd function $\cos pt, f(t)$ is an odd function and $\sin pt f(t)$ is an even function. So the first integral in the right side becomes zero. Therefore we get
- $f(x) = 2/\pi \int \sin px$

FOURIER COSINE AND SINE INTEGRAL

- When $f(t)$ is an odd function $\cos pt, f(t)$ is an odd function and $\sin pt f(t)$ is an even function. So the first integral in the right side becomes zero. Therefore we get
- $$f(x) = \frac{2}{\pi} \int \sin px \int f(t) \sin pt \, dt \, dp$$
- which is known as FOURIER SINE INTEGRAL.

- When $f(t)$ is an even function, the second integral in the right side becomes zero. Therefore we get
- $f(x) = 2/\pi \int \cos px \int f(t) \cos pt \, dt \, dp$
- which is known as FOURIER COSINE INTEGRAL.

FOURIER INTEGRAL IN COMPLEX FORM

- Since $\cos p(t-x)$ is an even function of p , we have
- $f(x) = \frac{1}{2\pi} \iint e^{ip(t-x)} f(t) dt dp$
- which is the required complex form.

INFINITE FOURIER TRANSFORM

- The fourier transform of a function $f(x)$ is given by
- $F\{f(x)\} = F(p) = \int f(x) e^{ipx} dx$
- The inverse fourier transform of $F(p)$ is given by
- $f(x) = 1/2\pi \int F(p) e^{-ipx} dp$

FOURIER SINE TRANSFORM

- The finite Fourier sine transform of $f(x)$ when $0 < x < l$, is defined as
- $F_s\{f(x)\} = F_s(n) \sin(n\pi x)/l \, dx$ where n is an integer and the function $f(x)$ is given by
- $f(x) = 2/l \sum F_s(n) \sin(n\pi x)/l$ is called the Inverse finite Fourier sine transform $F_s(n)$

FOURIER COSINE TRANSFORM

- We have $f(x) = \frac{2}{\pi} \int \cos px \int f(t) \cos pt \, dt \, dp$
- Which is the fourier cosine integral .Now
- $F_c(p) = \int f(x) \cos px \, dx$ then
- $f(x)$ becomes $f(x) = \frac{2}{\pi} \int F_c(p) \cos px \, dp$ which is the fourier cosine transform.

PROPERTIES

- Linear property of Fourier transform
- Change of Scale property
- Shifting property
- Modulation property

UNIT-3

Interpolation and Curve Fitting

Finite difference methods

Let $(x_i, y_i), i=0, 1, 2, \dots, n$ be the equally spaced data of the unknown function $y=f(x)$ then much of the $f(x)$ can be extracted by analyzing the differences of $f(x)$.

$$\text{Let } x_1 = x_0 + h$$

$$x_2 = x_0 + 2h$$

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$x_n = x_0 + nh$ be equally spaced points where the function value of $f(x)$

be $y_0, y_1, y_2, \dots, y_n$

Symbolic operators

Forward shift operator(E) :

It is defined as $Ef(x)=f(x+h)$ (or) $Ey_x = y_{x+h}$

The second and higher order forward shift operators are defined in similar manner as follows

$$E^2f(x)= E(Ef(x))= E(f(x+h))= f(x+2h)= y_{x+2h}$$

$$E^3f(x)= f(x+3h)$$

•

•

$$E^kf(x)= f(x+kh)$$

Backward shift operator(E^{-1}) :

It is defined as $E^{-1}f(x)=f(x-h)$ (or) $Ey_x = y_{x-h}$

The second and higher order backward shift operators are defined in similar manner as follows

$$E^{-2}f(x)= E^{-1}(E^{-1}f(x))= E^{-1}(f(x-h))= f(x-2h)= y_{x-2h}$$

$$E^{-3}f(x)= f(x-3h)$$

.

.

$$E^{-k}f(x)= f(x-kh)$$

Forward difference operator (Δ) :

The first order forward difference operator of a function $f(x)$ with increment h in x is given by

$$\Delta f(x) = f(x+h) - f(x) \quad (\text{or}) \quad \Delta f_k = f_{k+1} - f_k ; k=0,1,2,\dots$$

$$\Delta^2 f(x) = \Delta[\Delta f(x)] = \Delta[f(x+h) - f(x)] = \Delta f_{k+1} - \Delta f_k ; k=0,1,2,\dots$$

.....

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Relation between E and Δ :

$$\Delta f(x) = f(x+h) - f(x)$$

$$= Ef(x) - f(x)$$

$$= (E-1)f(x)$$

$$[Ef(x) = f(x+h)]$$

$$\Delta = E - 1 \quad E = 1 + \Delta$$

Backward difference operator (∇) :

The first order backward difference operator of a function $f(x)$ with increment h in x is given by

$$\nabla f(x) = f(x) - f(x-h) \quad (\text{or}) \quad \nabla f_k = f_{k+1} - f_k ; k=0,1,2,\dots$$

$$\nabla f(x) = \nabla[\nabla f(x)] = [f(x+h) - f(x)] = \nabla f_{k+1} - \nabla f_k ; k=0,1,2,\dots$$

.....
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Relation between E and ∇ :

$$\nabla f(x) = f(x+h) - f(x)$$

$$= Ef(x) - f(x)$$

$$= (E-1)f(x)$$

$$[Ef(x) = f(x+h)]$$

$$\nabla = E-1 \quad E = 1 + \nabla$$

Central difference operator (δ) :

The central difference operator is defined as

$$\delta f(x) = f(x+h/2) - f(x-h/2)$$

$$\delta f(x) = E^{1/2}f(x) - E^{-1/2}f(x)$$

$$= [E^{1/2} - E^{-1/2}]f(x)$$

$$\delta = E^{1/2} - E^{-1/2}$$



INTERPOLATION : The process of finding a missed value in the given table values of X, Y.

FINITE DIFFERENCES : We have three finite differences

1. Forward Difference
2. Backward Difference
3. Central Difference

RELATIONS BETWEEN THE OPERATORS IDENTITIES:

1. $\Delta = E - 1$ or $E = 1 + \Delta$

2. $\nabla = 1 - E^{-1}$

3. $\delta = E^{1/2} - E^{-1/2}$

4. $\mu = \frac{1}{2} (E^{1/2} - E^{-1/2})$

5. $\Delta = E \nabla = \nabla E = \delta E^{1/2}$

6. $(1 + \Delta)(1 - \nabla) = 1$

Newtons Forward interpolation formula :

$$y=f(x)=f(x_0+ph)= y_0 + p\Delta y_0 + \frac{p(p-1)}{2!} \Delta^2 y_0 + \frac{p(p-1)(p-2)}{3!} \Delta^3 y_0 + \dots + \frac{p(p-1)(p-2) \dots [p-(n-1)]}{n!} \Delta^n y_0.$$

Newtons Backward interpolatin formula :

$$y=f(x)=f(x_n+ph)= y_n + p\nabla y_n + \frac{p(p+1)}{2!} \nabla^2 y_n + \frac{p(p+1)(p+2)}{3!} \nabla^3 y_n + \dots + \frac{p(p+1)(p+2) \dots [p+(n-1)]}{n!} \nabla^n y_n.$$

GAUSS INTERPOLATION

The Gauss forward interpolation is given by $y_p = y_0 + p \Delta y_0 + \frac{p(p-1)}{2!} \Delta^2 y_{-1} + \frac{(p+1)p(p-1)}{3!} \Delta^3 y_{-1} + \frac{(p+1)p(p-1)(p-2)}{4!} \Delta^4 y_{-2} + \dots$

The Gauss backward interpolation is given by

$$y_p = y_0 + p \Delta y_{-1} + \frac{p(p+1)}{2!} \Delta^2 y_{-1} + \frac{(p+1)p(p-1)}{3!} \Delta^3 y_{-2} + \frac{(p+2)(p+1)p(p-1)}{4!} \Delta^4 y_{-2} + \dots$$

INTERPOLATION WITH UNEQUAL INTERVALS:

The various interpolation formulae Newton's forward formula, Newton's backward formula possess can be applied only to equal spaced values of argument. It is therefore, desirable to develop interpolation formula for unequally spaced values of x . We use Lagrange's interpolation formula.

The Lagrange's interpolation formula is given by

$$Y = \frac{(X-X_1)(X-X_2)\dots\dots(X-X_n)}{(X_0-X_1)(X_0-X_2)\dots\dots(X_0-X_n)} Y_0 +$$


$$\frac{(X-X_0)(X-X_2)\dots\dots(X-X_n)}{(X_1-X_0)(X_1-X_2)\dots\dots(X_1-X_n)} Y_1 + \dots\dots\dots$$

$$\frac{(X-X_0)(X-X_1)\dots\dots(X-X_{n-1})}{(X_n-X_0)(X_n-X_1)\dots\dots(X_n-X_{n-1})} Y_n +$$



CURVE FITTING

INTRODUCTION : In interpolation, We have seen that when exact values of the function $Y=f(x)$ is given we fit the function using various interpolation formulae. But sometimes the values of the function may not be given. In such cases, the values of the required function may be taken experimentally. Generally these expt. Values contain some errors. Hence by using these experimental values . We can fit a curve just approximately which is known as approximating curve.



Now our aim is to find this approximating curve as much best as through minimizing errors of experimental values this is called best fit otherwise it is a bad fit.

In brief by using experimental values the process of establishing a mathematical relationship between two variables is called CURVE FITTING.

METHOD OF LEAST SQUARES

Let $y_1, y_2, y_3, \dots, y_n$ be the experimental values of y .
 $f(x_1), f(x_2), \dots, f(x_n)$ be the exact values of the function $y=f(x)$.
Corresponding to the values of $x=x_0, x_1, x_2, \dots, x_n$.
Now error = experimental values – exact value. If we denote the corresponding errors of y_1, y_2, \dots, y_n as $e_1, e_2, e_3, \dots, e_n$, then $e_1 = y_1 - f(x_1)$, $e_2 = y_2 - f(x_2)$

- $e_3 = y_3 - f(x_3) \dots e_n = y_n - f(x_n)$. These errors $e_1, e_2, e_3, \dots, e_n$, may be either positive or negative. For our convenient to make all errors into +ve to the square of errors i.e $e_1^2, e_2^2, \dots, e_n^2$. In order to obtain the best fit of curve we have to make the sum of the squares of the errors as much minimum i.e $e_1^2 + e_2^2 + \dots + e_n^2$ is minimum.

METHOD OF LEAST SQUARES

- Let $S = e_1^2 + e_2^2 + \dots + e_n^2$, S is minimum. When S becomes as much as minimum. Then we obtain a best fitting of a curve to the given data, now to make S minimum we have to determine the coefficients involving in the curve, so that S minimum. It will be possible when differentiating S with respect to the coefficients involving in the curve and equating to zero.

FITTING OF STRAIGHT LINE

Let $y = a + bx$ be a straight line

By using the principle of least squares for solving the straight line equations.

The normal equations are

$$\sum y = na + b \sum x$$

$$\sum xy = a \sum x + b \sum x^2$$

solving these two normal equations we get the values of a & b , substituting these values in the given straight line equation which gives the best fit.

FITTING OF PARABOLA

Let $y = a + bx + cx^2$ be the parabola or second degree polynomial.

By using the principle of least squares for solving the parabola

The normal equations are

$$\sum y = na + b \sum x + c \sum x^2$$

$$\sum xy = a \sum x + b \sum x^2 + c \sum x^3$$

$$\sum x^2y = a \sum x^2 + b \sum x^3 + c \sum x^4$$

solving these normal equations we get the values of a, b & c , substituting these values in the given parabola which gives the best fit.

FITTING OF AN EXPONENTIAL CURVE

The exponential curve of the form

$$y = a e^{bx}$$

taking log on both sides we get

$$\log_e y = \log_e a + \log_e e^{bx}$$

$$\log_e y = \log_e a + bx \log_e e$$

$$\log_e y = \log_e a + bx$$

- $Y = A + bx$
- Where $Y = \log_e y$, $A = \log_e a$
- This is in the form of straight line equation and this can be solved by using the straight line normal equations we get the values of A & b , for $a = e^A$, substituting the values of a & b in the given curve which gives the best fit.

EXPONENTIAL CURVE

The equation of the exponential form is of the form $y = ab^x$

taking log on both sides we get

$$\log_e y = \log_e a + \log_e b^x$$

$$Y = A + Bx$$

where $Y = \log_e y$, $A = \log_e a$, $B = \log_e b$

this is in the form of the straight line equation which can be solved by using the normal equations we get the values of A & B for $a = e^A$

$b = e^B$ substituting these values in the equation which gives the best fit.

FITTING OF POWER CURVE

Let the equation of the power curve be

$$y = a x^b$$

taking log on both sides we get

$$\log_e y = \log_e a + \log_e x^b$$

$$Y = A + Bx$$

this is in the form of the straight line equation which can be solved by using the normal equations we get the values of A & B, for $a = e^A$ $b = e^B$, substituting these values in the given equation which gives the best fit.

UNIT-4

Solutions of Algebraic and
Transcendental equations and
Linear system of equations

Method 1: Bisection method

If a function $f(x)$ is continuous b/w x_0 and x_1 and $f(x_0)$ & $f(x_1)$ are of opposite signs, then there exist at least one root b/w x_0 and x_1

- Let $f(x_0)$ be $-ve$ and $f(x_1)$ be $+ve$, then the root lies b/w x_0 and x_1 and its approximate value is given by $x_2 = (x_0 + x_1)/2$
- If $f(x_2) = 0$, we conclude that x_2 is a root of the eqn $f(x) = 0$
- Otherwise the root lies either b/w x_2 and x_1 (or) b/w x_2 and x_0 depending on whether $f(x_2)$ is $+ve$ or $-ve$

Then as before, we bisect the interval and repeat the process until the root is known to the desired accuracy

Method 2: Iteration method or successive approximation

Consider the equation $f(x)=0$ which can take in the form

$$x = \phi(x) \text{ -----(1)}$$

where $|\phi'(x)| < 1$ for all values of x .

Taking initial approximation is X_0

we put $x_1 = \phi(x_0)$ and take x_1 is the first approximation

$x_2 = \phi(x_1)$, x_2 is the second approximation

$x_3 = \phi(x_2)$, x_3 is the third approximation

.

.

$x_n = \phi(x_{n-1})$, x_n is the n^{th} approximation

Such a process is called an iteration process

Method 3: Newton-Raphson method or Newton iteration method

Let the given equation be $f(x)=0$

Find $f'(x)$ and initial approximation x_0

The first approximation is $x_1 = x_0 - f(x_0) / f'(x_0)$

The second approximation is $x_2 = x_1 - f(x_1) / f'(x_1)$

.

.

The n^{th} approximation is $x_n = x_{n-1} - f(x_{n-1}) / f'(x_{n-1})$

LU Decomposition Method

$$a_{11}x_1 + a_{12}x_2 + a_{13}x_3 = b_1$$

$$a_{21}x_1 + a_{22}x_2 + a_{23}x_3 = b_2$$

$$a_{31}x_1 + a_{32}x_2 + a_{33}x_3 = b_3$$

$$A = \begin{bmatrix} a_{11} & a_{12} & a_{13} \\ a_{21} & a_{22} & a_{23} \\ a_{31} & a_{32} & a_{33} \end{bmatrix}, X = \begin{bmatrix} x_1 \\ x_2 \\ x_3 \end{bmatrix}, B = \begin{bmatrix} b_1 \\ b_2 \\ b_3 \end{bmatrix}$$

This is in the form $AX=B$ Where

Let $A=LU$ where

$$L = \begin{bmatrix} l_{11} & 0 & 0 \\ l_{21} & l_{22} & 0 \\ l_{31} & l_{32} & l_{33} \end{bmatrix}, U = \begin{bmatrix} u_{11} & u_{12} & u_{13} \\ 0 & u_{22} & u_{23} \\ 0 & 0 & u_{33} \end{bmatrix}$$

Hence $LUX=B$

(a) $LY=B$ (b) $UX=Y$

Solve for Y from (a) then Solve for X from (b)

$LY=B$ Can be solved by forward substitution and $UX=Y$ can be solved by backward substitution

Jacobi's Iteration Method

$$a_1x + b_1y + c_1z = d_1$$

$$a_2x + b_2y + c_2z = d_2$$

$$a_3x + b_3y + c_3z = d_3$$

$$x = \frac{1}{a_1}[d_1 - b_1y - c_1z] = k_1 - l_1y - m_1z$$

$$y = \frac{1}{b_2}[d_2 - a_2x - c_2z] = k_2 - l_2x - m_2z$$

$$z = \frac{1}{c_3}[d_3 - a_3x - b_3y] = k_3 - l_3x - m_3y$$

Substituting these on the right hand side we get second approximations this process is repeated till the difference between two consecutive approximations is negligible or same

Gauss-Seidel iteration Method

$$a_1x + b_1y + c_1z = d_1$$

$$a_2x + b_2y + c_2z = d_2$$

$$a_3x + b_3y + c_3z = d_3$$

$$x = \frac{1}{a_1}[d_1 - b_1y - c_1z] = k_1 - l_1y - m_1z$$

$$y = \frac{1}{b_2}[d_2 - a_2x - c_2z] = k_2 - l_2x - m_2z$$

$$z = \frac{1}{c_3}[d_3 - a_3x - b_3y] = k_3 - l_3x - m_3y$$

as soon as new approximation for an unknown is found it is immediately used in the next step

UNIT-5

Numerical Integration and
Numerical solution of First
order differential equations

NUMERICAL INTEGRATION

- Numerical Integration is a process of finding the value of a definite integral. When a function $y = f(x)$ is not known explicitly. But we give only a set of values of the function $y = f(x)$ corresponding to the same values of x . This process when applied to a function of a single variable is known as a quadrature.

NUMERICAL INTEGRATION

- For evaluating the Numerical Integration we have three important rules i.e
- Trapezoidal Rule
- Simpsons $1/3$ Rule
- Simpsons $3/8$ th Rule

NUMERICAL INTEGRATION

- Trapezoidal Rule : The Trapezoidal Rule of the function $y = f(x)$ is given by
- $\int f(x) dx = \frac{h}{2} (y_0 + y_n) + 2 (y_1 + y_2 + y_3 + \dots + y_{n-1})$
- $\int f(x) dx = \frac{h}{2} (\text{sum of the first and last terms}) + (\text{sum of the remaining terms})$

NUMERICAL INTEGRATION

- Simpson's $1/3$ rd Rule : The Simpson's $1/3$ rd Rule of the function $f(x)$ is given by
- $\int f(x) dx = h/3 (y_0 + y_n) + 4(y_1 + y_3 + y_5 + \dots y_{n-1}) + 2(y_2 + y_4 + y_6 + \dots)$

NUMERICAL INTEGRATION

- Simpson's 3/8 th Rule : The Simpson's 3/8 th rule for the function $f(x)$ is given by
- $\int f(x) dx = \frac{3h}{8} (y_0 + y_n) + 2 (y_3 + y_6 + y_9 + \dots) + 3 (y_1 + y_2 + y_4 + \dots)$
- $\int f(x) dx = \frac{3h}{8} (\text{sum of the first and the last term}) + 2 (\text{multiples of three}) + 3 (\text{sum of the remaining terms})$

INTRODUCTION

- There exists large number of ordinary differential equations, whose solution cannot be obtained by the known analytical methods. In such cases, we use numerical methods to get an approximate solution of a given differential equation with given initial condition.

NUMERICAL DIFFERENTIATION

- Consider an ordinary differential equation of first order and first degree of the form
- $dy/dx = f(x, y) \dots\dots\dots(1)$
- with the initial condition $y(x_0) = y_0$ which is called initial value problem.
- To find the solution of the initial value problem of the form (1) by numerical methods, we divide the interval (a, b) on which the solution is derived in finite number of sub-intervals by the points

TAYLOR'S SERIES METHOD

- Consider the first order differential equation
- $$dy/dx = f(x, y) \dots\dots\dots(1)$$

with initial conditions $y(x_0) = y_0$ then expanding $y(x)$ i.e $f(x)$ in a Taylor's series at the point x_0 we get

$$y(x_0 + h) = y(x_0) + hy'(x_0) + h^2/2! Y''(x_0) + \dots\dots\dots$$

Note : Taylor's series method can be applied only when the various derivatives of $f(x,y)$ exist and the value of $f(x-x_0)$ in the expansion of $y = f(x)$ near x_0 must be very small so that the series is convergent.

PICARDS METHOD OF SUCCESSIVE APPROXIMATION

- Consider the differential equation
- $dy/dx = f(x,y)$ with initial conditions
- $y(x_0) = y_0$ then
- The first approximation y_1 is obtained by
- $y_1 = y_0 + f(x, y_0) dx$
- The second approximation y_2 is obtained by $y_2 = y_1 + f(x, y_1) dx$

PICARDS APPROXIMATION METHOD

- The third approximation of y_3 is obtained by y_2 is given by
- $y_3 = y_0 + f(x, y_2) dx \dots \dots \dots$ and so on
- $\dots \dots \dots$
- $\dots \dots \dots$
- $y_n = y_0 + f(x, y^{n-1}) dx$
- The process of iteration is stopped when any two values of iteration are approximately the same.
- Note : When x is large , the convergence is slow.

EULER'S METHOD

- Consider the differential equation
- $dy/dx = f(x,y).....(1)$
- With the initial conditions $y(x_0) = y_0$
- The first approximation of y_1 is given by
- $y_1 = y_0 + h f(x_0, y_0)$
- The second approximation of y_2 is given by
- $y_2 = y_1 + h f(x_0 + h, y_1)$

EULER'S METHOD

- The third approximation of y_3 is given by
- $y_3 = y_2 + h f (x_0 + 2h, y_2)$
-
-
-
- $y_n = y_{n-1} + h f [x_0 + (n-1)h, y_{n-1}]$
- This is Eulers method to find an approximate solution of the given differential equation.

IMPORTANT NOTE

- Note : In Euler's method, we approximate the curve of solution by the tangent in each interval i.e by a sequence of short lines. Unless h is small there will be large error in y_n . The sequence of lines may also deviate considerably from the curve of solution. The process is very slow and the value of h must be smaller to obtain accuracy reasonably.

MODIFIED EULER'S METHOD

- By using Euler's method, first we have to find the value of $y_1 = y_o + hf(x_o, y_o)$
- WORKING RULE
- Modified Euler's formula is given by
- $y_{k+1}^i = y_k + h/2 [f(x_k, y_k) + f(x_{k+1}, y_{k+1})$
- when $i=1, y(o)_{k+1}$ can be calculated
- from Euler's method.

MODIFIED EULER'S METHOD

- When $k=0,1,2,3,\dots$ gives number of iterations
- $i = 1,2,3,\dots$ gives number of times a particular iteration k is repeated when
- $i=1$
- $Y^1_{k+1} = y_k + h/2 [f(x_k, y_k) + f(x_{k+1}, y_{k+1})], \dots$

RUNGE-KUTTA METHOD

The basic advantage of using the Taylor series method lies in the calculation of higher order total derivatives of y . Euler's method requires the smallness of h for attaining reasonable accuracy. In order to overcome these disadvantages, the Runge-Kutta methods are designed to obtain greater accuracy and at the same time to avoid the need for calculating higher order derivatives. The advantage of these methods is that the functional values only required at some selected points on the subinterval.

R-K METHOD

- Consider the differential equation
- $dy/dx = f(x, y)$
- With the initial conditions $y(x_0) = y_0$
- First order R-K method :
- $y_1 = y(x_0 + h)$
- Expanding by Taylor's series
- $y_1 = y_0 + h y'_0 + \frac{h^2}{2} y''_0 + \dots$

R-K METHOD

- Also by Euler's method
- $y_1 = y_o + h f(x_o, y_o)$
- $= y_o + h y_o'$
- It follows that the Euler's method agrees with the Taylor's series solution upto the term in h . Hence, Euler's method is the Runge-Kutta method of the first order.

- The second order R-K method is given by

- $y_1 = y_0 + \frac{1}{2} (k_1 + k_2)$

- where $k_1 = h f(x_0, y_0)$

- $k_2 = h f(x_0 + h, y_0 + k_1)$

- Third order R-K method is given by

- $y_1 = y_o + 1/6 (k_1 + k_2 + k_3)$
- where $k_1 = h f (x_o , y_o)$
- $k_2 = hf(x_o + 1/2h , y_o + 1/2 k_1)$
- $k_3 = hf(x_o + 1/2h , y_o + 1/2 k_2)$

The Fourth order R – K method :

This method is most commonly used and is often referred to as Runge – Kutta method only. Proceeding as mentioned above with local discretisation error in this method being $O(h^5)$, the increment K of y corresponding to an increment h of x by Runge – Kutta method from

$dy/dx = f(x,y), y(x_0)=Y_0$ is given by

- $K_1 = h f (x_o , y_o)$
- $k_2 = h f (x_o + \frac{1}{2} h, y_o + \frac{1}{2} k_1)$
- $k_3 = h f (x_o + \frac{1}{2} h, y_o + \frac{1}{2} k_2)$
- $k_4 = h f (x_o + \frac{1}{2} h, y_o + \frac{1}{2} k_3)$
- and finally computing
- $k = \frac{1}{6} (k_1 + 2 k_2 + 2 k_3 + k_4)$
- Which gives the required approximate value as
- $y_1 = y_o + k$

R-K METHOD

Note 1 : k is known as weighted mean of k_1, k_2, k_3 and k_4 .

Note 2 : The advantage of these methods is that the operation is identical whether the differential equation is linear or non-linear.

FINITE DIFFERENCE METHOD

$$y'_i = \frac{1}{2h}(y_{i+1} - y_{i-1})$$

$$y''_i = \frac{1}{h^2}(y_{i+1} - 2y_i + y_{i-1})$$

Similarly, $y'''_i = \frac{1}{2h^3}(y_{i+2} - 2y_{i+1} + 2y_{i-1} - y_{i-2})$

$$y^{iv}_i = \frac{1}{h^4}(y_{i+2} - 4y_{i+1} + 6y_i - 4y_{i-1} + y_{i-2})$$

Characteristic matrix: Let A be a square matrix of order n then the matrix $(A-\lambda I)$ is called Characteristic matrix of A. where I is the unit matrix of order n and λ is any scalar

Ex: Let $A = \begin{pmatrix} 1 & 2 \\ 3 & 4 \end{pmatrix}$ then

$A - \lambda I = \begin{pmatrix} 1-\lambda & 2 \\ 3 & 4-\lambda \end{pmatrix}$ is the characteristic matrix of A

Characteristic polynomial: Let A be square matrix of order n then $|A - \lambda I|$ is called characteristic polynomial of A

Ex: Let $A = \begin{pmatrix} 1 & -2 \\ 4 & 1 \end{pmatrix}$

then $|A - \lambda I| = \lambda^2 - 2\lambda + 9$ is the characteristic polynomial of A

Characteristic equation: Let A be square matrix of order n then $|A - \lambda I| = 0$ is called characteristic equation of A

Ex: Let $A = \begin{pmatrix} 1 & -2 \\ 4 & 1 \end{pmatrix}$

then $|A - \lambda I| = \lambda^2 - 2\lambda + 9 = 0$ is the characteristic equation of A

Eigen values: The roots of the characteristic equation $|A - \lambda I| = 0$ are called the eigen values

Ex: Let $A = \begin{pmatrix} 2 & 1 \\ 3 & 4 \end{pmatrix}$ then $|A - \lambda I| = \lambda^2 - 6\lambda + 5 = 0$

Therefore $\lambda = 1, 5$ are the eigen values of the matrix A

- Eigen vector: If λ is an eigen value of the square matrix A . If there exists a non- zero vector X such that $AX=\lambda X$ is said to be eigen vector corresponding to eigen value λ of a square matrix A
- Eigen vector must be a non-zero vector
- If λ is an eigen value of matrix A if and only if there exists a non-zero vector X such that $AX=\lambda X$
- If X is an eigen vector of a matrix A corresponding to the eigen value λ , then kX is also an eigen vector of A corresponding to the same eigen value λ . K is a non zero scalar.

Properties of eigen values and eigen vectors

- The matrices A and A^T have the same eigen values.
- If $\lambda_1, \lambda_2, \dots, \lambda_n$ are the eigen values of A then $1/\lambda_1, 1/\lambda_2, \dots, 1/\lambda_n$ are the eigen values of A^{-1} .
- If $\lambda_1, \lambda_2, \dots, \lambda_n$ are the eigen values of A then $\lambda_1^k, \lambda_2^k, \dots, \lambda_n^k$ are the eigen values of A^k .
- If λ is the eigen value of a non singular matrix A , then $|A|/\lambda$ is the eigen value of A .

- The sum of the eigen values of a matrix is the trace of the matrix

- If λ is the eigen value of A then the eigen values of $B = a_0 A^2 + a_1 A + a_2 I$ is $a_0 \lambda^2 + a_1 \lambda + a_2$.

Similar Matrices : Two matrices A & B are said to be similar if there exists an invertible matrix P such that $B = P^{-1}AP$.

➤ Eigen values of two similar matrices are same

➤ If A & B are square matrices and if A is invertible then the matrices $A^{-1}B$ & BA^{-1} have the same eigen values

Power Method

Power method for finding eigenvalues

- 1. Start with an initial guess for x**
- 2. Calculate $w = Ax$**
- 3. Largest value (magnitude) in w is the estimate of eigenvalue**
- 4. Get next x by rescaling w (to avoid the computation of very large matrix A^n)**
- 5. Continue until converged**

Power Method

- Start with initial guess $\mathbf{z} = \mathbf{x}_0$

$$\begin{aligned} \mathbf{w}^{(1)} &= A\mathbf{z}^{(1)} && \Rightarrow \lambda_k^{(1)} = w_k^{(1)} \\ \mathbf{z}^{(2)} &= \frac{\mathbf{w}^{(1)}}{\lambda_k^{(1)}} = \frac{A\mathbf{z}^{(1)}}{\lambda_k^{(1)}} && \leftarrow \text{rescaling} \\ \mathbf{w}^{(2)} &= A\mathbf{z}^{(2)} && \Rightarrow \lambda_k^{(2)} = w_k^{(2)} \\ \mathbf{z}^{(3)} &= \frac{\mathbf{w}^{(2)}}{\lambda_k^{(2)}} = \frac{A\mathbf{z}^{(2)}}{\lambda_k^{(2)}} && \leftarrow \lambda_k \text{ is the dominant eigenvalue} \\ &\vdots \end{aligned}$$

If $|\lambda_1| > |\lambda_2| > |\lambda_3| > \dots > |\lambda_n|$, then

$$|\lambda_1|^k \gg |\lambda_2|^k \gg |\lambda_3|^k \gg \dots \gg |\lambda_n|^k$$

Power Method

1. Initial guess $z^{(1)} = x_0 = [1, 1, \dots, 1]^T$
2. Calculate $Az^{(1)} = w^{(1)} = \lambda z^{(2)}$; (*normalize* z by biggest w_k)
3. Calculate $Az^{(2)} = w^{(2)} = \lambda z^{(3)}$; (*normalize* z by biggest w_k)
- \vdots
- ... Calculate $Az^{(k)} = w^{(k)} = \lambda z^{(k+1)}$

- For large number of iterations, λ should converge to the largest eigenvalue
- The normalization make the right hand side converge to λ , rather than λ^n

THANK YOU

