
APPENDIX**A**

Solution of Linear Systems by Matrix Methods

A.1 OVERVIEW OF MATRICES

A *matrix* is a rectangular array of *elements* arranged in rows and columns. If the matrix contains m rows and n columns, the matrix is said to be of size $m \times n$. The element in the i th row and j th column of the matrix A is denoted by a_{ij} . For example,

$$A = \begin{bmatrix} a_{11} & a_{12} & a_{13} \\ a_{21} & a_{22} & a_{23} \\ a_{31} & a_{32} & a_{33} \\ a_{41} & a_{42} & a_{43} \end{bmatrix}$$

is a 4×3 matrix.

Types of Matrices

Matrices may belong to one of the many types discussed here.

1. **Square Matrix** A matrix in which rows m is equal to columns n .

2. **Identity Matrix** A square matrix in which all the diagonal elements are one and other elements are zero. That is

$$\begin{aligned} a_{ij} &= 1 & \text{for} & \quad i = j \\ a_{ij} &= 0 & \text{for} & \quad i \neq j \end{aligned}$$

Identity matrices are denoted by I . For example, a 3×3 identity is written as

$$I = \begin{bmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{bmatrix}$$

3. Row Vector A matrix with one row and n columns.

4. Column Vector A matrix with m rows and one column.

5. Transpose Matrix The matrix A^T is called the transpose of A if the element a_{ij} in A is equal to element a_{ji} in A^T for all i and j . For example, if

$$A = \begin{bmatrix} 1 & 2 \\ 3 & 4 \\ 5 & 6 \end{bmatrix}$$

then,

$$A^T = \begin{bmatrix} 1 & 3 & 5 \\ 2 & 4 & 6 \end{bmatrix}$$

It is clear that A^T is obtained by interchanging the rows and the columns of A .

6. Zero Matrix A matrix in which every element is zero.

7. Equal Matrix A matrix is said to be equal to another matrix if and only if they have the same order and the corresponding elements are equal. That is,

$$\text{matrix } A = \text{matrix } B$$

if $a_{ij} = b_{ij}$ for all i and j .

Matrix Algebra

We can perform only three operations, namely, addition, subtraction and multiplication on the matrices. The division, although not defined, is replaced by the concept of inversion discussed later.

Two matrices A and B can be added together (or subtracted from each other) if they are of the same order. Then

$$\text{sum } C = A + B$$

can be obtained by adding the corresponding elements. That is,

$$c_{ij} = a_{ij} + b_{ij} \quad \text{for all } i \text{ and } j$$

Similarly,

$$\text{difference } E = A - B$$

can be obtained by subtracting the corresponding elements. That is,

$$e_{ij} = a_{ij} - b_{ij} \quad \text{for all } i \text{ and } j$$

Two matrices A and B can be multiplied in the order AB if and only if the number of columns of A is equal to the number of rows of B . That is, if A is of order $m \times r$, then B should be of order $r \times n$, where m and n are arbitrary values. In such cases, we may obtain

$$P = AB$$

which is of order $m \times n$. Elements of the matrix P is given by

$$p_{ij} = \sum_{k=1}^r a_{ik} b_{kj} \quad \text{for all } i \text{ and } j$$

The following general properties apply to matrix algebra.

1. $A \pm B = B \pm A$
2. $A \pm (B \pm C) = (A \pm B) \pm C$
3. $(A \pm B)^T = A^T \pm B^T$
4. $IA = AI$, where I is identity matrix
5. $(AB)C = A(BC)$
6. $C(A + B) = CA + CB$
7. $(A + B)C = AC + BC$
8. $\alpha(AB) = (\alpha A)B = A(\alpha B)$, where α is a scalar

Traces and Determinants

The *trace* of a matrix is the number obtained by adding its diagonal elements (from upper left corner to the lower right corner). For example, the trace of the matrix

$$A = \begin{bmatrix} a_{11} & a_{12} & a_{13} \\ a_{21} & a_{22} & a_{23} \\ a_{31} & a_{32} & a_{33} \end{bmatrix} \text{ is } a_{11} + a_{22} + a_{33}$$

The *determinant* of a 2×2 matrix, say A , is written in the form

$$\begin{vmatrix} a_{11} & a_{12} \\ a_{21} & a_{22} \end{vmatrix}$$

and the determinant is computed as

$$a_{11} a_{22} - a_{12} a_{21}$$

Similarly, for a 3×3 matrix the determinant is given by

$$\begin{aligned} |A| &= \begin{vmatrix} a_{11} & a_{12} & a_{13} \\ a_{21} & a_{22} & a_{23} \\ a_{31} & a_{32} & a_{33} \end{vmatrix} \\ &= a_{11} \begin{vmatrix} a_{22} & a_{23} \\ a_{32} & a_{33} \end{vmatrix} - a_{12} \begin{vmatrix} a_{21} & a_{23} \\ a_{31} & a_{33} \end{vmatrix} + a_{13} \begin{vmatrix} a_{21} & a_{22} \\ a_{31} & a_{32} \end{vmatrix} \end{aligned}$$

$$\begin{aligned}
 &= a_{11}(a_{22}a_{33} - a_{23}a_{32}) - a_{12}(a_{21}a_{33} - a_{31}a_{23}) \\
 &\quad + a_{13}(a_{21}a_{32} - a_{22}a_{31}) \\
 &= a_{11}a_{22}a_{33} + a_{12}(-a_{21}a_{33} + a_{31}a_{23}) \\
 &\quad - a_{21}a_{12}a_{33} - a_{31}a_{22}a_{13}
 \end{aligned}$$

Note that there are 6 product terms added together. For larger matrices, the determinant is much more difficult to define and compute manually. In general, for an $n \times n$ matrix, the determinant will contain a sum of $n!$ signed product terms, each having n elements.

Some of the important properties of determinants that would be helpful in computing their values are:

1. Interchanging two rows (or two columns) of a matrix changes the sign of the determinant but not the value.
2. If two rows (or columns) of a matrix are identical then its determinant is zero.
3. If a matrix contains a row (or column) of all zeros, then its determinant is zero.
4. Value of the determinant of a matrix does not change when a scalar multiple of one row (or column) is added to another row (or column).
5. If every element of a row (or column) is multiplied by a scalar α , the value of the determinant is multiplied by α .
6. If A and B are square matrices of same size, then $|AB| = |A||B|$.
7. For a triangular matrix (in which all the elements below (or above) the diagonal are zero), the determinant is the product of the diagonal elements.

Minors and Cofactors

The minor M_{ij} of the element a_{ij} of the determinant $|A|$ is obtained by striking out the i th row and j th column. That is,

$$M_{11} = \begin{vmatrix} a_{22} & a_{23} \\ a_{32} & a_{33} \end{vmatrix}, \quad M_{22} = \begin{vmatrix} a_{11} & a_{13} \\ a_{31} & a_{33} \end{vmatrix}$$

and so on. That is, the minor of a particular element is the determinant that remains after the row and the column that contain the element have been deleted.

The cofactor of an element is its minor with a sign attached. The cofactor d_{ij} of an element a_{ij} is given by

$$d_{ij} = (-1)^{i+j} M_{ij}$$

The value of the determinant of a matrix can be obtained by expanding the determinant by cofactors. This is done by choosing any column or row and determining the sum of the product of each element in the chosen row or column and its cofactor.

Adjoint Matrix

If d_{ij} is the cofactor of the element a_{ij} of the square matrix A , then, by definition, the *adjoint matrix* of A is given by

$$\text{adj}(A) = D^T$$

where

$$D = \begin{bmatrix} d_{11} & d_{12} & \dots & d_{1n} \\ d_{21} & d_{22} & \dots & d_{2n} \\ \vdots & \vdots & & \vdots \\ d_{n1} & d_{n2} & \dots & d_{nn} \end{bmatrix}$$

Inverse of a Matrix

If B and C are two $n \times n$ square matrices such that

$$BC = CB = I \text{ (Identity matrix)}$$

then, B is called the *inverse* of C and C the inverse of B . The common notation for inverses is B^{-1} and C^{-1} . That is,

$$B^{-1} B = I$$

$$C^{-1} C = I$$

The concept of matrix inversion is useful in solving linear systems of equations.

A.2 SOLUTION OF LINEAR SYSTEMS BY DETERMINANTS

We can solve a system of linear equations by determinants using a method called *Cramer's Rule*. For the sake of simplicity, we consider a 2×2 system such as

$$ax_1 + bx_2 = c$$

$$dx_1 + ex_2 = f$$

We can solve for the variable x_1 by eliminating the variable x_2 . Thus,

$$x_1 = \frac{ce - bf}{ae - bd}$$

Similarly,

$$x_2 = \frac{af - cd}{ae - bd}$$

Alternatively, we can express x_1 and x_2 using determinants. That is,

$$x_i = \frac{\begin{vmatrix} c & b \\ f & e \end{vmatrix}}{\begin{vmatrix} a & b \\ d & e \end{vmatrix}} \quad \text{and} \quad x_2 = \frac{\begin{vmatrix} a & c \\ d & e \end{vmatrix}}{\begin{vmatrix} a & b \\ d & e \end{vmatrix}}$$

This is known as *Cramer's rule*.

Example A.1

Solve the following system of equations using Cramer's rule:

$$2x_1 + 3x_2 = 12$$

$$4x_1 - x_2 = 10$$

$$x_1 = \frac{\begin{vmatrix} 12 & 3 \\ 10 & -1 \end{vmatrix}}{\begin{vmatrix} 2 & 3 \\ 4 & -1 \end{vmatrix}} = \frac{-12 - 30}{-2 - 12} = \frac{42}{14} = 3$$

$$x_2 = \frac{\begin{vmatrix} 2 & 12 \\ 4 & 10 \end{vmatrix}}{\begin{vmatrix} 2 & 3 \\ 4 & -1 \end{vmatrix}} = \frac{20 - 48}{-14} = \frac{28}{14} = 2$$

A.4 SOLUTION OF LINEAR SYSTEMS BY MATRIX INVERSION

A linear system of n equations in n unknowns can be represented in matrix form as

$$AX = B$$

where A , X and B are matrices and are given by

$$A = \begin{bmatrix} a_{11} & a_{12} & \dots & a_{1n} \\ a_{21} & a_{22} & \dots & a_{2n} \\ \vdots & \vdots & & \vdots \\ a_{n1} & a_{n2} & \dots & a_{nn} \end{bmatrix}$$

$$X = \begin{bmatrix} x_1 \\ x_2 \\ \vdots \\ x_n \end{bmatrix} \quad B = \begin{bmatrix} b_1 \\ b_2 \\ \vdots \\ b_n \end{bmatrix}$$

A is called *coefficient matrix* and B is known as *constant vector*. X is the required solution and, therefore, it is called the *solution vector*. If we multiply the matrix equation

$$AX = B$$

By A^{-1} on both sides, we get

$$A^{-1}AX = A^{-1}B$$

A^{-1} is the *inverse matrix* of A . We know that $A^{-1}A = I$ is the *identity matrix* and is given by

$$I = \begin{bmatrix} 1 & 0 & 0 & \dots & 0 \\ 0 & 1 & 0 & \dots & 0 \\ 0 & 0 & 1 & \dots & 0 \\ \vdots & & & & \\ 0 & 0 & 0 & \dots & 1 \end{bmatrix}$$

Therefore,

$$A^{-1}AX = IX = X$$

is the solution of the system of equations and is obtained from

$$X = A^{-1}B = CB$$

If we know the inverse of the matrix A , we can obtain the solution vector X by post-multiplying it by B .

The inverse A^{-1} of a square matrix A exists, if and only if, A is *nonsingular* (i.e. $\det A \neq 0$). A^{-1} is the matrix obtained from A by replacing each element a_{ij} by its *cofactor* d_{ij} and then transposing the resulting matrix and dividing it by the *determinant* of A .

$$C = A^{-1} = \frac{\text{adj}(A)}{\det A}$$

where $\text{adj}(A)$ is the *adjoint* of matrix A and is given by the transpose of the *cofactor* matrix of A .

Then

$$\text{adj}(A) = D^T$$

where

$$D = \begin{bmatrix} d_{11} & d_{12} & \dots & d_{1n} \\ d_{21} & d_{22} & \dots & d_{2n} \\ \vdots & \vdots & & \\ d_{n1} & d_{n2} & \dots & d_{nn} \end{bmatrix}$$

d_{ij} is the cofactor of the element a_{ij} and is given by

$$d_{ij} = (-1)^{i+j} M_{ij}$$

M_{ij} is called the minor of a_{ij} and is taken as the determinant of matrix A after deleting i th row and j th column.

Solve the following system using matrix inversion method

$$2x_1 + x_2 + x_3 = 7$$

$$x_1 - x_2 + x_3 = 0$$

$$4x_1 + 2x_2 - 3x_3 = 4$$

Given, $A = \begin{bmatrix} 2 & 1 & 1 \\ 1 & -1 & 1 \\ 4 & 2 & -3 \end{bmatrix}$ and $B = \begin{bmatrix} 7 \\ 0 \\ 4 \end{bmatrix}$

$$M_{11} = \begin{vmatrix} -1 & 1 \\ 2 & -3 \end{vmatrix} = 1 \quad d_{11} = 1$$

$$M_{12} = \begin{vmatrix} 1 & 1 \\ 4 & -3 \end{vmatrix} = -7 \quad d_{12} = 7$$

$$M_{13} = \begin{vmatrix} 1 & -1 \\ 4 & -2 \end{vmatrix} = 6 \quad d_{13} = 6$$

$$M_{21} = \begin{vmatrix} 1 & 1 \\ 2 & -3 \end{vmatrix} = -5 \quad d_{21} = 5$$

$$M_{22} = \begin{vmatrix} 2 & 1 \\ 4 & -3 \end{vmatrix} = -10 \quad d_{22} = -10$$

$$M_{23} = \begin{vmatrix} 2 & 1 \\ 4 & 2 \end{vmatrix} = 0 \quad d_{23} = 0$$

$$M_{31} = \begin{vmatrix} 1 & 1 \\ -1 & 1 \end{vmatrix} = 2 \quad d_{31} = 2$$

$$M_{32} = \begin{vmatrix} 2 & 1 \\ 1 & 1 \end{vmatrix} = 1 \quad d_{32} = -1$$

$$M_{33} = \begin{vmatrix} 2 & 1 \\ 1 & -1 \end{vmatrix} = -3 \quad d_{33} = -3$$

$$D = \begin{bmatrix} 1 & 7 & 6 \\ 5 & -10 & 0 \\ 2 & -1 & -3 \end{bmatrix}$$

$$\text{Adj}A = D^T = \begin{bmatrix} 1 & 5 & 2 \\ 7 & -10 & -1 \\ 6 & 0 & -3 \end{bmatrix}$$

$$\det A = a_{11}d_{11} + a_{12}d_{12} + a_{13}d_{13} = 15$$

$$A^{-1} = C = \begin{bmatrix} \frac{1}{15} & \frac{5}{15} & \frac{2}{15} \\ \frac{7}{15} & \frac{-10}{15} & \frac{-1}{15} \\ \frac{6}{15} & 0 & \frac{-3}{15} \end{bmatrix}$$

$$B = \begin{bmatrix} 7 \\ 0 \\ 4 \end{bmatrix}$$

We know that $X = CB$ and therefore

$$x_1 = \frac{7}{15} + 0 + \frac{8}{15} = 1$$

$$x_2 = \frac{49}{15} - 0 - \frac{4}{15} = 3$$

$$x_3 = \frac{42}{15} + 0 - \frac{12}{15} = 2$$

A.4 GAUSS-JORDAN MATRIX INVERSION

The Gauss-Jordan elimination technique (discussed in Chapter 7) can be used to invert a matrix effectively. The square matrix A which is to be inverted is augmented by a unit matrix of the same order as follows.

$$\left[\begin{array}{cccc|ccc} a_{11} & a_{12} & \dots & a_{1n} & 1 & 0 & \dots & 0 \\ a_{21} & a_{22} & \dots & a_{2n} & 0 & 1 & \dots & 0 \\ \vdots & \vdots & & \vdots & \vdots & & & \\ a_{n1} & a_{n2} & \dots & a_{nn} & 0 & 0 & \dots & 1 \end{array} \right]$$

If we carry out Gauss-Jordan elimination using the first row as a pivot row we get

$$\left[\begin{array}{cccc|cccc} 1 & a_{12} & \dots & a_{1n} & a_{1,n+1} & 0 & \dots & 0 \\ 0 & a_{21} & \dots & a_{2n} & a_{2,n+1} & 1 & \dots & 0 \\ \vdots & \vdots & & \vdots & \vdots & & & \\ 0 & a_{n2} & \dots & a_{nn} & a_{n,n+1} & 0 & \dots & 1 \end{array} \right]$$

When we repeat the process using the second row as the pivot row, the result is

$$\left[\begin{array}{ccccc|ccccc} 1 & 0 & a_{13} & \dots & a_{1n} & a_{1,n+1} & a_{1,n+2} & 0 & \dots & 0 \\ 0 & 1 & a_{23} & \dots & a_{2n} & a_{2,n+1} & a_{2,n+2} & 0 & \dots & 0 \\ 0 & 0 & a_{33} & \dots & a_{3n} & a_{3,n+1} & a_{3,n+2} & 1 & \dots & 0 \\ 0 & 0 & a_{n3} & \dots & a_{nn} & a_{n,n+1} & a_{n,n+2} & 0 & \dots & 1 \end{array} \right]$$

This elimination process, if continued for all the n rows, yields the final result as follows:

$$\left[\begin{array}{ccccc|ccccc} 1 & 0 & 0 & \dots & 0 & a_{1,n+1} & a_{1,n+2} & \dots & a_{1,n+n} \\ 0 & 1 & 0 & \dots & 0 & a_{2,n+1} & a_{2,n+2} & \dots & a_{2,n+n} \\ 0 & 0 & 1 & \dots & 0 & \vdots & & & \\ \vdots & & & & & & & & \\ 0 & 0 & 0 & \dots & 1 & a_{n,n+1} & a_{n,n+2} & \dots & a_{n,n+n} \end{array} \right]$$

The matrix a_{ij} for $i = 1$ to n and $j = n + 1$ to $2n$ is the inverse matrix of A .

$$C = A^{-1} = [a_{ij}] \quad \begin{cases} i = 1, \dots, n \\ j = n + 1, \dots, 2n \end{cases}$$

where the element

$$c_{ij} = a_{i,n+j} \text{ for } i = 1, \dots, n \quad \text{and} \quad j = 1, \dots, n.$$

Example A-3

Find the inverse of the coefficient matrix of the system given in Example A2 using Gauss-Jordan elimination technique.

$$\text{Augmented } A = \left[\begin{array}{ccc|ccc} 2 & 1 & 1 & 1 & 0 & 0 \\ 1 & -1 & 1 & 0 & 1 & 0 \\ 4 & 2 & -3 & 0 & 0 & 1 \end{array} \right]$$

Pivot row-1

$$\left[\begin{array}{ccc|ccc} 1 & 1/2 & 1/2 & 1/2 & 0 & 0 \\ 0 & -3/2 & 1/2 & -1/2 & 1 & 0 \\ 0 & 0 & -5 & -2 & 0 & 1 \end{array} \right]$$

Pivot row-2

$$\left[\begin{array}{ccc|ccc} 1 & 0 & 2/3 & 1/3 & 1/3 & 0 \\ 0 & 1 & -1/3 & 1/3 & -2/3 & 0 \\ 0 & 0 & -5 & -2 & 0 & 1 \end{array} \right]$$

Pivot row-3

$$\left[\begin{array}{ccc|ccc} 1 & 0 & 0 & 1/15 & 1/3 & 2/15 \\ 0 & 1 & 0 & 7/15 & -2/3 & -1/15 \\ 0 & 0 & 1 & 2/5 & 0 & -1/5 \end{array} \right]$$

The last three columns represent the inverse of the matrix

$$\begin{matrix} 2 & 1 & 1 \\ 1 & -1 & 1 \\ 4 & 2 & -3 \end{matrix}$$

Compare this with the result obtained in the previous section.

Once the inverse of A is known, the solution vector X can be obtained by simple matrix multiplication. That is

$$X = A^{-1}B = CB$$

$$x_1 = c_{11}b_1 + c_{12}b_2 + \dots + c_{1n}b_n$$

$$x_2 = c_{21}b_1 + c_{22}b_2 + \dots + c_{2n}b_n$$

and so on.

In general,

$$x_i = \sum_{j=1}^n c_{ij} b_j \quad i = 1, \dots, n$$

Note that $c_{ij} = a_{i, n+j}$ of augmented A in the final step.

Key Terms

- Adjoint matrix*
- Coefficient matrix*
- Cofactor*
- Column vector*
- Cramer's rule*
- Determinant*
- Equal matrix*
- Identity matrix*

- Inverse matrix*
- Minor*
- Nonsingular matrix*
- Row vector*
- Square matrix*
- Trace of a matrix*
- Transpose matrix*
- Zero matrix*

Solution of Polynomials by Graeffe's Root Squaring Method

Graeffe's root squaring method is a direct method of finding roots of a polynomial with real coefficients. This method deserves attention both for its historical interest and a novel idea involved.

Graeffe's method transforms a polynomial $p_n(x)$ into another polynomial of the same degree but whose roots are the squares of the roots of the original polynomial. Because of the squaring property, the roots of the new polynomial will be spread apart more widely than in the original one, when the roots are greater than 1 in absolute value. Repeating this process until the roots are really far apart, we can compute the roots directly from the coefficients.

We consider here a simple example to illustrate the root squaring technique.

Let

$$p_0(x) = (x - 1)(x - 2)(x - 3) \quad (B.1)$$

Then, we consider another function $p_1(y)$ such that

$$\begin{aligned} p_1(y) &= -p_0(x) p_0(-x) \\ &= -(x - 1)(x - 2)(x - 3)(-x - 1)(-x - 2)(-x - 3) \\ &= (x - 1)(x - 2)(x - 3)(x + 1)(x + 2)(x + 3) \\ &= (x^2 - 1)(x^2 - 4)(x^2 - 9) \\ &= (y - 1)(y - 4)(y - 9) \end{aligned} \quad (B.2)$$

where

$$y = x^2$$

We know that the roots of $p_0(x) = 0$ are $x_1 = 1$, $x_2 = 2$, and $x_3 = 3$. And from Eq. (B.2) the roots of $p_1(y) = 0$ are $y_1 = 1$, $y_2 = 4$, and $y_3 = 9$. It shows that roots of $p_1(y) = 0$ are the squares of the roots of $p_0(y) = 0$. This implies that if we compute the roots of $p_1(y)$, then we can obtain the roots of $p_0(x)$ from

$$x_1 = \sqrt{y_1}$$

$$x_2 = \sqrt{y_2}$$

$$x_3 = \sqrt{y_3}$$

Now, let us repeat the procedure for finding the roots of $p_1(y)$. Consider a third polynomial

$$\begin{aligned} p_2(z) &= -p_1(y) p_1(-y) \\ &= (z - 1)(z - 16)(z - 81) \end{aligned} \quad (\text{B.3})$$

The roots of Eq. (B.3) are

$$\begin{array}{ccccccc} z_1 = 1 (= y_1^2) & \longleftarrow & y_1 (= x_1^2) & \longleftarrow & x_1 \\ z_2 = 16 (= y_2^2) & \longleftarrow & y_2 (= x_2^2) & \longleftarrow & x_2 \\ z_3 = 81 (= y_3^2) & \longleftarrow & y_3 (= x_3^2) & \longleftarrow & x_3 \\ \text{Iteration 2} & & & & \text{Iteration 1} \end{array}$$

That is, after the second iteration, we can estimate the roots of original equation $p_0(x)$ from the relation

$$z_i = (x_i^2)^2 = x_i^4, \quad i = 1, 2, 3$$

Suppose we have done the squaring process k times and the roots of the final equation are r_i , then

$$r_i = x_i^{(2^k)} = (x_i)^{2^k} \quad (\text{B.4})$$

Remember that we never have $p_0(x)$ in factored form as given in Eq. (B.1), but the result is the same.

Now, let us consider a third degree polynomial in standard form as

$$p_0(x) = a_3x^3 + a_2x^2 + a_1x + a_0 \quad (\text{B.5})$$

Then,

$$\begin{aligned} p_1(y) &= -p_0(-x) p(x) \\ &= (a_3x^3 + a_2x^2 + a_1x + a_0)(a_3x^3 - a_2x^2 + a_1x - a_0) \\ &= a_3^2x^6 - (a_2^2 - 2a_1a_3)x^4 + (a_1^2 - 2a_0a_2)x^2 - a_0^2 \\ &= b_3y^3 + b_2y^2 + b_1y_1 + b_0 \end{aligned} \quad (\text{B.6})$$

where

$$y = x^2$$

and

$$\boxed{\begin{aligned} b_3 &= + a_3^2 \\ b_2 &= -(a_2^2 - 2a_1a_3) \\ b_1 &= + (a_1^2 - 2a_0a_2) \\ b_0 &= -a_0^2 \end{aligned}} \quad (\text{B.7})$$

We can thus show that for a general polynomial of degree n ,

$$p_0(x) = a_nx^n + a_{n-1}x^{n-1} + \dots + a_0 \quad (\text{B.8})$$

after first squaring process,

$$\boxed{\begin{aligned} b_n &= + a_n^2 \\ b_{n-1} &= -(a_{n-1}^2 - 2a_n a_{n-2}) \\ b_{n-2} &= + (a_{n-2}^2 - 2a_{n-1} a_{n-3} + 2a_n a_{n-4}) \\ &\vdots \\ b_0 &= (-1)^n a_0^2 \end{aligned}} \quad (\text{B.9})$$

This process can be repeated replacing 'a' values by 'b' values in Eq. (B.9) each time. Let us suppose that our final equation after k iterations (i.e. squaring k times) is

$$B_n y^n + B_{n-1} y^{n-1} + \dots + B_0 = 0 \quad (\text{B.10})$$

Assuming that the roots of Eq. (B.10) are now more widely separated, we have

$$|y_1| >> |y_2| >> |y_3| \dots >> y_n$$

Then,

$$\begin{aligned} y_1 &\approx -\frac{B_{n-1}}{B_n} \\ y_2 &\approx -\frac{B_{n-2}}{B_{n-1}} \\ &\vdots \\ &\vdots \\ y_n &\approx -\frac{B_0}{B_1} \end{aligned}$$

That is,

$$\begin{aligned} y_i &\approx -\frac{B_{n-i}}{B_{n-i+1}}, \quad i = 1, 2, \dots, n \\ &= (x_i)^{2^k} \end{aligned} \quad (\text{B.11})$$

Thus,

$$x_i = 2^k \text{th root of } y_i = 2^k \sqrt[k]{\left| \frac{B_{n-i}}{B_{n-i+1}} \right|} \quad (\text{B.12})$$

The main advantage of Graeffe's root squaring method over other methods is that it does not require any initial guessing about roots. The method is also capable of giving all the roots but the limitation is that the polynomial should have only real coefficients.

Example 8.1

Apply root squaring technique to estimate the roots of

$$x^3 - 3x^2 - 6x + 8 = 0$$

Table below shows the coefficients of successive polynomials (using Eq. (B.7)) as well as the roots estimated (using Eq. (B.12)).

k	Coefficients				Roots estimated		
	a_3	a_2	a_1	a_0	x_1	x_2	x_3
0	1	-3	-6	8			
1	1	-21	84	-64	4.5826	2.0	0.8729
2	1	-273	4368	-4096	4.0648	2.0	0.9841
3	1	-65793	16843008	-16777216	4.0020	2.0	0.9995

The exact values are 1, -2, and 4. Signs must be determined by substituting the estimates in the original polynomial. When we substitute an estimated root, if the value of the polynomial is zero, then the root is positive; otherwise negative.

Difference Operators and Central Difference Interpolation Formulae

C.1

INTRODUCTION

In Chapter 9 we have already discussed briefly the application of finite differences for interpolating the function values. Here we consider again the finite differences of functions and discuss in detail various operators used on them. We also discuss here some of the central difference formulae used for interpolation.



FINITE DIFFERENCES

Suppose we have a function, $f(x)$, whose values are known (or tabulated) at a set of points $x_0, x_1, x_2, \dots, x_n$. Let us denote the function values $f(x_0), f(x_1), \dots, f(x_n)$ by f_0, f_1, \dots, f_n . The difference between any two consecutive function values is called the *finite difference*. The difference in function values

$$f(x_{i+1}) - f(x_i) = f_{i+1} - f_i \quad (C.1)$$

is known as the *first forward difference* at $x = x_i$. We may denote this first difference at $x = x_i$ as

$$\Delta f_i = f_{i+1} - f_i \quad (C.2)$$

where Δ is an operator called the *forward difference operator*.

By applying the process repeatedly we generate the second forward difference, third forward difference, and so on. Thus the k th forward difference of $f(x)$ at $x = x_i$ is given by

$$\Delta^k f_i = \Delta(\Delta^{k-1} f_i) = \Delta^{k-1} f_{i+1} - \Delta^{k-1} f_i \quad (C.3)$$

Example C.1

Find expression for $\Delta^2 f_i$ and $\Delta^3 f_i$.

$$\begin{aligned}\Delta^2 f_i &= \Delta(\Delta f_i) \\&= \Delta(f_{i+1} - f_i) \\&= \Delta f_{i+1} - \Delta f_i \\&= f_{i+2} - f_{i+1} - f_{i+1} + f_i \\&= f_{i+2} - 2f_{i+1} + f_i \\ \Delta^3 f_i &= \Delta(\Delta^2 f_i) \\&= \Delta(f_{i+2} - 2f_{i+1} + f_i) \\&= f_{i+3} - 3f_{i+2} + 3f_{i+1} - f_i\end{aligned}$$

The difference in function values

$$f(x_i) - f(x_{i-1}) = f_i - f_{i-1} \quad (C.4)$$

is known as the *first backward difference* at $x = x_i$. This is denoted by

$$\nabla f_i = f_i - f_{i-1} \quad (C.5)$$

where ∇ is an operator called the *backward difference operator*.

The k th backward difference of $f(x)$ at $x = x_i$ is defined by

$$\nabla^k f_i = \nabla(\nabla^{k-1} f_i) = \nabla^{k-1} f_i - \nabla^{k-1} f_{i-1} \quad (C.6)$$

Forward and backward differences introduce asymmetry. Sometimes we may need formulae which are symmetrical about the points of interest. Such formulae are based on central differences.

The difference in function values

$$f_{i+1/2} - f_{i-1/2} \quad (C.7)$$

is known as the *first central difference* of $f(x)$ at $x = x_i$. This is denoted by

$$\delta f_i = f_{i+1/2} - f_{i-1/2} \quad (C.8)$$

where δ is called the *central difference operator*.

The k th central difference at $x = x_i$ is defined as

$$\delta^k f_i = \delta(\delta^{k-1} f_i) = \delta^{k-1} f_{i+1/2} - \delta^{k-1} f_{i-1/2} \quad (C.9)$$

Note that if the function values are only known at x_1, x_2, \dots, x_n , then $f(x_{i+1/2})$ is indeterminate and therefore $\delta f(x_i)$ is not computable. However, we can evaluate $\delta^2 f(x_i)$ as follows:

$$\delta^2 f_i = \delta(f_{i+1/2} - f_{i-1/2})$$

$$\begin{aligned} &= f_{i+1} - f_i - f_i + f_{i-1} \\ &= f_{i+1} - 2f_i + f_{i-1} \end{aligned} \quad (\text{C.10})$$

In general, we can compute $\delta^k f_i$ for all positive integers k .



DIFFERENCE OPERATORS

We have seen three difference operators, namely, forward operator Δ , backward operator ∇ , and central operator δ . We consider here some more operators that are often used in the manipulation of interpolation formulae.

Shift Operator

We have an operator known as *shift* or *displacement* or *translation operator* denoted by E . The displacement or shift operator is defined as

$$Ef_i = f_{i+1} \quad (\text{C.11})$$

$$E^k f_i = f_{i+k} \quad (\text{C.12})$$

If the values of x_i are equally spaced such that

$$x_{i+1} - x_i = h$$

then

$$x_{i+1} = x_i + h$$

Eq. (C.12) may be written as

$$E^k f(x_i) = f(x_i + kh) \quad (\text{C.13})$$

Inverse Operator

An operator opposite of E is known as *inverse operator* and is defined as

$$E^{-1} f(x) = g(x) \quad (\text{C.14})$$

Note that

$$Eg(x) = EE^{-1} f(x) = f(x)$$

$$Eg(x) = g(x + h)$$

Then

$$g(x + h) = f(x)$$

Therefore,

$$g(x) = f(x - h)$$

That is,

$$E^{-1} f(x) = f(x - h) \quad (\text{C.15})$$

Similarly,

$$E^{-k} f(x) = f(x - kh) \quad (\text{C.16})$$

Averaging Operator

The averaging operator μ is defined as

$$\delta f_i = \frac{1}{2} (f_{i+1/2} + f_{i-1/2}) \quad (\text{C.17})$$

That is,

$$\mu f(x_i) = \frac{1}{2} \left[f\left(x_i + \frac{h}{2}\right) + f\left(x_i - \frac{h}{2}\right) \right]$$

C.4

RELATIONS BETWEEN THE OPERATORS

The difference operators are related to one another in a number of ways. We consider here a few of them.

Relation between Δ and E

We know that

$$\begin{aligned} \Delta f(x) &= f(x + h) - f(x) \\ &= Ef(x) - 1.f(x) \\ &= (E - 1)f(x) \end{aligned}$$

Therefore,

$$\boxed{\Delta = E - 1 \quad \text{or} \quad E = \Delta + 1} \quad (\text{C.18})$$

Relation between ∇ and E

$$\begin{aligned} \nabla f(x) &= f(x) - f(x - h) \\ &= 1.f(x) - E^{-1}f(x) \\ &= (1 - E^{-1})f(x) \end{aligned}$$

Therefore

$$\boxed{\nabla = 1 - E^{-1}} \quad (\text{C.19})$$

or

$$E = (1 - \nabla)^{-1}$$

(since $(E^{-1})^{-1} = E$)

Relation between Δ , ∇ and E

We have

$$\begin{aligned} E\nabla &= E(1 - E^{-1}) \\ &= E - 1 \\ &= \Delta \end{aligned}$$

Therefore

$$\boxed{E\nabla = \Delta} \quad (\text{C.20})$$

Similarly, we can show that

$$E^{-1}\Delta = \nabla$$

Relation between E , δ and μ

We have

$$\begin{aligned}\delta f(x) &= f\left(x + \frac{h}{2}\right) - f\left(x - \frac{h}{2}\right) \\ &= E^{1/2}f(x) - E^{-1/2}f(x)\end{aligned}$$

Therefore,

$$\delta = E^{1/2} - E^{-1/2} = E^{-1/2}\Delta - E^{1/2}\nabla \quad (\text{C.21})$$

Similarly, we can show that

$$\mu = \frac{1}{2}(E^{1/2} + E^{-1/2}) \quad (\text{C.22})$$

From Eqs (C.21) and (C.22),

$$\mu \delta = \frac{1}{2}(E - E^{-1})$$

We know

$$\Delta + \nabla = (E - 1)(1 - E^{-1}) = E - E^{-1}$$

Then,

$$\mu \delta = \frac{1}{2}(\Delta + \nabla)$$

Example C.2

Prove that $\Delta - \nabla = \Delta\nabla$.

$$\begin{aligned}\Delta - \nabla &= (E - 1) - (1 - E^{-1}) \\ &= E - 2 + E^{-1}\end{aligned}$$

$$\begin{aligned}\Delta\nabla &= (E - 1)(1 - E^{-1}) \\ &= E - 2 + E^{-1}\end{aligned}$$

Therefore

$$\Delta - \nabla = \Delta\nabla$$

Example C.3

Show that $\Delta^2 = E^2 - 2E + 1$

$$\begin{aligned}\Delta^2 &= (E - 1)(E - 1) \\ &= E^2 - 2E + 1\end{aligned}$$

Example C.4.

Prove $\delta^2 = \Delta - \nabla$

$$\begin{aligned}\delta^2 &= [E^{1/2} - E^{-1/2}]^2 \\ &= E + E^{-1} - 2 \\ \Delta - \nabla &= (E - 1) - (1 - E^{-1}) \\ &= E + E^{-1} - 2\end{aligned}$$

Hence,

$$\delta^2 = \Delta - \nabla$$

From Example C.2, we know

$$\Delta - \nabla = \Delta \nabla$$

and, therefore,

$$\delta^2 = \Delta \nabla$$

From Example C.3, we can prove that

$$\begin{aligned}\Delta^2 &= E^2 - 2E + 1 \\ &= E(E - 2 + E^{-1}) \\ &= E(\Delta - \nabla) = E \Delta \nabla\end{aligned}$$

Similarly, the readers may attempt to prove the following relations.

$$1. \quad \mu^2 = 1 + \frac{\delta^2}{4} \quad \text{or} \quad \delta^2 = 4(\mu^2 - 1)$$

$$2. \quad \frac{\Delta}{\nabla} - \frac{\nabla}{\Delta} = \Delta + \nabla$$

$$3. \quad \frac{\Delta}{\nabla} - \frac{\nabla}{\Delta} = 2\mu\delta$$

C.5 CENTRAL DIFFERENCE INTERPOLATION FORMULAE

We have already discussed in Chapter 9 the forward and backward difference formulae. They are useful when the value required is at the beginning or at the end of the table. However, if the point of interest is located at the middle of the table, then we may use the formulae based on central differences.

Let x_0 be the middle point and f_0 the corresponding function value. Consider the values of x on either side x_0 such that

$$f_1 = f(x_0 + h), \quad f_{-1} = f(x_0 - h)$$

$$f_2 = f(x_0 + 2h), \quad f_{-2} = f(x_0 - 2h)$$

and so on.

We can now form a difference table with the values $f(x)$ on either side of x_0 , as shown in Table C.1.

Table C.1 Central difference table

x	f	<i>First difference</i>	<i>Second difference</i>	<i>Third difference</i>	<i>Fourth difference</i>
$x_0 - 2h$	f_{-2}				
		$\Delta f_{-2} = \delta f_{-3/2}$			
$x_0 - h$	f_{-1}		$\Delta^2 f_{-2} = \delta^2 f_{-1}$		
		$\Delta f_{-1} = \delta f_{-1/2}$		$\Delta^3 f_{-2} = \delta^3 f_{-1/2}$	
x_0	f_0		$\Delta^2 f_{-1} = \delta^2 f_0$		$\Delta^4 f_{-2} = \delta^4 f_0$
$x_0 + h$	f_1	$\Delta f_0 = \delta f_{1/2}$		$\Delta^3 f_{-1} = \delta^3 f_{1/2}$	
			$\Delta^2 f_0 = \delta^2 f_1$		
$x_0 + 2h$	f_2	$\Delta f_1 = \delta f_{3/2}$			

The entries in the table are related using the relation between Δ and δ operators. We know that

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

and therefore we can show that

$$\delta^2 f_{-1} = (\Delta E^{-1/2})^2 f_{-1} = \Delta^2 E^{-1} f_{-1} = \Delta^2 f_{-2}$$

Similarly, we can prove for all the entries.

We present here the following central difference interpolation formulae which use differences close to the centre of the table.

1. Gauss forward formula
2. Gauss backward formula
3. Stirling formula
4. Bessel formula
5. Laplace-Everett formula

Gauss Forward Interpolation Formula

We have used the Newton-Gregory forward interpolation formula in Chapter 9. This is given by

$$f(x) = f_0 + p\Delta f_0 + \frac{p(p-1)}{2!} \Delta^2 f_0 + \frac{p(p-1)(p-3)}{3!} \Delta^3 f_0 + \dots \quad (C.23)$$

where

$$x = x_0 + ph \quad \text{or} \quad p = \frac{x - x_0}{h}$$

We can show that

$$\Delta^2 f_0 = \Delta^2 E f_{-1} = \Delta^2 (1 + \Delta) f_{-1} = \Delta^2 f_{-1} + \Delta^3 f_{-1}$$

$$\Delta^3 f_0 = \Delta^3 E f_{-1} = \Delta^3 (1 + \Delta) f_{-1} = \Delta^3 f_{-1} + \Delta^4 f_{-1}$$

$$\Delta^4 f_0 = \Delta^4 f_{-1} + \Delta^5 f_{-1} \quad \text{and so on.}$$

Substituting for $\Delta^2 f_0$, $\Delta^3 f_0$, ..., the above equivalents in Eq. (C.23) and after simplification, we obtain

$$f(x) = f_0 + \binom{p}{1} \Delta f_0 + \binom{p}{2} \Delta^2 f_{-1} + \binom{p+1}{3} \Delta^3 f_{-1} + \binom{p+1}{4} \Delta^4 f_{-1} + \dots \quad (\text{C.24})$$

where

$$\binom{m}{n} = \frac{m(m-1)(m-2)\dots(m-n+1)}{n!}$$

Equation (C.24) is known as *Gauss forward formula*. Note that this formula involves 'odd' differences below the centre line and 'even' differences on the central line.

Gauss Backward Interpolation Formula

Gauss backward formula is obtained using the odd differences above the centre line and even differences on the central line. This is obtained by using the following substitutions in Eq. (C.23).

$$\Delta f_0 = \Delta f_{-1} + \Delta^2 f_{-1}$$

$$\Delta^2 f_0 = \Delta^2 f_{-1} + \Delta^3 f_{-1}$$

$$\Delta^3 f_0 = \Delta^3 f_{-1} + \Delta^4 f_{-1}$$

$$\Delta^3 f_{-1} = \Delta^3 f_{-2} + \Delta^4 f_{-2}$$

$$\Delta^4 f_{-1} = \Delta^4 f_{-2} + \Delta^5 f_{-2}$$

Thus, we obtain the Gauss backward interpolation formula as

$$f(x) = f_0 + \binom{p}{1} \Delta f_{-1} + \binom{p+1}{2} \Delta^2 f_{-1} + \binom{p+1}{3} \Delta^3 f_{-2} + \binom{p+2}{4} \Delta^4 f_{-2} + \dots \quad (\text{C.25})$$

Equation (C.25) is popularly known as *Gauss backward formula*.

Stirling Formula

Stirling formula is obtained by taking the average of the two Gauss formulae. Therefore, adding Eqs (C.24) and (C.25) and dividing by 2, we obtain

$$f(x) = f_0 + p \left(\frac{\Delta f_0 + \Delta f_{-1}}{2} \right) + \frac{p^2}{2!} \Delta^2 f_{-1} + \frac{p(p^2 - 1)}{3!} \left(\frac{\Delta^3 f_{-1} + \Delta^3 f_{-2}}{2} \right) + \frac{p^2(p^2 - 1)}{4!} \Delta^4 f_{-2} + \dots \quad (\text{C.26})$$

Equation (C.26) is known as *Stirling formula*. Note that Eq. (C.26) involves the means of the 'odd' differences just above and just below the central line and 'even' differences on the line. To use this formula, p must satisfy the condition

$$-\frac{1}{2} < p < \frac{1}{2}$$

Bessel Formula

Bessel formula is a modified form of Gauss forward formula, which is given below.

$$f(x) = f_0 + p\Delta f_0 + \frac{p(p-1)}{2!} \Delta^2 f_{-1} + \frac{(p+1)p(p-1)}{3!} \Delta^3 f_{-1} + \dots \quad (\text{C.27})$$

We know

$$\Delta f_0 = f_1 - f_0 \quad \text{and, therefore,} \quad f_0 = f_1 - \Delta f_0 \quad (\text{C.28})$$

$$\Delta^3 f_{-1} = \Delta^2 f_0 - \Delta^2 f_{-1}, \quad \text{then,} \quad \Delta^2 f_{-1} = \Delta^2 f_0 - \Delta^3 f_{-1} \quad (\text{C.29})$$

Similarly,

$$\Delta^4 f_{-2} = \Delta^4 f_{-1} - \Delta^5 f_{-2} \quad (\text{C.30})$$

Now, we rewrite Eq. (C.27) as

$$\begin{aligned} f(x) = & \left(\frac{f_0}{2} + \frac{f_0}{2} \right) + p\Delta f_0 + \frac{1}{2} \frac{p(p-1)}{2!} \Delta^2 f_{-1} + \frac{1}{2} \frac{p(p-1)}{2!} \Delta^2 f_{-1} \\ & + \frac{(p+1)p(p-1)}{3!} \Delta^3 f_{-1} + \dots \end{aligned} \quad (\text{C.31})$$

Now, substituting Eqs (C.28), (C.29), and (C.30) in (C.31), we get

$$\begin{aligned} f(x) = & \frac{f_0}{2} + \frac{f_1 - \Delta f_0}{2} + p\Delta f_0 + \frac{1}{2} \frac{p(p-1)}{2!} \Delta^2 f_{-1} \\ & + \frac{1}{2} \frac{p(p-1)}{2!} (\Delta^2 f_0 - \Delta^3 f_{-1}) + \frac{(p+1)p(p-1)}{3!} \Delta^3 f_{-1} + \dots \\ = & \frac{f_0 + f_1}{2} + \left(p - \frac{1}{2} \right) \Delta f_0 + \frac{p(p-1)}{2!} \frac{(\Delta^2 f_0 + \Delta^2 f_{-1})}{2} \\ & + \frac{\left(p - \frac{1}{2} \right) p(p-1)}{3!} (\Delta^3 f_{-1}) + \dots \end{aligned} \quad (\text{C.32})$$

Equation (C.32) is called *Bessel formula*. Note that the Bessel formula involves odd differences below the centre line and averages of the even differences on and below the central line.

Observe that when $p = \frac{1}{2}$ all odd order differences vanish and we get

$$\begin{aligned} f(x) &= \frac{1}{2}(f_0 + f_1) - \frac{1}{8}\left(\frac{\Delta^2 f_{-1} + \Delta^2 f_0}{2}\right) + \frac{3}{128}\left(\frac{\Delta^4 f_{-2} + \Delta^4 f_{-1}}{2}\right) \\ &\quad - \frac{5}{1024}\left(\frac{\Delta^6 f_{-3} + \Delta^6 f_{-2}}{2}\right) + \dots \end{aligned} \quad (\text{C.33})$$

Equation (C.33) is known as *formula for interpolating to haves*.

Laplace-Everett Formula

Again consider the Gauss forward formula

$$f(x) = f_0 + p\Delta f_0 + \frac{p(p-1)}{2!} \Delta^2 f_{-1} + \frac{(p+1)p(p-1)}{3!} \Delta^3 f_{-1} + \dots \quad (\text{C.34})$$

If we eliminate the odd differences from Eq. (C.34) the result will give us the Laplace-Everett formula. We know

$$\Delta f_0 = f_1 - f_0$$

$$\Delta^2 f_{-1} = \Delta^2 f_0 - \Delta^2 f_{-1}$$

$$\Delta^3 f_{-2} = \Delta^4 f_{-1} - \Delta^4 f_{-2}$$

and so on.

Substituting these in Eq. (C.34) and rearranging the terms, we get

$$\begin{aligned} f(x) &= (1-p)f_0 + \binom{p}{1}f_1 - \binom{p}{1}\Delta^2 f_{-1} + \binom{p+1}{3}\Delta^2 f_0 \\ &\quad - \binom{p+1}{5}\Delta^4 f_{-2} + \binom{p+2}{7}\Delta^4 f_{-1} + \dots \end{aligned} \quad (\text{C.35})$$

Letting

$$p = 1 - q \quad \text{or} \quad 1 - p = q$$

and simplifying, Eq. (C.35) becomes

$$f(x) = \left[qf_0 + \binom{q+1}{3}\Delta^2 f_{-1} + \dots \right] + \left[pf_1 + \binom{p+1}{3}\Delta^2 f_0 + \dots \right] \quad (\text{C.36})$$

Equation (C.36) is known as Laplace-Everett formula. This formula involves only even differences on and below the central line. This can be used when $0 < p < 1$.

Key Terms

<i>Averaging operator</i>	<i>Forward difference operator</i>
<i>Backward difference</i>	<i>Gauss backward formula</i>
<i>Backward difference operator</i>	<i>Gauss forward formula</i>
<i>Bessel formula</i>	<i>Inverse operator</i>
<i>Central difference</i>	<i>Laplace-Everett formula</i>
<i>Central difference operator</i>	<i>Shift operator</i>
<i>Displacement operator</i>	<i>Stirling formula</i>
<i>Forward difference</i>	<i>Translation operator</i>

APPENDIX D

C Programs

D.1

INTRODUCTION

C programs are basically made of functions. As we know, a function is a unit of a program that performs a particular task. A C function is similar to a subroutine in FORTRAN. A function begins running when its name is used in the program.

All C programs need a *main* function which is named as **main()**. Therefore, the programs in this appendix begin with the main function. Since the main function does not use any arguments, its name is followed by a set of empty parentheses.

Everything inside a C function is enclosed between two curly brackets, like `{...}`. The left curly bracket signals the beginning of the function and the right one signals the end of the function.

Some features of C that are distinctly different from FORTRAN are:

- C is a free-form language and therefore we may follow any format of statements that may please us.
- A comment or remark can be inserted anywhere that a space can appear in a C program provided it is preceded by `/*` and followed by `*/`
- As a convention (and not as a rule), everything in C is written using lower-case letters. Thus, names of all variables and functions are written in lower-case.
- All C statements must end with a semicolon.
- C does not support statement numbers. However, we can assign a label to a statement as follows:

```
begin : sum = 0.0;
```

We may direct the control to this statement by using the **goto** statement as

```
goto begin;
```

- C language does not have STOP and END statements. While the closing bracket } represents the "END" of a function, STOP may be replaced by invoking the **exit()** function available in C.
- Unlike FORTRAN, all variables and functions that return values must explicitly be declared for their types.

```
int i,j,a,b; /* declaring integer variables */
float x,y,a[10]; /*declaring real variables */
```

- C supports what are known as preprocessor directives **#include** and **#define**. The **#include** directive is used to include in the program library functions and other files. We may use the **#define** directive to define constants and functions that are used in the program.
- C does not support any operator for exponentiation. Instead, it uses a function **pow(x,y)** to compute x^y .
- C uses square brackets to represent arrays variables, like, $a[i]$, $b[i][j]$, etc.
- Unlike FORTRAN, C array elements are numbered from ZERO, not ONE. That is, the array $a[3]$ will be represented in the memory as $a[0]$, $a[1]$, and $a[2]$.
- C uses an operator & known as address operator to read the values from the keyboard. We may use this operator to obtain the address of a variable in the memory. For example, & x gives the address of the location of x .
- C defines % as the modulus operator. This operator divides the first operand by the second and gives the remainder, (not the quotient) as the result.
- C defines another operator known as "star" operator * to define a pointer. The statements

```
int a;
int *p; /* p is declared as pointer */
p=&a; /* address of a is assigned to p */
```

make **p** point to **a**. Now, the statements $*p+5$ and $a+5$ both would give the same result. That is, ***p** means the value stored at the address pointed to by **p**.

- C uses the following relational and logical operators:

= =	is equal to
<	is less than
>	is greater than
<=	is less than or equal to
>=	is greater than or equal to
!=	is not equal to
	logical OR
&&	logical AND

- C has two more control statements **continue** and **break** in addition to the conventional **goto** statement. These two are used inside a loop. While the break transfers the control to the first statement outside the loop, the continue statement skips the remaining part of the loop.

D.2 FORTRAN TO C CONVERSION

The FORTRAN programs in the text have been translated almost line by line using C equivalents given in Table D.1

Table D.1 C equivalents

<i>FORTRAN Statements</i>	<i>C Equivalents</i>
INTEGER	int
REAL	float, double
PARAMETERS (M = 100)	#define M 100
INTRINSIC	#include <math.h>
READ(*,*) statement	scanf() statement
WRITE (*,*) statement	printf() statement
Function subprograms	Preprocessor macro definition
Statement number	Label name
GOTO <number>	goto <label>
GOTO inside a loop	break or continue
STOP inside a program	exit() function
IF (...) THEN...	if (...) ...;
IF (...) THEN ... ELSE	if (...) ... else
DO ... CONTINUE	for (...) {}
END	}

The preprocessor directives **#define** and **#include** are always placed before the main() function. Any constants and variables declared before the main function are considered to be global and are available to all the functions defined after that point in the file.

It should be noted that the programs have been translated line by line so as to enable the readers to understand the logic easily. It is quite possible to simplify or improve the efficiency of C programs in this appendix using some of the unique features of C. This requires a complete understanding of the language.

Readers new to C language must read an introductory book on C such as *Programming in ANSI C* by E Balagurusamy before attempting to analyse and apply the C programs given in this appendix.

Table D.2 List of programs

<i>Program No.</i>	<i>Name</i>	<i>Description</i>
1	POLY	Program POLY evaluates a polynomial of degree n at any point x using Horner's rule.
2	BISECT	This program finds a root of a nonlinear equation using the bisection method.
3	FALSE	This program finds a root of a nonlinear equation by false position method.
4	NEWTON	This program finds a root of a nonlinear equation by Newton-Raphson method.
5	SECANT	This program finds a root of a nonlinear equation by secant method.
6	FIXEDP	This program finds a root of a function using the fixedp point iteration method.
7	MULTIR	The program finds all the real roots of a polynomial.
8	COMPR	This program locates all the roots, both real and complex, using the Bairstow method.
9	MULLER*	This program evaluates root of a polynomial using Muller's method.
10	LEG1	This program solves a system of linear equations using simple Gaussian elimination method.
11	LEG2	This program solves a system of linear equations using Gaussian elimination with partial pivoting.
12	DOLIT	This program solves a system of linear equations using Dolittle LU decomposition.
13	JACIT	This program uses the subprogram JACOBI to solve a system of equations by Jacobi iteration method.
14	GASIT	This program uses the subprogram GASEID to solve a system of equations by Gauss-Seidel iteration method.
15	LAGRAN	This program computes the interpolation value at a specified point, given a set of data points, using the Lagrange interpolation representation.
16	NEWINT	This program constructs the Newton interpolation polynomial for a given set of data points and then computes interpolation value at a specified value.
17	SPLINE	This program computes the interpolation value at a specified value, given a set of table points, using the natural cubic spline interpolation.
18	LINREG	This program fits a line $Y = A + BX$ to a given set of data points by the method of least squares.
19	POLREG	This program fits a polynomial curve to a given set of data points by the method of least squares.
20	NUDIF*	This program computes the derivative of a tabulated function at a specified value using the Newton interpolation approach.

(Contd.)

Table D.2 (Contd.)

Program No.	Name	Description
21	TRAPE1	This program integrates a given function using the trapezoidal rule.
22	SIMS1	This program integrates a given function using the Simpson's 1/3 rule.
23	ROMBRG	This program performs Romberg integration by bisecting the intervals N times.
24	TRAPE2*	This program integrates a tabulated function using the trapezoidal rule.
25	SIMS2*	This program integrates a tabulated function using the Simpson's 1/3 rule. If the number of segments is odd, the trapezoidal rule is used for the last segment.
26	EULER	This program estimates the solution of the first order differential equation $y' = f(x, y)$ at the given point using Euler's method.
27	HEUN	This program solves the first order differential equation $y' = f(x, y)$ using the Heun's method.
28	POLYGN	This program solves the differential equation of type $y' = f(x, y)$ by polygon method.
29	RUNGE4	This program computes the solution of first order differential equation of type $y' = f(x, y)$ using the 4th order Runge-Kutta method.
30	MILSIM	This program solves the first order differential equation $y' = f(x, y)$ using the Milne-Simpson method.

* These are new programs in C and are not available in FORTNIN version in the text. Readers may try to develop FORTRAN equivalents of these programs.

Program 1 POLY

```
/*
 * Main program
 * Program POLY evaluates a polynomial of degree n
 * at any point X using Horner's rule
 *
 * Subroutines used
 * horner
 *
 * Variables used
 * n - Degree of polynomial
 * a - Array of polynomial coefficients
 * x - Point of evaluation
 * p - Value of polynomial at x
 *
 * Constants used
 * NIL
 */

```

536 Numerical Methods

```
main( )
{
    int n,i;      /* Declaration of variables */
    float x,p,a[10];
    float horner(int n, float a[], float x);

    printf("Input degree of polynomial, n\n");
    scanf("%d", &n);
    printf("Input polynomial coefficients a(0) to a(n) \n");
    for(i=0;i<=n;i++)
        scanf("%f", &a[i]);
    printf("Input value of x (point of evaluation) \n");
    scanf("%f", &x);

/* Evaluating polynomial at x using Horner's rule */

    p = horner(n,a,x); /* Calling the function horner */

/* Writing the result */
    printf("\n");
    printf("f(x) = %f at x = %f \n", p, x);
    printf("\n");
}

/* End of main program POLY */
/* -----
/* Defining the function horner */
float horner(int n, float a[], float x)
/*
    horner computes the value of a polynomial of order n
    at any given point x. */
{
    int i; /* Local variables */
    float p;

    p = a[n];
    for(i=n-1;i>=0;i--)
    {
        p = p*x + a[i];
    }
    return(p);
}
/* End of Function horner */
/* ----- */
```

Program 2 BISECT

```

/*
 * Main program
 * This program finds a root of a nonlinear
 * equation using the bisection method
 *
 * Functions invoked
 * Macro F(x)
 *
 * Subroutines used
 * bim()
 *
 * Variables used
 * a - Left endpoint of interval
 * b - Right endpoint of interval
 * s - Status
 * root - Final Solution
 * count - Number of Iterations done
 *
 * Constants used
 * EPS - Error bound
 */

#include <stdio.h>
#include <math.h>
#define EPS      0.000001
#define F(x)     (x)*(x)+(x)-2

main( )
{
    int s, count;
    float a,b,root;

    printf("\n");
    printf("SOLUTION BY BISECTION METHOD \n");
    printf("\n");
    printf("Input starting values \n");
    scanf("%f %f",&a,&b);

/* Calling the subroutine bim() */
    bim(&a, &b, &root, &s, &count);
    if(s==0)
    {
        printf("\n");
        printf("Starting points do not bracket any root \n");
        printf("(Check whether they bracket EVEN roots \n");
        printf("\n");
    }
}

```

```

    }
else
{
    printf("\nRoot = %f \n", root);
    printf("F(root) = %f \n", F(root));
    printf("\n");
    printf("Iterations = %d \n", count);
    printf("\n");
}
/* End of main program */
/* -----
/* Defining the subroutine bim() */
bim(float *a, float *b, float *root, int *s, int *count)
/* This subroutine finds a root of nonlinear equation
   in the interval [a,b] using the bisection method */
{
    float x1,x2,x0,f0,f1,f2;
/* Function values at initial points */
    x1 = *a;
    x2 = *b;
    f1 = F(x1);
    f2 = F(x2);
/* Test if initial values bracket a SINGLE root */
    if(f1*f2 > 0)
    {
        *s = 0;
        return; /* Program terminated */
    }
    else
/* Bisect the interval and locate the root iteratively */
    {
        *count = 0;

        begin: /* Iteration begins */
        x0 = (x1 + x2)/2.0;
        f0 = F(x0);
        if(f0==0)
        {
            *s = 1;
            *root = x0;
            return;
        }
        if(f1*f0 < 0)

```

```

{
    x2 = x0;
}
else
{
    x1 = x0;
    f1 = f0;
}
/* Test for accuracy and repeat the process, if necessary */
if(fabs((x2-x1)/x2) < EPS)
{
    *s = 1;
    *root = (x1+x2)/2.0;
    return; /* Iteration ends */
}
else
{
    *count = *count + 1;
    goto begin;
}
}
/* End of subroutine bim() */
/* ----- */

```

Program 3 FALSE

```

/* -----
 * Main program
 * This program finds a root of a nonlinear equation
 * by false position method
 * -----
 * Functions invoked
 * Macro F(x)
 * -----
 * Subroutines used
 * fal()
 * -----
 * Variables used
 * a - Left endpoint of interval
 * b - Right endpoint of interval
 * s - Status
 * root - Final solution
 * count - Number of iterations completed
 * -----
 * Constants used
 * EPS - Error bound
 * -----
 * /

```

540 Numerical Methods

```
#include <math.h>
#define EPS      0.000001
#define F(x)      (x)*(x)+(x)-2
main( )
{
    int s,count;
    float a,b,root;

    printf("\n");
    printf("SOLUTION BY FALSE POSITION METHOD \n");
    printf("\n");
    printf("Input starting values \n");
    scanf("%f %f", &a, &b);
/* Calling the function fal() */

    fal(&a,&b,&s,&root,&count);

    if(s==0)
    {
        printf("\n");
        printf("Starting points do not bracket
              any root \n");
        printf("\n");
    }
    else
    {
        printf("\n");
        printf("Root = %f \n", root);
        printf("F(root) = %f \n", F(root));
        printf("\n NO. OF ITERATIONS = %d \n", count);
    }
}
/* End of main() program */

/* -----
/* Defining the subroutine fal() */
fal(float *a, float *b, int *s, float *root, int *count)
/* fal finds a root of a nonlinear equation */

{
    float x1,x2,x0,f0,f1,f2;
    x1 = *a;
    x2 = *b;
    f1 = F(x1);
    f2 = F(x2);

    /* Test if a and b bracket a root */
```

```

if(f1*f2 > 0)
{
    *s = 0;
    return;           /* Program terminated */
}
else
{
    printf("      n      x1      x2 \n");
}

*count = 1;
begin:                   /* Iteration begins */
x0 = x1-f1*(x2-x1)/(f2-f1);
f0 = F(x0);
if(f1*f2 < 0)
{
    x2 = x0;
    f2 = f0;
}
else
{
    x1 = x0;
    f1 = f0;
}
printf("%5d %15.6f %15.6f \n", *count, x1, x2);
/* Test whether the desired accuracy has been achieved */
if(fabs((x2-x1)/x2) < EPS)
{
    *s = 1;
    *root = (x1+x2)*0.5;
    return;           /* Iteration ends */
}
else
{
    *count = *count + 1;
    goto begin;
}
}

/* End of subroutine fal() */

/* -----
 * Main program
 * This program finds a root of a nonlinear equation
 * by Newton-Raphson method
 */

```

Program 4 NEWTON

```

/*
 * -----
 * Main program
 * This program finds a root of a nonlinear equation
 * by Newton-Raphson method
 */

```

```

* -----
* Functions invoked
*   Macros F(x), FD(x), Library function fabs()
*
* -----
* Subroutines used
*   NTL
*
* -----
* Variables used
*   x0 - Initial value of x
*   xn - New value of x
*   fx - Function value at x
*   fdx - Value of function derivative at x
*   count - Number of iterations done
*
* -----
* Constants used
*   EPS - Error bound
*   MAXIT - Maximum number of iterations permitted
* -----
* /

```

```

#include <math.h>
#define EPS 0.000001
#define MAXIT 20
#define F(x) (x)*(x)+(x)-2
#define FD(x) 2*(x)+1

main( )
{
    int count;
    float x0, xn, fx, fdx;

    printf("\n");
    printf("Input initial value of x \n");
    scanf("%f", &x0);
    printf("\n");
    printf("      SOLUTION BY NEWTON-RAPHSON METHOD \n");
    printf("\n");

    count = 1;
    begin:                               /* Iteration begins */
    fx = F(x0);
    fdx = FD(x0);
    xn = x0 - fx/fdx;

    if(fabs((xn-x0)/xn) < EPS)          /* Iteration ends */
    {
        printf("Root = %f \n", xn);
        printf("Function value = %f \n", F(xn));
        printf("Number of iterations = %d \n", count);
        printf("\n");
    }
}

```

```

    }
else
{
    x0 = xn;
    count = count + 1;
    if(count < MAXIT)
    {
        goto begin;
    }
    else
    {
        printf("\n SOLUTION DOES NOT CONVERGE \n");
        printf("IN %d ITERATIONS \n", MAXIT);
    }
}
}

/* End of main() program */
/* -----

```

Program 5 SECANT

```

/*
 * Main program
 *   This program finds a root of a nonlinear
 *     equation by secant method
 *
 * Functions invoked
 *   Macro F(x)
 *
 * Subroutines used
 *   sec()
 *
 * Variables used
 *   a - Left endpoint of interval
 *   b - Right endpoint of interval
 *   x1 - New left point
 *   x2 - New right point
 *   root - Final solution
 *   count - Number of iterations completed
 *
 * Constants used
 *   EPS - Error bound
 *   MAXIT - Maximum number of iterations permitted
 */

#include <math.h>
#define EPS      0.000001

```

544 Numerical Methods

```
#define MAXIT 50
#define F(x) (x)*(x)+(x)-2
main()
{
    float a,b,root,x1,x2;
    int count,status;

    printf("\n");
    printf("          SOLUTION BY SECANT METHOD \n");
    printf("\n");
    printf("Input two starting points \n");
    scanf("%f %f", &a, &b);

    sec(&a, &b, &x1, &x2, &root, &count, &status);

    if(status == 1)
    {
        printf("\n");
        printf("          DIVISION BY ZERO \n");
        printf("\nLast x1 = %f \n", x1);
        printf("\nLast x2 = %f \n", x2);
        printf("\nNO. OF ITERATIONS = %d \n", count);
        printf("\n");
    }
    else if(status == 2)
    {
        printf("\n");
        printf("NO CONVERGENCE IN %d ITERATIONS \n", MAXIT);
        printf("\n");
    }
    else
    {
        printf("\n");
        printf("Root = %f \n", root);
        printf("Function value at root = %f \n", F(root));
        printf("\n");
        printf("NO. OF ITERATIONS = %d \n", count);
        printf("\n");
    }
}
/* End of main() program */
/* -----
/* Defining subroutine sec() */
sec(float *a, float *b, float *x1, float *x2, float
    *root, int *count, int *status)
```

```
/* This subroutine computes a root of an equation using
   the secant method. */
{
    float x3,f1,f2,error;
/* Function values at initial points */
    *x1 = *a;
    *x2 = *b;
    f1 = F(*a);
    f2 = F(*b);

/* Compute the root iteratively */
    *count = 1;
    begin:           /* Iteration process begins */
    if(fabs(f1-f2) <= 1.E-10)
    {
        *status = 1;
        return;          /* Program terminated */
    }
    x3 = *x2 - f2*(x2-*x1)/(f2-f1);
    error = fabs((x3-*x2)/x3);

/* Test for accuracy */
    if(error > EPS)
    {
/* Test for convergence */
        if(*count == MAXIT)
        {
            *status = 2;
            return;          /* Program terminated */
        }
        else
        {
            *x1 = *x2;
        }
        *x2 = x3;
        f1 = f2;
        f2 = F(x3);
        *count = *count + 1;
        goto begin;      /* Compute next approximation */
    }
    else
    {
        *root = x3;
        *status = 3;
        return;          /* Iteration ends */
    }
}
```

```

    }
}
/* End of subroutine sec() */
/* ----- */
```

Program 6 FIXEDP

```

/* -----
 * Main program
 *   This program finds a root of a function using
 *   the fixedp point iteration method
 *
 * -----
 * Functions invoked
 *   Library function fabs() and macro G(x)
 *
 * -----
 * Subroutines used
 *   NIL
 *
 * -----
 * Variables used
 *   x0 - Initial guess
 *   x - Estimated root
 *   error - Relative error
 *
 * -----
 * Constants used
 *   EPS - Error bound
 *   MAXIT - Maximum iterations allowed
 *
# include <math.h>
# define EPS  0.000001
# define G(x)  2.0-(x)*(x)

main( )
{
    int MAXIT, i;
    float x0, x, error;
    printf("\n      SOLUTION BY FIXED-POINT METHOD \n");
    printf("\n");
    printf("Input initial estimate of a root \n");
    scanf("%f", &x0);
    printf("Maximum iterations allowed \n");
    scanf("%d", &MAXIT);
    printf("\n      ITERATION      VALUE OF X      ERROR \n");
/*Iteration process begins*/
    for(i=1;i<=MAXIT;i++)
    {
        x = G(x0);
```

```

    error = fabs((x-x0)/x);
    printf("%10d %10.7f %10.7f \n", i, x, error);
    if (error < EPS)
        goto end; /*Iteration process ends*/
    else
        x0 = x;
    }
    printf("\nProcess does not converge to a root.\n");
    printf("Exit from iteration loop.\n");
    end:
    ;
}
/* End of main() program/
/* ----- */

```

Program 7 MULTIR

```

/* -----
 * Main program
 * The program finds all the real roots of a polynomial
 * -----
 * Functions invoked
 * NIL
 * -----
 * Subroutines used
 * newton()
 * dflat()
 * -----
 * Variables used
 * n - Degree of polynomial
 * a - Polynomial coefficients A(N+1)
 * x0 - Initial guess
 * xr - Root obtained by Newton method
 * root - Root Vector
 * status - Solution status
 * -----
 * Constants used
 * EPS - Error bound
 * MAXIT - Maximum iterations permitted
 * -----
#include <math.h>
#define EPS 0.000001
#define MAXIT 50

void main( )
{
    int n, status, i, j;

```

```

float a[11], root[10], x0, xr;
void dflat(int n, float a[11], float xr);
void newton(int n, float a[11], float x0, int *status,
            float *xr);

printf("\n");
printf("\n      EVALUATION OF MULTIPLE ROOTS \n");
printf("\n");
printf("Input degree of polynomial, n. \n");
scanf("%d", &n);

printf("\nInput poly coefficients, a(1) to a(n+1). \n");
for(i = 1; i <= n+1; i++)
    scanf("%f", &a[i]);
printf("\nInput initial guess of x \n");
scanf ("%f", &x0);
printf("\n");

/* Process of root searching begins */
for(i=n;i>=2;i--)
/* Find ith root by Newton Method */
{
    newton(n,a,x0,&status,&xr);

    if(status == 2)
    {
        for(j=n;j>=i+1;j--)
            printf("root %d = %f \n",j, root[j]);

        printf("\nNext root does not converge in \n");
        printf("%d iterations \n", MAXIT);
        printf("\n");
        goto end; /* Processing ends */

    }
    root[i] = xr;
}

/* Deflate the polynomial by division (x-xr) */
dflat(n,a,xr);
x0 = xr;
/* Proceed to next root */
} /* End of for loop */

/* Compute the last root */
root[1] = -a[1]/a[2];

/* Write results */
printf("\n  ROOTS OF POLYNOMIAL ARE: \n");
printf("\n");
for (i=1;i<=n;i++)
    printf("ROOT %d = %f \n", i, root[i]);

```

```
    printf("\n");
  end:
  printf("END");
}
/* End of main() program */
/* -----
/* Defining the subroutine newton() */
void newton(int n, float a[11], float x0, int *status,
            float *xr)
/* This subroutine finds a root of the polynomial using
   the Newton-Raphson method */
{
    int i, count;
    float fx; /* Value of polynomial at x0 */
    float fdx; /* Value of polynomial derivative at x0 */
    count = 1;
/* Compute the value of function at x0 */
    begin:
    fx = a[n+1];
    for(i=n;i>=1;i--)
        fx = fx * x0 + a[i];
/* Compute the value of derivative at x0 */
    fdx = a[n+1] * n;
    for(i=n;i>=2;i--)
        fdx = fdx * x0 + a[i] * (i-1);
/* Compute a root xr */
    *xr = x0-fx/fdx;
/* Test for accuracy */
    if(fabs((*xr-x0)/(*xr)) <= EPS)
    {
        *status = 1;
        return;
    }
/* Test for convergence */
    if(count<MAXIT)
    {
        x0 = *xr;
        count = count + 1;
        goto begin;
    }
    else
    {
```

```

        *status = 2;
        return;
    }
}

/* End of subroutine newton() */

/* -----
/* Defining the subroutine dflat() */
void dflat(int n, float a[11], float xr)
/* This subroutine reduces the degree of polynomial by
one using synthetic division */
{
    float b[11];
    int i;

/* Evaluate the coefficients of the reduced polynomial */
    b[n+1] = 0;
    for(i=n;i>-1;i--)
        b[i] = a[i+1] + xr * b[i+1];

/* Change coefficients from b array to a array */
    for(i=1;i<=n+1;i++)
        a[i] = b[i];
}
/* End of subroutine dflat() */
/* ----- */

```

Program 8 COMPR

```

* -----
* Main program
*   This program locates all the roots, both real
*   and complex, using the Bairstow method
* -----
* Functions invoked
*   NIL
* -----
* Subroutines used
*   input,bstow,quad,output
* -----
* Variables used
*   n - Degree of polynomial
*   a - Array of coefficients of polynomial
*   u0,v0 - Initial values of coefficients of the
*           quadratic factor

```

```

*   u,v   - Computed values of coefficients of the   *
*           quadratic factor
*   b - Coefficients of the reduced polynomial   *
*   x1,x2 - Roots of the quadratic factor       *
*   type - Type of roots (real,imaginary or equal) *
* -----
* Constants used                                *
*   EPS - Error bound                          *
* -----
* /



#include <math.h>
#define EPS    0.000001
#define image  1
#define equal  2
#define unequal 3

main( )
{
    int n, i;
    float a[11], b[11], u0, v0, u, v, x1, x2, d0, d1,
          d2, root, type, status;

    printf("\n");
    printf("      EVALUATION OF COMPLEX ROOTS \n");
    printf("\n");

/* Get input data */
    printf("Input degree of polynomial, n \n");
    scanf("%d",&n);
    printf("\n Input coefficients a(n+1) to a(1) \n");
    for(i=n+1;i>=1;i--)
        scanf("%f",&a[i]);
    printf("\n Give initial values u0 and v0 \n");
    scanf("%f %f",&u0,&v0);

begin:
    if(n > 2)
    {
        /* Obtain a quadratic factor */

bstow(n,a[11],b[11],u0,v0,&u,&v,&status);
        if(status == 1)
        {
            d2 = 1;
            d1 = -u;
            d0 = -v;
        }
        else

```

```

{
    printf("\n No Convergence in 100 iterations \n");
    goto end;
}
/* Find roots of the quadratic factor */
quad(d2,d1,d0,&x1,&x2,&type);
/* Print the roots */
output(n,type,x1,x2);
/* Set the coefficients of the factor polynomial */
n = n-2;
for(i=1;i<=n+1;i++)
    a[i] = b[i+2];
/* Set initial values for next quadratic factor */
u0 = u;
v0 = v;
goto begin;
} /* endif */
if(n == 2) /* polynomial is quadratic */
{
    quad(a[3],a[2],a[1],&x1,&x2,&type);
    output(n,type,x1,x2);
}
else /* last root of an odd order polynomial */
{
    root = -a[1]/a[2];
    printf("\n");
    printf("Final root = %f \n", root);
    printf("\n");
}
end:
printf("End");
}

/* End of main() program */
/* -----
/* Defining the subroutine bstow() */
bstow(int n,float a[11],float b[11],float u0,float v0,
      float *u,float *v,float *status)
/* This subroutine finds the quadratic factor using
multivariable Newton's method and also finds the
reduced polynomial */

```

```

{
    float d,delu,delv,c[11];
    int count,i;

    count = 1;
    begin:
    b[n+1] = a[n+1];
    b[n] = a[n] + u0*b[n+1];
    for(i=n-1;i>=1;i--)
        b[i] = a[i] + u0 * b[i+1] + v0 * b[i+2];
    c[n+1] = 0;
    c[n] = b[n+1];
    for(i=n-1;i>=1;i--)
        c[i] = b[i+1] + u0 * c[i+1] + v0 * c[i+2];
    d = c[2] * c[2] - c[1] * c[3];
    delu = -(b[2] * c[2] - b[1] * c[3])/d;
    delv = -(b[1] * c[2] - b[2] * c[3])/d;
    *u = u0 + delu;
    *v = v0 + delv;

    if(fabs(delu/*u) <= EPS && fabs(delv/*v) <= EPS)
    {
        *status = 1;
        return;
    }

    if(count < 100)
    {
        u0 = *u;
        v0 = *v;
        count = count + 1;
        goto begin;
    }
    else
    {
        *status = 2;
        return;
    }
}
/* End of subroutine bstow() */
/* -----
/* Define the subroutine quad() */
quad(float a,float b,float c,float *x1,float *x2,float
 *type)

```

```

/* This subroutine solves a quadratic equation of type
   a(x*x) + bx + c */
{
    float q;
    q = b*b - 4*a*c;
    if(q < 0.0) /* roots are complei */
    {
        *x1 = -b/(2*a);
        *x2 = sqrt(fabs(q))/(2*a);
        *type = image;
    }
    else if (q == 0.0) /* roots are real and equal */
    {
        *x1 = -b/(2*a);
        *x2 = *x1;
        *type = equal;
    }
    else /* roots are real and unequal */
    {
        *x1 = (-b+sqrt(q))/(2*a);
        *x2 = (-b-sqrt(q))/(2*a);
    }
    return;
}
/* End of subroutine quad() */
/* ----- */

/* Defining output() routine */
output(int n, int type, float x1, float x2)
/* This subroutine displays the roots of the quadratic
   equation */
{
    printf("\n");
    printf("Roots of quadratic factor at n = %d \n", n);
    printf("\n");
    if(type == image)
    {
        printf("Root1 = %f%lj \n", x1, x2);
        printf("Root2 = %f%lj \n", x1, x2);
    }
    else if(type == equal)
    {
        printf("Root1 = %f \n", x1);
        printf("Root2 = %f \n", x2);
    }
    else /* Type == unequal */

```

```

    {
        printf("Root1 = %f \n", x1);
        printf("Root2 = %f \n", x2);
    }
    return;
}
/* End of subroutine output() */
/* ----- */

```

Program 9 MULLER

```

/* -----
 * Main program
 *   This program evaluates root of a polynomial using
 *   Muller's method
 * -----
 * Functions invoked
 *   NIL
 * -----
 * Subroutines used
 *   F(x)
 * -----
 * Variables used
 *   x1,x2,x3 - initial values
 *   f1,f2,f3 - function values at x1,x2,x3
 *   a0,a1,a2 - coefficients of quadratic polynomial
 *   hi - xi-x3
 *   di - function difference fi-f3
 * -----
 * Constants used
 *   EPS - Error bound
 * -----
 * / */

#include <math.h>
#define EPS 0.000001

main( )
{
    float F(float x);
    float x1,x2,x3,x4,f1,f2,f3,f4,h1,h2,d1,d2,a0,a1,a2,h;
    printf("\nInput three initial points \n");
    scanf("%f %f %f", &x1,&x2,&x3);

    f1 = F(x1);
    f2 = F(x2);
    f3 = F(x3);

    begin:
    h1 = x1-x3;

```

```

    h2 = x2-x3;
    d1 = f1-f3;
    d2 = f2-f3;
/* Compute parameters a0,a1,a2 */
    a0 = f3;
    a1 = (d2*h1*h1-d1*h2*h2)/(h1*h2*(h1-h2));
    a2 = (d1*h2-d2*h1)/(h*h2*(h1-h2));

/* Compute h */
    if(a1>0.0)
        h=(-2.0*a0)/(a1+sqrt(a1*a1-4*a2*a0));
    else
        h=(-2.0*a0)/(a1-sqrt(a1*a1-4*a2*a0));
/* Compute x4 and f4 */
    x4 = x3+h;
    f4 = F(x4);

/* Test for accuracy */
    if(f4<=EPS) /* root obtained */
    {
        printf("\n\nROOT BY MULLER'S METHOD\n\n");
        printf("Root=%f\n", x4);
        printf("\n");
    }
    else
    {
        x1 = x2;
        x2 = x3;
        x3 = x4;
        f1 = f2;
        f2 = f3;
        f3 = f4;
        goto begin;
    }
}
/* End of main() program */
/* ----- */
/* Defining the subroutine F(x) */
float F(float x)
{
    float f;
    f = x*x*x+2*x*x+10*x-20;
    return (f);
}
/* End of subroutine () */
/* ----- */

```

Program 10 LEG1

```

/* -----
 * Main program
 * This program solves a system of linear equations
 * using simple Gaussian elimination method
 *
 * -----
 * Functions invoked
 * NIL
 *
 * -----
 * Subroutines used
 * GAUSS1
 *
 * -----
 * Variables used
 * n - Number of equations in the system
 * a - Matrix of coefficients
 * b - Right side vector
 * x - Solution vector
 *
 * -----
 * status - Solution status
 * ----- */
main( )
{
    int status,n,i,j;
    float a[10][10], b[10], x[10];
    printf("\n      SOLUTION BY SIMPLE GAUSS METHOD \n");
    printf("What is the size of the system (n)? \n");
    scanf("%d", &n);
    printf("Input coefficients a(i,j), row-wise, \n");
    printf("one row on each line. \n");
    for(i=1;i<=n;i++)
        for(j=1;j<=n;j++)
            scanf("%f", &a[i][j]);
    printf("\nInput vector b \n");
    for(i=1;i<=n;i++)
        scanf("%f", &b[i]);
/* obtain solution by simple Gauss elimination method */
/* call the subroutine gauss1() */
    gauss1(n,a,b,x,&status);
    if(status != 0)
    {
        printf("\nSOLUTION VECTOR X \n");
    }
}

```

```

        for(i=1;i<=n;i++)
            printf("%10.6f", x[i]);
        printf("\n");
    }
    else
    {
        printf("Singular matrix, reorder equations. \n");
    }
}
/* End of main() program */
/* -----
/* Defining subroutine gauss1() */
gauss1(int n, float a[10][10], float b[10], float x[10],
       int *status)
{
    /* This subroutine solves a set of n linear equations
       by Gauss elimination method */
    int i,j,k;
    float pivot, factor, sum;
    /* ----- Elimination begins ----- */
    for(k=1;k<=n-1;k++)
    {
        pivot = a[k][k];
        if(pivot < 0.000001)
        {
            *status = 0;
            return;
        }
        *status = 1;
        for(i=k+1;i<=n;i++)
        {
            factor = a[i][k] /pivot;
            for(j=k+1;j<=n;j++)
            {
                a[i][j] = a[i][j] - factor * a[k][j];
            }
            b[i] = b[i] - factor * b[k];
        }
    }
    /* ----- Back substitution begins ----- */
    x[n] = b[n] / a[n][n];
    for(k=n-1;k>=1;k--)

```

```

    {
        sum = 0.0;
        for(j=k+1;j<=n;j++)
            sum = sum + a[k][j] * x[j];
        x[k] = (b[k] - sum) / a[k][k];
    }
    return;
}
/* End of subroutine gauss1 () */
/* ----- */

```

Program 11 LEG2

```

/* -----
 * Main program
 * This program solves a system of linear equations
 * using Gaussian elimination with partial pivoting
 * -----
 * Functions invoked
 *     NIL
 * -----
 * Subroutines used
 *     Gauss2
 * -----
 * Variables used
 *     n - Number of equations
 *     a - Coefficients matrix
 *     b - Right side vector
 *     x - Solution vector
 * -----
 * Constants used
 *     NIL
 * -----
main( )
{
    int i,j,n;
    float a[10][10], b[10], x[10];
    printf("\n  GAUSS METHOD WITH PARTIAL PIVOTING \n");
    printf("\nWhat is the size n of the system? \n");
    scanf("%d", &n);
    printf("\nInput coefficients a(i,j), row-wise \n");
    printf("one row on each line \n");
    for(i=1;i<=n;i++)
        for(j=1;j<=n;j++)
            scanf("%f", &a[i][j]);

```

```

printf("\nEnter vector b \n");
for(i=1;i<=n;i++)
    scanf("%f", &b[i]);
gauss2(n,a,b,x);

printf("\n      SOLUTION VECTOR X \n");
printf("\n");
for(i=1;i<=n;i++)
    printf("\t %f", x[i]);
printf("\n");

} /* End of main() program */

/* -----
/* Defining subroutine gauss2() */
gauss2(int n, float a[10][10],float b[10], float x[10])
/* This subroutine solves a system of linear equations using
   Gauss elimination method with partial pivoting */

{
/* Forward elimination */
elim(n,a,b);

/* Solution by back substitution */
bsub(n,a,b,x);

return;
} /* Endof routine gauss2() */

/* -----
/* Defining the subroutine elim() */
elim(int n, float a[10][10],float b[10])
/* This subroutine performs forward elimination
incorporating partial pivoting technique */

{
    int i,j,k;
    float factor;
    for(k=1;k<=n-1;k++)
    {
        pivot(n,a,b,k);
        for(i=k+1;i<=n;i++)
        {
            factor = a[i][k] / a[k][k];
            for(j=k+1;j<=n;j++)
            {

```

```

        a[i][j] = a[i][j]-factor * a[k][j];
    }
    b[i] = b[i]-factor*b[k];
}
}
return;
} /* End of elim() routine */

/* -----
/* Defining subroutine pivot() */
# include <math.h>
pivot(int n, float a[10][10], float b[10], int k)
/* This subroutine performs the task of partial pivoting
   (reordering of equations) */
{
    int p,i,j;
    float large, temp;

/* Find pivot p */
    p = k;
    large = fabs(a[k][k]);
    for(i=k+1;i<=n;i++)
    {
        if(fabs(a[i][k])>large)
        {
            large = fabs(a[i][k]);
            p = i;
        }
    }

/* Exchange rows p and k */
    if(p!=k)
    {
        for(j=k;j<=n;j++)
        {
            temp = a[p][j];
            a[p][j] = a[k][j];
            a[k][j] = temp;
        }
        temp = b[p];
        b[p] = b[k];
        b[k] = temp;
    }
    return;
}
/* End of subroutine pivot() */
/* -----

```

562 Numerical Methods

```
/* Defining subroutine bsub() */
bsub(int n, float a[10][10], float b[10], float x[10])
/* This subroutine obtains the solution vector x by back
substitution */

{
    int i,j,k;
    float sum;
    x[n] = b[n] / a[n][n];
    for(k=n-1;k>=1;k--)
    {
        sum = 0.0;
        for(j=k+1;j<=n;j++)
            sum = sum + a[k][j] * x[j];
        x[k] = (b[k]-sum) / a[k][k];
    }
    return;
}
/* End of subroutine bsub() */
/* ----- */
```

Program 12 DOLIT

```
/* -----
* Main program
*   This program solves a system of linear equations
*   using Dolittle LU decomposition
* -----
* Functions invoked
*   NIL
* -----
* Subroutines used
*   LUD, SOLVE
* -----
* Variables used
*   n - System size
*   a - Coefficient matrix of the system
*   b - Right side vector
*   l - Lower triangular matrix
*   u - Upper triangular matrix
*   fact - Factorization status
* -----
* Constants used
*   YES,NO
* ----- */
```

```

#define YES 1
#define NO 0

main()
{
    int n, fact, i,j;
    float a[10][10], u[10][10], l[10][10], b[10], x[10];
    printf("\n      SOLUTION BY DOLITTLE METHOD \n\n");

/* Read input data */
    printf("\nWhat is size of A? \n");
    scanf("%d", &n);
    printf("Type coefficients a(i,j), row by row \n");
    for(i=1;i<=n;i++)
        for(j=1;j<=n;j++)
            scanf("%f", &a[i][j]);
    printf("\nType vector b on one line \n");
    for(i=1;i<=n;i++)
        scanf("%f", &b[i]);
/* LU factorization */
    lud(n,a,u,l,&fact);
    if(fact == YES)           /* Print LU matrices */
    {
/* Print U matrix */
        printf("\nMATRIX U \n");
        for(i=1;i<=n;i++)
        {
            for(j=1;j<=n;j++)
            {
                printf("%15.6f", u[i][j]);
            }
            printf("\n");
        }
/* Print L matrix */
        printf("\nMATRIX L \n");
        for(i=1;i<=n;i++)
        {
            for(j=1;j<=n;j++)
            {
                printf("%15.6f", l[i][j]);
            }
            printf("\n");
        }
    }
}

```

```

/* Solve for x */
    solve(n,u,l,b,x);
    printf("\nSOLUTION VECTOR X \n\n");
    for(i=1;i<=n;i++)
        printf("%15.6f \n", x[i]);
    printf("\n");
}
else
{
    printf("\n      FACTORIZATION NOT POSSIBLE \n");
}
}

/* End of main() program */

/* -----
/* Defining the subroutine lud() */

lud(int n, float a[10][10], float u[10][10], float l[10][10], int *fact)
{
    /* This subroutine decomposes the matrix A into L and U
       matrices using DOLITTLE algorithm */

    {
        int i,j,k;
        float sum;
    /* Initialize U and L matrices */
        for(i=1;i<=n;i++)
            for(j=1;j<=n;j++)
            {
                u[i][j] = 0.0;
                l[i][j] = 0.0;
            }
    /* Compute the elements of U and L */
        for(j=1;j<=n;j++)
            u[1][j] = a[1][j];
        for(i=1;i<=n;i++)
            l[i][i] = 1.0;
        for(i=1;i<=n;i++)
            l[i][1] = a[i][1] / u[1][1];
        for(j=2;j<=n;j++)
        {
            for(i=2;i<=j;i++)
            {
                sum = a[i][j];
                for(k=1;k<=i-1;k++)
                    sum = sum - l[i][k] * u[k][j];
            }
        }
    }
}

```

```

        u[i][j] = sum;
    }
    if(u[j][j]<=1.E-6)
    {
        *fact = NO;
        return;
    }
    for(i=j+1;i<=n;i++)
    {
        sum = a[i][j];
        for(k=1;k<=j-1;k++)
            sum = sum - l[i][k] * u[k][j];
        l[i][j] = sum / u[j][j];
    }
}
*fact = YES;
return;
} /* End of subroutine lud() */
/*
/* Defining the subroutine solve() */
solve(int n, float u[10][10], float l[10][10], float b[10], float x[10])
/* This subroutine obtains the solution vector X using the
coefficients of L and U matrices */
{
    int i,j;
    float sum, z[10];
/* Forward substitution */
    z[1] = b[1];
    for(i=2;i<=n;i++)
    {
        sum = 0.0;
        for(j=1;j<=i-1;j++)
            sum = sum + l[i][j] * z[j];
        z[i] = b[i] - sum;
    }
/* Back substitution */
    x[n] = z[n]/u[n][n];
    for(i=n-1;i>=1;i--)
    {
        sum = 0.0;
        for(j=i+1;j<=n;j++)
            sum = sum + u[i][j] * x[j];
        x[i] = (z[i] - sum)/u[i][i];
    }
}
```

```

        x[i] = (z[i]-sum)/u[i][i];
    }
    return;
}
/* End of subroutine solve() */
/* -----
 * /
```

Program 13 JACIT

```

/* -----
 * Main program
 * This program uses the subprogram JACOBI to solve
 * a system of equations by Jacobi iteration method
 *
 * -----
 * Functions invoked
 *     NIL
 *
 * -----
 * Subroutines used
 *     JACOBI
 *
 * -----
 * Variables used
 *     a - Coefficient matrix
 *     b - Right side vector
 *     n - System size
 *     x - Solution vector
 *     count - Number of iterations completed
 *     status - Convergence status
 *
 * -----
 * Constants used
 *     EPS - Error bound
 *     MAXIT - Maximum iterations permitted
 * /
```

```

#include <math.h>
#define EPS 0.000001
#define MAXIT 100

main( )
{
    int i,j,n,count,status;
    float a[10][10], b[10], x[10];

    printf("\n      SOLUTION BY JACOBI ITERATION \n");
    printf("\nWhat is the size n of the system? \n");
    scanf("%d", &n);
    printf("\nInput coefficients a(i,j), row by row \n");
    for(i=1;i<=n;i++)
        for(j=1;j<=n;j++)
            /*
```

```

        scanf("%f", &a[i][j]);
scanf("\nInput vector b\n");
for(i=1;i<=n;i++)
    scanf("%f", &b[i]);
jacobi(n,a,b,x, &count, &status);
if(status==2)
{
    printf("\nNo convergence in %d iterations", MAXIT);
    printf("\n\n");
}
else
{
    printf("\n                                     SOLUTION VECTOR X \n\n");
    for(i=1;i<=n;i++)
        printf("%15.6f", x[i]);
    printf("\n\nIterations = %d", count);
}
/* End of main() program */
/* -----
/* Defining the routine jacobi() */
jacobi(int n, float a[10][10], float b[10], float x[10], int
*count, int *status)
/* This subroutine solves a system of n linear equations using
   the Jacobi iteration method */
{
    int i,j,key;
    float sum, x0[10];
/* Initial values of x */
    for(i=1;i<=n;i++)
        x0[i] = b[i] / a[i][i];
/* Jacobi iteration begins */
    *count = 1;
    begin:
    key = 0;
/* Computing values of x(i) */
    for(i=1;i<=n;i++)
    {
        sum = b[i];
        for(j=1;j<=n;j++)
        {

```

```

        if(i==j)
            continue;
        sum = sum - a[i][j] * x0[j];
    }
    x[i] = sum / a[i][i];
    if(key == 0)
    {
/* Testing for accuracy */
    if(fabs((x[i]-x0[i])/x[i]) > EPS)
        key = 1;
    }
/* Testing for convergence */
if(key == 1)
{
    if(*count==MAXIT)
    {
        *status = 2;
        return;
    }
    else
    {
        *status = 1;
        for(i=1;i<=n;i++)
            x0[i] = x[i];
    }
    *count = *count+1;
    goto begin;
}
return;
}
/* End of subroutine jacobi() */
*/

```

Program 14 GASIT

```

/*
 * Main program
 * This program uses the subprogram GASEID to solve a
 * system of equations by Gauss-Seidel iteration method
 *
 * Functions invoked
 *     NIL
 *

```

```

* Subroutines used
* GASEID
*
* -----
* Variables used
*   a - Coefficient matrix
*   b - Right side vector
*   n - System size
*   x - Solution vector
*   count - Number of iterations completed
*   status - Convergence status
*
* -----
* Constants used
*   EPS - Error bound
*   MAXIT - Maximum iterations permitted
* -----
* -
# define MAXIT 50
# define EPS 0.000001

main( )
{
    float a[10][10], b[10], x[10];
    int i,j,n,count,status;
    printf("\n      SOLUTION BY GAUSS-SEIDEL ITERATION \n");
    printf("\nWhat is the size n of the system? \n");
    scanf("%d", &n);
    printf("\nInput coefficients a(i,j), row by row \n");
    printf("one row on each line \n");
    for(i=1;i<=n;i++)
        for(j=1;j<=n;j++)
            scanf("%f", &a[i][j]);
    printf("\nInput vector b\n");
    for(i=1;i<=n;i++)
        scanf("%f", &b[i]);
    gaseid(n,a,b,x, &count, &status);
    if(status==2)
    {
        printf("\nNo convergence in %d iterations \n", MAXIT);
        printf("\n\n");
    }
    else
    {
        printf("\n                      SOLUTION VECTOR X \n\n");
        for(i=1;i<=n;i++)
            printf("%15.6f", x[i]);
        printf("\n\niterations = %d", count);
    }
}

```

```

    }
}

/* End of main() program */
/* -----
/* Defining the routine gaseid() */
gaseid(int n, float a[10][10], float b[10], float x[10], int
       *count, int *status)
/* This subroutine solves a system of linear equations using
   the Gauss-Seidel iteration algorithm */
{
    int i,j,key;
    float sum, x0[10];
/* Initial values of x */
    for(i=1;i<=n;i++)
        x0[i] = b[i] / a[i][i];
/* Gaseid iteration begins */
    *count = 1;
    begin:
    key = 0;
/* Computing values of x(i) */
    for(i=1;i<=n;i++)
    {
        sum = b[i];
        for(j=1;j<=n;j++)
        {
            if(i==j)
                continue;
            sum = sum - a[i][j] * x0[j];
        }
        x[i] = sum / a[i][i];
        if(key==0)
        {
            /* Testing for accuracy */
            if(fabs((x[i]-x0[i])/x[i])>EPS)
                key = 1;
        }
    }
/* Testing for convergence */
    if(key==1)
    {
        if(*count==MAXIT)
        {

```

```

        *status = 2;           /* Program terminated */
        return;
    }
    else
    {
        *status = 1;
        for(i=1;i<=n;i++)
            x0[i] = x[i];
    }
    *count = *count+1;
    goto begin;
}
return;
}
* End of routine gaseid() */
/* ----- */

```

Program 15 LAGRAN

```

/* -----
 * -Main program
 *   This program computes the interpolation value at a
 *   specified point, given a set of data points, using
 *   the Lagrange interpolation representation
 *
 * Functions invoked
 *   NIL
 *
 * Subroutines used
 *   NIL
 *
 * Variables used
 *   n   - Number of data sets
 *   x(i) - Data points
 *   f(i) - Function values at data points
 *   xp  - Point at which interpolation is required
 *   fp  - Interpolated value at XP
 *   lf  - Lagrangian factor
 *
 * Constants used
 *   MAX - Maximum number of data points permitted
 *
#define MAX 10
main( )
{

```

```

int n,i,j;
float x[MAX],f[MAX],fp,lf,sum,xp;

printf("\nInput number of data points, n \n");
scanf("%d", &n);

printf("\nInput data points x(i) and values f(i) \n");
printf("(one set in each line)\n")
for(i=1;i<=n;i++)
    scanf("%f %f", &x[i], &f[i]);

printf("\nInput x at which interpolation
      is required. \n");
scanf("%f", &xp);

sum = 0.0;
for(i=1;i<=n;i++)
{
    lf = 1.0;
    for(j=1;j<=n;j++)
    {
        if(i!=j)
            lf = lf*(xp-x[j])/(x[i]-x[j]);
    }
    sum = sum + lf * f[i];
}
fp = sum;
printf("\n      LAGRANGIAN INTERPOLATION \n\n");
printf("Interpolated function value \n");
printf("at x=%f is %f \n", xp, fp);
}

/* End of main() program */
/* -----
   * Main program
   * This program constructs the Newton interpolation
   * polynomial for a given set of data points and then
   * computes interpolation value at a specified value
   *
   * Functions invoked
   *     NIL
   *
   * -----
   * Subroutines used
   *     NIL
   * -----
*/

```

Program 16 NEWINT

```

/*
   * Main program
   * This program constructs the Newton interpolation
   * polynomial for a given set of data points and then
   * computes interpolation value at a specified value
   *
   * Functions invoked
   *     NIL
   *
   * -----
   * Subroutines used
   *     NIL
   * -----
*/

```

```

* Variables used
*   n - Number of data points
*   x - Array of independent data points
*   f - Array of function values
*   xp - Desired point for interpolation
*   fp - Interpolation value at XP
*   f - Difference table
*   a - Array of coefficients of interpolation
*       polynomial
* -----
* Constants used
*   NIL
* -----
main( )
{
    int i,j,n;
    float xp, fp, sum, pi, x[10], f[10], a[10], d[10][10];
    printf("\nInput number of data points \n");
    scanf("%d", &n);
    printf("\nInput values of x and f(x), \n");
    printf("One set on each line \n");
    for(i=1;i<=n;i++)
        scanf("%f %f", &x[i], &f[i]);
/* Construct difference table */
    for(i=1;i<=n;i++)
        d[i][1] = f[i];
    for(j=2;j<=n;j++)
        for(i=1;i<=n-j+1;i++)
            d[i][j] = (d[i+1][j-1]-d[i][j-1])/(x[i+j-1]-x[i]);
/* Set the coefficients of interpolation polynomial */
/* Set the coefficients of interpolation polynomial */
    for(j=1;j<=n;j++)
        a[j] = d[1][j];
/* Compute interpolation value */
    printf("\nInput xp where interpolation is required \n");
    scanf("%f", &xp);
    sum = a[1];
    for(i=2;i<=n;i++)
    {
        pi = 1.0;
        for(j=1;j<=i-1;j++)
            pi = pi * (xp-x[j]);
        sum = sum + a[i] * pi;
    }
}

```

574 Numerical Methods

```
    }
    fp = sum;

/* Write results */
    printf("\n");
    printf("      NEWTON INTERPOLATION \n");
    printf("\n");
    printf("Interpolated value \n");
    printf("at x = %f is %f \n", xp, fp);
}

/* End of main() program */
/* ----- */
```

Program 17 SPLINE

```
/* -----
 * Main program
 * This program computes the interpolation value at
 * a specified value, given a set of table points,
 * using the natural cubic spline interpolation
 *
 * Functions invoked
 * NIL
 *
 * Subroutines used
 * GAUSS
 *
 * Variables used
 *   n - Number of data points.
 *   x - N by 1 array of data points.
 *   f - N by 1 array of function values
 *   xp - Point at which interpolation is required
 *   fp - Interpolation value at XP
 *   a - Array of second derivatives (N-2 by 1)
 *   d - Array representing (N-2 by 1)
 *   c - Matrix (N-2 by N-2) representing the
 *       coefficients of second derivatives
 *   h - Array of distances between data points
 *       ( h(i) = x(i) - x(i-1) )
 *   df - Array of differences of functions
 *
 * Constants used
 *   MAX - Maximum number of table points permitted
 * ----- */

#define S 10

main( )
{
```

```

int i,j,n,m;
float x[S],f[S],a[S],d[S],c[S][S],h[S],df[S],u[S],
      xp,fp,q1,q2,q3;

/* Reading input data */

printf("\nInput number of data points, n \n");
scanf("%d", &n);

printf("\nInput data points x(i) and function \n");
printf("values f(i), one set on each line \n");
for(i=1;i<=n;i++)
    scanf("%f %f", &x[i],&f[i]);

printf("\nInput point of interpolation \n");
scanf("%f", &xp);

/* Compute distances between data points and function
   differences */

    for(i=2;i<=n;i++)
    {
        h[i] = x[i] - x[i-1];
        df[i] = f[i] - f[i-1];
    }

/* Initialize C matrix */
    for(i=2;i<=n-1;i++)
        for(j=2;j<=n-1;j++)
            c[i][j] = 0.0;

/* Compute diagonal elements of c */
    for(i=2;i<=n-1;i++)
        c[i][i] = 2.0 * (h[i] - h[i+1]);

/* Compute off-diagonal elements of c */
    for(i=3;i<=n-1;i++)
    {
        c[i-1][i] = h[i];
        c[i][i-1] = h[i];
    }

/* Compute elements of d array */
    for(i=2;i<=n-1;i++)
        d[i] = (df[i+1]/h[i+1] - df[i]/h[i]) * 6.0;

/* Compute elements of a using Gaussian elimination.
   Change array subscripts from 2 to n-1 to 1 to n-2
   before calling gauss() subroutine. */

    m = n-2;
    for(i=1;i<=m;i++)
    {

```

576 Numerical Methods

```
d[i] = d[i+1];
for(j=1;j<=m;j++)
    c[i][j] = c[i+1][j+1];
}
gauss(m,c,d,a);

/* Compute the coefficients of natural cubic spline */
for(i=n-1;i>=2;i--)
    a[i] = a[i-1];
a[1] = 0.0;
a[n] = 0.0;

/* Locate the domain of xp */
for(i=2;i<=n;i++)
{
    if(xp <= x[i])
        break;
}

/* Compute interpolation value at xp */
u[i-1] = xp - x[i-1];
u[i] = xp - x[i];
q1 = h[i]*h[i] * u[i] - u[i]*u[i]*u[i];
q2 = u[i-1]*u[i-1]*u[i-1] - h[i]*h[i] * u[i-1];
q3 = f[i]*u[i-1] - f[i-1]*u[i];
fp = (a[i-1]*q1 + a[i]*q2)/(6.0 * h[i]) + q3/h[i];

/* Write results */
printf("\n          SPLINE INTERPOLATION \n \n");
printf("Interpolation value = %f \n", fp);
}

/* End of main() program */

/* ----- */

/* Defining gauss() subroutine */
gauss(int n, float a[10][10], float b[10], float x[10])
/* This subroutine solves a set of n linear equations using
   Gauss elimination method */

{
    int i,j,k;
    float pivot, factor, sum;
}

/* ----- Elimination begins ----- */
for(k=1;k<=n-1;k++)
{
    pivot = a[k][k];
    for(i=k+1;i<=n;i++)
        a[i][k] = a[i][k]/pivot;
    for(i=k+1;i<=n;i++)
        b[i] = b[i]/pivot;
    for(i=k+1;i<=n;i++)
        for(j=k+1;j<=n;j++)
            a[i][j] = a[i][j] - a[i][k]*a[j][k];
}
```

```

    }
    factor = a[i][k] / pivot;
    for(j=k+1;j<=n;j++)
        a[i][j] = a[i][j] - factor * a[k][j];
    b[i] = b[i] - factor * b[k];
}
}

/* Back substitution begins */
x[n] = b[n]/a[n][n];
for(k=n-1;k>=1;k--)
{
    sum = 0.0;
    for(j=k+1;j<=n;j++)
        sum = sum + a[k][j] * x[j];
    x[k] = (b[k]-sum)/a[k][k];
}
return;
}

/* End of subroutine gauss() */
/* ----- */

```

Program 18 LINREG

```

/* -----
* Main program
*   This program fits a line Y = A + BX to a given
*   set of data points by the method least squares
* -----
* Functions invoked
*   Library function fabs()
* -----
* Subroutines used
*   NIL
* -----
* Variables used
*   x,y - Data arrays
*   n - Number of data sets
*   sumx - Sum of x values
*   sumy - Sum of y values
*   sumxx - Sum of squares of x values
*   sumxy - Sum of products of x and y
*   xmean - Mean of x values
*   ymean - Mean of y values
*   a - y intercept of the line
*   b - Slope of the line
* -----
*
```

```

* Constants used
* MAX - Limit for number of data points
* EPS - Error limit
* -----
#include <math.h>
#define MAX 10
#define EPS 0.000001

main( )
{
    int i,n;
    float x[10], y[10];
    float sumx, sumy, sumxx, sumxy, xmean, ymean,
        denom, a, b;

    printf("\n      LINEAR REGRESSION \n \n");

/* Reading data values */
    printf("\nInput number of data points, n \n");
    scanf("%d", &n);
    printf("\nInput x and y values \n");
    printf("One set on each line \n");
    for(i=1;i<=n;i++)
        scanf("%f %f", &x[i], &y[i]);

/* Computing constants a and b of the linear equation */
    sumx = 0.0;
    sumy = 0.0;
    sumxx = 0.0;
    sumxy = 0.0;
    for(i=1;i<=n;i++)
    {
        sumx = sumx + x[i];
        sumy = sumy + y[i];
        sumxx = sumxx + x[i] * x[i];
        sumxy = sumxy + x[i] * y[i];
    }
    xmean = sumx/n;
    ymean = sumy/n;
    denom = n*sumxx - sumx * sumx;
    if(fabs(denom) > EPS)
    {
        b = (n*sumxy - sumx * sumy)/denom;
        a = ymean - b * xmean;
        printf("\n      LINEAR REGRESSION LINE y = a+bx \n");
        printf("\nThe coefficients are: \n");
    }
}

```

```

        printf("a = %f \n", a);
        printf("b = %f \n", b);
    }
    else
    {
        printf("\n NO SOLUTION \n");
    }
}
/* End of main() program */
/* ----- */

```

Program 19 POLREG

```

/* -----
 * Main program
 *   This program fits a polynomial curve to a given
 *   set of data points by the method of
 *   least squares
 *
 * -----
 * Functions invoked
 *   NIL
 *
 * -----
 * Subroutines used
 *   normal, gauss
 *
 * -----
 * Variables used
 *   x,y - Arrays of data values
 *   n - Number of data points
 *   mp - Order of the polynomial under construction
 *   m - Number of polynomial coefficients
 *   c - Coefficient matrix of normal equations
 *   b - Right side vector of normal equations
 *   a - Array of coefficients of the polynomial
 *
 * -----
 * Constants used
 *   MAX - Maximum number of data points
 *
 * -----
 * /

```

```

#include <math.h>
#define MAX 10

main( )
{
    int n,mp,m,i;
    float x[MAX],y[MAX],c[MAX][MAX],a[MAX],b[MAX];
    printf("\n      POLYNOMIAL REGRESSION \n\n");

```

580 Numerical Methods

```
/* Reading values */
    printf("Input number of data points n \n");
    scanf("%d", &n);
    printf("Input order of polynomial, mp required \n");
    scanf("%d", &mp);
    printf("Input data values, x and y \n");
    printf("one set on each line \n");
    for(i=1;i<=n;i++)
        scanf("%f %f", &x[i], &y[i]);

/* Testing the order */
    if(n <= mp)
    {
        printf("\n REGRESSION IS NOT POSSIBLE \n");
        goto stop;
    }

/* Number of polynomial coefficients */
    m = mp + 1;

/* Computation of elements of c and b */
    normal(x,y,c,b,n,m);

/* Computation of coefficients a(1) to a(m) */
    gauss(m,c,b,a);

/* Printing of coefficients a(i) */
    printf("\n      POLYNOMIAL COEFFICIENTS \n\n");
    for(i=1;i<=m;i++)
        printf("%15.6f", a[i]);
    printf("\n");
    stop:
    printf("END");

}
/* End of main() program */
/* -----
/* Defining the subroutine normal() */
normal(float x[MAX], float y[MAX], float c[MAX][MAX],
       float b[MAX], int n, int m)

/* This subroutine computes the coefficients of normal
   equations */
{
    int i,j,k,l1,l2;
    for(j=1;j<=m;j++)
    {
        for(k=1;k<=m;k++)
            l1 =
```

```

    {
        c[j][k] = 0.0;
        l1 = k+j-2;
        for(i=1;i<=n;i++)
            c[j][k] = c[j][k] + pow(x[i],l1);
    }
    for(j=1;j<=m;j++)
    {
        b[j] = 0.0;
        l2 = j-1;
        for(i=1;i<=n;i++)
            b[j] = b[j]+y[i] * pow(x[i],l2);
    }
    return;
}
/* End of subroutine normal() */
/* -----
/* Defining gauss() subroutine */
gauss(int n, float a[MAX][MAX], float b[MAX], float x[MAX])
/* This subroutine solves a set of n linear equations by Gauss
   Elimination method */
{
    int i,j,k;
    float pivot, factor, sum;
/* Elimination begins */
    for(k=1;k<=n-1;k++)
    {
        pivot = a[k][k];
        for(i=k+1;i<=n;i++)
        {
            factor = a[i][k]/pivot;
            for(j=k+1;j<=n;j++)
                a[i][j] = a[i][j] - factor * a[k][j];
            b[i] = b[i] - factor * b[k];
        }
    }
/* Back substitution begins */
    x[n] = b[n]/a[n][n];
    for(k=n-1;k>=1;k--)
    {
        sum = 0.0;
        for(j=k+1;j<=n;j++)

```

```

        sum = sum + a[k][j] * x[j];
        x[k] = (b[k]-sum)/a[k][k];
    }
    return;
}
/* End of subroutine gauss() */
*/

```

Program 20 NUDIF

```

/* -----
 * Main program
 * This program computes the derivative of a tabulated
 * function at a specified value using the Newton
 * interpolation approach
 *
 * Functions invoked
 * NIL
 *
 * -----
 * Subroutines used
 * NIL
 *
 * -----
 * Variables used
 * n - Number of function values given
 * x - Array of values
 * f - Array of function values
 * a - Array of coefficients of Newton polynomial
 * d - Difference table
 * xp - Desired point of differentiation
 * dif - Derivative of the function at XP
 *
 * Constants used
 * NIL
 * -----
 */

main( )
{
    int i,j,k,n;
    float x[10],f[10],a[10],d[10][10],xp,dif,sum,p;

    /* Reading input data */
    printf("\nInput number of data points \n");
    scanf("%d", &n);
    printf("\nInput values of x and f(x), \n");
    printf("one set on each line \n");
    for(i=1;i<=n;i++)
        scanf("%f %f", &x[i], &f[i]);

```

```

/* Constructing difference table d */
    for(i=1;i<=n;i++)
        d[i][1] = f[i];
    for(j=2;j<=n;j++)
        for(i=1;i<=n-j+1;i++)
            d[i][j] = (d[i+1][j-1] - d[i][j-1])/(x[i+j-1] -
x[i]);
/* Set the coefficients of Newton interpolating polynomial */
    for(j=1;j<=n;j++)
        a[j] = d[1][j];
/* Compute derivative at a given x = xp */
    printf("\nInput xp where derivative is required \n");
    scanf("%f", &xp);
    dif = a[2];
    for(k=3;k<=n;k++)
    {
        sum = 0.0;
        for(i=1;i<=k-1;i++)
        {
            p = 1.0;
            for(j=1;j<=k-1;j++)
            {
                if(i == j)
                    continue;
                p = p * (xp - x[j]);
            }
            sum = sum + p;
        }
        dif = dif + a[k] * sum;
    }
/* Write results */
    printf("\nNUMERICAL DIFFERENTIATION USING ");
    printf("NEWTON POLYNOMIAL \n\n");
    printf("DERIVATIVE at x = %f is %f \n", xp,dif);
}
/* End of main() program */
/* -----

```

Program 21 TRAPE1

```

/* -----
 * Main program
 * This program integrates a given function
 * using the trapezoidal rule
 * -----

```

584 Numerical Methods

```
* Functions invoked
* NIL
*
* -----
* Subroutines used
* F(x)
*
* -----
* Variables used
*     a - Lower limit of integration
*     b - Upper limit of integration
*     h - Segment width
*     n - Number of segments
*     ict - Value of integral
*
* -----
* Constants used
* NIL
* -----
*/
```

```
#include <math.h>

Main( )
{
    int n,i;
    float a,b,h,sum,ict;
    float F(float x);

    printf("\nGive initial value of x \n");
    scanf("%f",&a);
    printf("\nGive final value of x \n");
    scanf("%f", &b);
    printf("\nWhat is the segment width? \n");
    scanf("%f", &h);

    n = (b-a)/h;
    sum = (F(a) + F(b))/2.0;
    for(i=1;i<=n-1;i++)
    {
        sum = sum + F(a+i*h);
    }
    ict = sum * h;

    printf("\n");
    printf("Integration between %f and %f \n", a,b);
    printf("When h = %f is %f \n", h, ict);
    printf("\n");
}

/* End of main() program */
/* ----- */
```

```

float F(float x)
{
    float f;
    f = 1.0-exp(-x/2.0);
    return (f);
}

/* End of subroutine F(x) */

/* -----
   * Main program
   * This program integrates a given function
   * using the Simpson's 1/3 rule
   *
   * Functions invoked
   * Macro F(x)
   *
   * -----
   * Subroutines used
   * NIL
   *
   * -----
   * Variables used
   * a - Lower limit of integration
   * b - Upper limit of integration
   * h - Segment width
   * n - Number of segments
   * ics - Value of the integral
   *
   * -----
   * Constants used
   * NIL
   *
   * -----
   */

```

Program 22 SIMS1

```

#include <math.h>
#define F(x) 1 - exp(-(x)/2.0)
main( )
{
    int n,m,i;
    float a,b,h,sum,ics,x,f1,f2,f3;
    printf("\nInitial value of x \n");
    scanf("%f", &a);
    printf("\nFinal value of x \n");
    scanf("%f", &b);
    printf("\nNumber of segments (EVEN number) \n");
    scanf("%d", &n);

```

586 Numerical Methods

```
h = (b-a)/n;
m = n/2;

sum = 0.0;
x = a;
f1 = F(x);
for(i=1;i<=m;i++)
{
    f2 = F(x+h);
    f3 = F(x+2*h);
    sum = sum + f1 + 4*f2 + f3;
    f1 = f3;
    x = x + 2*h;
}
ics = sum * h/3.0;
printf("\nIntegral from %f to %f \n", a,b);
printf("When h = %f is %f \n", h, ics);
}

/* End of main() program */
/* --- */
```

Program 23 ROMBRG

```
/* -----
 * Main program
 * This program performs Romberg integration
 * by bisecting the intervals N times
 *
 * -----
 * Functions invoked
 * Macro F(x), Library function fabs()
 *
 * -----
 * Subroutines used
 * NIL
 *
 * -----
 * Variables used
 * a - Starting point of the interval
 * b - End point of the interval
 * h - Width of the interval
 * n - Number of times bisection is done
 * m - Number of trapezoids
 * r - Matrix of Romberg integral values
 *
 * -----
 * Constants used
 * EPS - Error bound
 * ----- */
```

```

#include <math.h>
#define EPS 0.000001
#define F(x) 1.0/(x)

main( )
{
    int i,j,k,m,n;
    float a,b,h,sum,x,r[10][10];

    printf("\nInput end points of the interval \n");
    scanf("%f %f", &a, &b);
    printf("\nInput maximum number of times");
    printf("\nthe subintervals are bisected \n");
    scanf("%d", &n);

/* Compute area using entire interval as one trapezoidal */
    h = b - a;
    r[1][1] = h * (F(a)+F(b))/2.0;
    printf("\n%15.6f \n", r[1][1]);

/* Process of Romberg integration begins */
    for(i=2;i<=n+1;i++)
    {
        m = pow(2,(i-2)); /* trapezoidal for ith refinement */
        h = h/2;           /* bisect step-size */

/* Use recursive trapezoidal rule for m strips */
        sum = 0.0;
        for(k=1;k<=m;k++)
        {
            x = a + (2*k-1)*h;
            sum = sum + F(x);
        }
        r[i][1] = r[i-1][1]/2.0 + h*sum;

/* Compute Richardson's improvements */
        for(j=2;j<=i;j++)
        {
            r[i][j]=r[i][j-1]+(r[i][j-1]-r[i-1][j-1])/(pow(4,j-1)-1);
        }

/* Write results of improvements for ith refinement */
        for(j=1;j<=i;j++)
            printf("%15.6f", r[i][j]);
        printf("\n");

/* Test for accuracy */
        if(fabs(r[i-1][i-1] - r[i][i]) < EPS)
/* Stop further refinement */
        {

```

```

    printf("\n");
    printf("ROMBERG INTEGRATION = %f \n", r[i][i]);
    printf("\n");
    goto stop;
}
else
{
    continue;
}
}

/* Write final result */
printf("\nROMBERG INTEGRATION = %f \n", r[n+1][n+1]);
printf("(Normal exit from loop) \n");

stop:
printf("End");
}

/* End of main() program */
/* ----- */

```

Program 24 TRAPEZOIDAL RULE

```

/* -----
 * Main program
 * This program integrates a tabulated function
 * using the trapezoidal rule
 * -----
 * Functions invoked
 * Library function fabs()
 * -----
 * Subroutines used
 * NIL
 * -----
 * Variables used
 * n - Number of table points
 * x - Array of independent data points
 * y - Array of function values
 * a - Lower limit of integration
 * b - Upper limit of integration
 * h - Distance between points
 * n1 - Position of A in the table
 * n2 - Position of B in the table
 * ict - Value of integral
 * -----
 * Constants used
 * NIL
 * -----
 * /

```

```

#include <math.h>
#define MAX 15

main( )
{
    int n,n1,n2,i;
    float a,b,h,sum,ict,x[MAX],y[MAX];

/* Reading table values */
    printf("Number of data points \n");
    scanf("%d", &n);
    printf("\nInput table values, set by set \n");
    for(i=1;i<=n;i++)
        scanf("%f %f", &x[i],&y[i]);

/* Reading the limits of integration */
    printf("Initial value of x \n");
    scanf("%f",&a);
    printf("Final value of x \n");
    scanf("%f", &b);
    printf("\nWhat is segment width? \n");
    scanf("%f", &h);

/* Computing the position of initial and final values */
    n1 = (int)(fabs(a-x[1])/h+1.5);
    n2 = (int)(fabs(b-x[1])/h+1.5);

/* Evaluating the integral */
    sum = 0.0;
    for(i=n1;i<=n2-1;i++)
        sum = sum + y[i] + y[i+1];
    ict = sum * h/2.0;
    printf("\nIntegral from %f to %f is %f\n", a,b,ict);
}

/* End of main() program */

/* -----
 * Main program
 * This program integrates a tabulated function
 * using the Simpson's 1/3 rule. If the number of
 * segments is odd, the trapezoidal rule is used
 * for the last segment.
 *
 * Functions invoked
 * Library function fabs()
 *
 */

```

Program 25 SIMS2

```

/* -----
 * Main program
 * This program integrates a tabulated function
 * using the Simpson's 1/3 rule. If the number of
 * segments is odd, the trapezoidal rule is used
 * for the last segment.
 *
 * Functions invoked
 * Library function fabs()
 *
 */

```

```

* Subroutines used *
* NIL
* -----
* Variables used
*   n - Number of table points
*   x - Array of independent values
*   y - Array of function values
*   a - Lower limit of integration
*   b - Upper limit of integration
*   h - Distance between independent values
*   n1 - Position of A in the table
*   n2 - Position of B in the table
*   i1 - Area computed using Simpson's rule
*   i2 - Area of the last segment (by trapezoidal rule)
*   ics - Value of integral
* -----
* Constants used
* NIL
* -----
* /

```

```

#include <math.h>
main( )
{
    int i,n,n1,n2,m,l;
    float x[15],y[15],a,b,h,sum,ics,i1,i2;
/* Reading table values */
    printf("Input number of data points \n");
    scanf("\%d", &n);
    printf("\nInput table values, set by set \n");
    for(i=1;i<=n;i++)
        scanf("\%f \%f", &x[i],&y[i]);
/* Reading the limits of integration */
    printf("Input initial value of x \n");
    scanf("\%f", &a);
    printf("Input final value of x \n");
    scanf("\%f", &b);
    printf("What is segment width? \n");
    scanf("\%f", &h);
/* Computing the position of initial and final values */
    n1 = (int)(fabs(a-x[1])/h+1.5);
    n2 = (int)(fabs(b-x[1])/h+1.5);
/* Testing for even intervals */
    m = n2 - n1;
    if(m&2 == 0) /* m is even */

```

```

    {
        i2 = 0.0;
        l = n2 - 2;
    }
    else /* m is odd */
    {
        /* Use trapezoidal rule for the last strip */
        i2 = (y[n2-1]+y[n2])*h/2.0;
        l = n2 - 3;
    }

/* Use Simpson's rule for l strips */
    sum = 0.0;
    for(i=n1;i<=l;i=i+2)
        sum = sum + y[i] + 4*y[i+1] + y[i+2];
    i1 = sum*h/3.0;

/* Integral is sum of i1 and i2 */
    ics = i1 + i2;

/* Writing the results */
    printf("\nIntegral from %f to %f is %f \n", a,b,ics);
}
/* End of main() program */
/* -----

```

Program 26 EULER

```

/* -----
 * Main program
 * This program estimates the solution of the first
 * order differential equation  $y' = f(x,y)$  at a given
 * point using Euler's method
 * -----
 * Functions invoked
 *     NIL
 * -----
 * Subroutines used
 *     func()
 * -----
 * Variables used
 *     x - Initial value of independent variable
 *     y - Initial value of dependent variable
 *     xp - Point of solution
 *     h - Incremental step-size
 *     n - Number of computational steps required
 *     dy - Incremental Y in each step
 * -----

```

```

* Constants used
* NIL
* -----
* /



main( )
{
    int i,n;
    float x,y,xp,h,dy;
    float func(float, float);

    printf("\n      SOLUTION BY EULER'S METHOD \n \n");

/* Reading initial data */
    printf("Input initial values of x and y \n");
    scanf("%f %f", &x,&y);
    printf("Input x at which y is required \n");
    scanf("%f", &xp);
    printf("Input step-size, h \n");
    scanf("%f", &h);

/* Compute number of steps required */
    n = (int)((xp-x)/h+0.5);

/* Compute y recursively at each step */
    for(i=1;i<=n;i++)
    {
        dy = h * func(x,y);
        x = x + h;
        y = y + dy;
        printf("%5d %10.6f %10.6f \n", i,x,y);
    }

/* Write the final result */
    printf("\nValue of y at x = %f is %f \n", x,y);
}

/* End of main() program */

/* -----
* /



/* Defining subroutine func() */
float func(float x, float y)
{
    float f;
    f = 2.0 * y/x;
    return(f);
}

/* End of subroutine func() */
/* -----
* /

```

Program 27 HEUN

```

/* -----
 * Main program
 * This program solves the first order differential
 * equation y' = f(x,y) using the Heun's method
 *
----- */
* Functions invoked
* NIL
*
----- */
* Subroutines used
* func()
*
----- */
* Variables used
* x - Initial value of independent variable
* y - Initial value of dependent variable
* xp - Point of solution
* h - Step-size
* n - Number of steps
*
----- */
* Constants used
* NIL
* -----
 */

main( )
{
    int i,n;
    float x,y,xp,h,m1,m2;
    float func(float, float);

    printf("\n      SOLUTION BY HEUN'S METHOD \n \n");

/* Reading initial data */
    printf("Input initial values of x and y \n");
    scanf("%f %f", &x,&y);
    printf("Input x at which y is required \n");
    scanf("%f", &xp);
    printf("Input step-size, h \n");
    scanf("%f", &h);

/* Compute number of steps required */
    n = (int)((xp-x)/h+0.5);

/* Compute y recursively at each step */
    for(i=1;i<=n;i++)
    {
        m1 = func(x,y);
        m2 = func(x+h,y+m1*h);
        x = x + h;
        y = y + 0.5*h*(m1+m2);
    }
}

```

594 Numerical Methods

```
        printf("%5d %10.6f %10.6f \n", i,x,y);
    }

/* Write the final result */
printf("\nValue of y at x = %f is %f \n", x,y);
}

/* End of main() program */

/* -----
/* Defining subroutine func() */
float func(float x, float y)
{
    float f;
    f = 2.0 * y/x;
    return(f);
}
/* End of subroutine func() */
/* ----- */
```

Program 28 POLYGN

```
/* -----
* Main program
*   This program solves the differential equation
*   of type  $y' = f(x,y)$  by polygon method
* -----
* Functions invoked
*   NIL
* -----
* Subroutines used
*   func()
* -----
* Variables used
*   x - Initial value of the independent variable
*   y - Initial value of the dependent variable
*   xp - Point of solution
*   h - Incremental step-size
*   n - Number of computational steps required
* -----
* Constants used
*   NIL
* ----- */

main( )
{
    int i,n;
    float x,y,xp,h,m1,m2;
```

```

float func(float, float);
printf("\n SOLUTION BY POLYGON METHOD \n \n");
/* Reading initial data */
printf("Input initial values of x and y \n");
scanf("%f %f", &x,&y);
printf("Input x at which y is required \n");
scanf("%f", &xp);
printf("Input step-size, h \n");
scanf("%f", &h);

/* Compute number of steps required */
n = (int)((xp-x)/h+0.5);

/* Compute y recursively at each step */
for(i=1;i<=n;i++)
{
    m1 = func(x,y);
    m2 = func(x+0.5*h,y+0.5*h*m1);
    x = x + h;
    y = y + m2*h;
    printf("%5d %10.6f %10.6f \n", i,x,y);
}

/* Write the final result */
printf("\nValue of y at x = %f is %f \n", x,y);

}
/* End of main() program */
/* -----
/* Defining subroutine func() */
float func(float x, float y)
{
    float f;
    f = 2.0 * y/x;
    return(f);
}
/* End of subroutine func() */
/* -----

```

Program 29 RUNGE4

```

/* -----
 * Main program
 * This program computes the solution of first order
 * differential equation of type y' = f(x,y) using
 * the 4th order Runge-Kutta method
 * -----

```

```

* Functions invoked
* Nil
* -----
* Subroutines used
* func()
*
* -----
* Variables used
*     x - Initial value of independent variable
*     y - Initial value of dependent variable
*     xp - Point of solution
*     h - Step-size
*     n - Number of steps
*
* -----
* Constants used
* NIL
* -----
* /
main( )
{
    int i,n;
    float x,y,xp,h,m1,m2,m3,m4;
    float func(float, float);

    printf("\n . SOLUTION BY 4th ORDER RK METHOD\n\n");

/* Input initial data */
    printf("Input initial values of x and y \n");
    scanf("%f %f", &x,&y);
    printf("Input x at which y is required \n");
    scanf("%f", &xp);
    printf("Input step-size, h \n");
    scanf("%f", &h);

/* Compute number of steps required */
    n = (int)((xp-x)/h+0.5);

/* Compute y at each step */
    printf("\n");
    printf("-----\n");
    printf(" STEP      X      Y      \n");
    printf("-----\n");
    for(i=1;i<=n;i++)
    {
        m1 = func(x,y);
        m2 = func(x+0.5*h, y+0.5*m1*h);
        m3 = func(x+0.5*h, y+0.5*m2*h);
        m4 = func(x+h, y+m3*h);
    }
}

```

```

        x = x+h;
        y = y + (m1+2.0*m2 + 2.0*m3 + m4) * h/6.0;
        printf("%5d %15.6f %15.6f \n", i,x,y);
    }
    printf(" ----- \n");

/* Write the final value of y */
    printf("\nValue of y at x = %f is %f \n", x,y);

}
/* End of main() program */

/* -----
/* Defining subroutine func() */
float func(float x, float y)
{
    float f;
    f = 2.0 * y/x;
    return(f);
}
/* End of subroutine func() */

* -----

```

Program 30 MILSIM

```

/* -----
* Main program
* This program solves the first order differential
* equation y' = f(x,y) using Milne-Simpson method
* -----
* Functions invoked
* NIL
* -----
* Subroutines used
* func()
* -----
* Variables used
* x(1) - Initial value of independent variable
* y(1) - Initial value of dependent variable
* xp - Point of solution
* n - Number of steps
* h - Step-size
* x - Array of independent variable
* y - Array of dependent variable
* -----
* Constants used
* NIL
* -----

```

```

main( )
{
    int i,n;
    float h,x[15],y[15],xp,m1,m2,m3,m4,f2,f3,f4,f5;
    float func(float, float);

/* Reading initial values */
    printf("Input initial values of x and y \n");
    scanf("%f %f", &x[1],&y[1]);
    printf("Input value of x at which y is required \n");
    scanf("%f", &xp);
    printf("Input step-size, h \n");
    scanf("%f", &h);

/* Compute number of computation required */
    n = (int)((xp-x[1])/h+0.5);

/* We require four starting points for Milne-Simpson
method.
Initial
values form the first point. Remaining three points are
obtained
using 4th order RK method */

    printf("\nINITIAL VALUES ARE %10.6f %10.6f", x[1],y[1]);

/* Computing three more points by RK method */
    printf("\n\nTHREE VALUES BY RK METHOD \n");
    for(i=1;i<=3;i++)
    {
        m1 = func(x[i],y[i]);
        m2 = func(x[i]+0.5*h, y[i]+0.5*m1*h);
        m3 = func(x[i]+0.5*h, y[i]+0.5*m2*h);
        m4 = func(x[i]+h, y[i]+m3*h);
        x[i+1] = x[i] + h;
        y[i+1] = y[i] + (m1+2.0*m2 + 2.0*m3 + m4)*h/6.0;
        printf("\n%5d %10.6f %10.6f", i, x[i+1],y[i+1]);
    }

/* Computing values by Milne-Simpson method */
    printf("\n\nVALUES OBTAINED BY MILNE-SIMPSON METHOD \n");
    for(i=4;i<=n;i++)
    {
        f2 = func(x[i-2], y[i-2]);
        f3 = func(x[i-1], y[i-1]);
        f4 = func(x[i], y[i]);

        /* Predicted value of y (by Milne's formula)*/
        y[i+1] = y[i-3] + 4.0*h/3.0 * (2.0*f2-f3+2.0*f4);
        x[i+1] = x[i] + h;
        f5 = func(x[i+1], y[i+1]);
    }
}

```

```
/* Corrected value of y (by Simpson's formula) */
y[i+1] = y[i-1]+h/3.0 * (f3+4.0*f4+f5);
printf("\n%5d %10.6f %10.6f", i, x[i+1], y[i+1]);
}
printf("\n\nValue of y at x = %f is %f", x[n+1], y[n+1]);
}
/* End of main() program */

/* -----
/* Defining subroutine func()
float func(float x, float y)
{
    float f;
    f = 2.0 * y/x;
    return(f);
}
/* End of subroutine func() */
/* ----- */
```

APPENDIX

E

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Index

- Abacus 14
Absolute error 66, 70
Accuracy 8, 63
Adams-Basforth predictor 450
Adams-Moulton corrector 450
Adjoint matrix 509
Algebraic equation 122
Algebraic expression 102
Algorithm 7, 30, 80
 Bairstow 186
 Basic Gauss elimination 213
 Bisection 133
 Cholesky 241
 Crout 240
 Dolittle 231, 233
 False position 139
 Fixed-point 169
 Gaussian integration 401
 Gauss-Jordan 228
 Gauss-Seidel 261
 Horner's rule 128
 Incremented search 167
 Iterative refinement 243
 Jacobi 254
 Linear regression 327
 Muller 196
 Multiple roots 178
 Natural cubic spline 307
 Newton-Raphson 146, 171
 Partial pivoting 220
 Polynomial fit 336
 Polynomial method 477
 Power method 479
 Secant 153
 Shooting method 469
Analog computer 2, 17
Analytical method 123
Application programs 26
Approximating functions 276
Arguments 109, 110
Arithmetic logic unit 22
Arithmetic operators 102
Array 99
Assembler 28
Assembly language 27
Assignment statement 102
Associative law 57
Asymptotic convergence factor 86
Back substitution 212
Backward difference derivative 361
Backward difference quotient 348
Backward difference table 298
Backward difference 298
Backward error analysis 87
Backward jump 35, 103
Bairstow's method 183
Base 41
Basic Gauss elimination 212
Bender-Schmidt equation 494
Binary chopping method 131
Binary code 27
Binary digits 40
Binary numbers 40

- Binary system 42
 Bisection algorithm 133
 Bisection method 131
 Bits 49
 Blunder 69
 Boole's rule 387
 Boundary value problem 412, 464
 Bracketing method 124
 Branching structure 33
 Bugs 113
 Built-in-functions 112
 C 9
 C++ 9
 Calling program 109, 110
 Calling statement 109, 110
 Central difference 512
 Central difference derivative 361
 Central difference operator 521
 Central difference quotient 349
 Central processing unit 22
 Character constants 98
 Character type 98
 Characteristic equation 472
 Characteristic matrix 472
 Characteristic polynomial 472
 Characteristic values 472
 Chebyshev basis polynomial 315
 Chebyshev interpolation 314
 Chebyshev points 314
 Chebyshev polynomial 315
 Cholesky's factorisation 241
 Cholesky's algorithm 241
 Chopping 65
 Closed form 371
 Coefficient matrix 244, 246, 511
 Cofactor 508
 Column vector 506
 Comment line 95
 Compiler 26, 28
 Complete pivoting 220
 Complex constants 98
 Complex number
 Complex root 173, 176, 183, 199
 Complex type 97
 Composite approach 374
 Composite Simpson's 1/3 rule 382
 Composite trapezoidal rule 374
 Computer memory 40
 Computer program 20, 24, 27
 Computer-aided design 13
 Computer-aided learning 13
 Computer-aided manufacture 13
 Computer-managed learning 13
 Condition number 81, 246, 247
 Conditional execution 103
 Conditioning 80
 Constants 97
 Continuous data 2
 Control statement 96
 Control unit 22
 Control variable 106
 Convergence 86
 Bisection method 134
 False position method 140
 Fixed-point method 163
 Gauss-Seidel 267
 Jacobi method 267
 Newton-Raphson 147
 Secant method 153
 Conversion error 64
 Corrector 427
 Counter 96
 Cramer's rule 509
 Crank-Nicholson formula 496
 Cross average 489
 Crout algorithm 240
 Crout LU decomposition 231
 Cubic spline 302
 Curve fitting 275, 323
 Data 20, 40
 Data error 64
 DATA statement 101
 Debugging 26, 113
 Decimal numbers 41
 Decimal system 41
 Decomposition 231
 Deflation 175
 Degree of equations 410
 Dependent equations 209
 Dependent variable 323, 410
 Descartes rule 174
 Determinant 507
 Diagonal average 489
 Diagonally dominant system 268
 Difference equation 484
 Difference tables 360
 Differential equations 408
 Digital computer 3, 17
 Dimension 100
 Direct method 207

- Direct substitution method 161
 Dirichlet boundary condition 488
 Discrete data 2
 Discretisation error 67
 Display screen 20
 Distributive law 57
 Divided difference table 290
 Divided differences 288
 DO loop 106
 DO..WHILE structure 106
 Documentation 114
 Dolittle algorithm 231
 Dolittle LU decomposition 231
 Double precision 99
 Drag coefficient 69
 Driver program 109
 Dummy variables 109, 110

 Efficiency 9
 Eigenvalue problem 465
 Eigenvalues 472, 475
 Eigenvector 472, 476
 Elimination approach 207
 Elliptic equation 484, 486
 Empirical error 64
 ENIAC 15
 Equal matrix 506
 Error propagation 73
 Euler's method 419
 Euler-Maclaurin formula 388
 Executable statement 96
 Experimental error analysis 88
 Expert systems 16
 Explicit method 495
 Exponent 51
 Exponential form 98
 Expressions 102
 Extrapolation method 124
 Extrapolation 362, 388

 Factorisation 230
 Fadeev-Leverrier method 474
 False position algorithm 139
 False position formula 138
 False position method 138
 Fifth generation 16
 Finite difference method 464, 469,
 484
 Finite-element method 484
 First generation 16
 Five-point formula 350

 Fixed loop 34
 Fixed point arithmetic 52
 Fixed point equation 161
 Fixed point method 160, 168
 Floating point arithmetic 53
 Floating point form 50, 98
 Floating point number 51
 Floppy disk 23
 Flowchart 32, 103, 107
 FORMAT statement 101
 Format-directed I/O statement 100
 FORTRAN 9
 FORTRAN expression 102
 Forward difference 295
 Forward difference derivative 360
 Forward difference quotient 348
 Forward difference table 296
 Forward difference 295
 Forward elimination 211, 212
 Forward error analysis 87
 Forward jump 34, 103
 Fourth generation 16
 Functions 109

 Gauss elimination 212, 219, 228
 Gaussian integration 394
 Gaussian quadrature 395
 Gauss-Jordan method 228
 Gauss-Legendre formula 396
 Gauss-Legendre rules 371
 Gauss-Seidel iteration 259, 489
 General purpose computer 3, 18
 General solution 411
 Global truncation error 421
 Graeffe's method 516
 Graphical method 123

 Half-interval method 131
 Hard disk 24
 Hardcopy 20
 Hardware 21
 Heat equation 493
 Heun's method 425
 Hexadecimal numbers 40, 42
 Hexadecimal system 42
 High-level language 28
 Homogeneous equations 209
 Horner's rule 129
 Human error 69
 Hybrid computer 17
 Hyperbolic equation 484, 497

- IC chips 16
 Identity matrix 246, 472, 505
 IF...ELSE structure 103
 Ill-conditioned problem 9, 81
 Ill-conditioned system 209, 244
 Ill-conditioning 342
 Implicit method 496
 Inconsistent equations 208
 Incremental search 167
 Independent variable 323, 410
 Induced errors 80
 Induced instability 85
 Infinite solutions 208
 Information 20
 Inherent error 64
 Inherently unstable 81
 Initialisation 96
 Initial-value problem 412, 464
 Inner loop 108
 Input devices 21
 Input error 64
 Input statement 100
 Integer arithmetic 52
 Integer constants 98
 Integer form 50
 Integer type 97
 Integration nodes 370
 Interactive computing 30
 Interpolation 276
 Interpolation function 281
 Interpolation method 124
 Interpolation polynomial 276
 Interpreter 26
 Interval arithmetic 52
 Intrinsic functions 112
 Inverse matrix 509
 Iterative method 123, 124
 Iterative process 8
 Iterative refinement 242
 Jacobi iteration 253
 Kirchhoff's law 409
 Knots 300, 312
 Knowledge-based systems 16
 Lagrange basis polynomial 282
 Lagrange interpolation 281
 Lagrange interpolation polynomial 282, 371
 Language processors 26
 Laplace's equation 486
 Laplacian operator 486
 Large-scale integration(LSI) 16
 Law of cooling 408
 Law of motion 408
 Least-squares polynomials 276
 Least-squares regression 276, 324, 325
 Liebmann's method 489
 Linear equation 206, 411
 Linear fitting 324
 Linear function 121
 Linear interpolation 138, 279
 Linear interpolation formula 280, 282
 Linear system 168
 Linearly convergent 134, 140, 164
 List-directed I/O statement 100
 Local truncation error 421
 Logarithm 14
 Logic errors 114
 Logical constants 98
 Logical expression 105
 Logical operators 105
 Logical type 98
 Looping structure 34, 106
 Low level language 27
 Lower triangular matrix 231
 LU decomposition 231, 240
 Machine code 27
 Machine epsilon 71
 Machine language 26, 27
 Main program 109
 Mainframe computer 18
 Mantissa 51
 Mark 114
 Mathematical model 3
 Matrix inversion 245
 Matrix norm 247
 Mean-value theorem 79
 Memory unit 22
 Method of square roots 241
 Microcomputer 18
 Microprocessor 18
 Midpoint method 433
 Milne's formula 445
 Milne-Simpson method 445, 451
 Minicomputer 18
 Minor 508
 Mixed-mode expression 102

- Modelling 13
 Modelling error 69
 Modifiers 452
 Modular programming 31, 109
 Modular structure 222
 Modules 109
 Monotone convergence 163
 Monotone divergence 163
 Muller's method 195
 Multidimensional array 99
 Multiple roots 174, 178, 199
 Multisegment approach 374
 Multistep methods 412, 444

 Named constant 96
 Natural cubic spline 306
 Nesting 107
 Newton form 278, 287
 Newton interpolation polynomial 288, 371
 Newton's interpolation 286
 Newton's three-eighths rule 385
 Newton-Cotes formula 371
 Newton-Cotes rules 371, 388
 Newton-Gregory formula 296, 298
 Newton-Gregory interpolation 296, 298
 Newton-Raphson algorithm 146
 Newton-Raphson formula 145
 Newton-Raphson method 145, 170, 178
 Nodes 300
 Non-linear equation 206, 411
 Non-linear fitting 329
 Non-linear function 121
 Non-linear system 168
 Non-singular matrix 511
 Normal equations 326
 Normalisation 52, 214
 Normalised numbers 52
 Numerical computing 1
 Numerical differentiation 347
 Numerical error 65
 Numerical instability 80
 Numerical integration 370
 Numerical method 3
 Numerical quadrature 370
 Numerical stability 8
 Numerically unstable 80

 Object program 26, 28

 Octal numbers 40, 44
 Octal system 44
 One-dimensional array 99
 One-step methods 412
 Open end method 124
 Open form 371
 Operation system 24
 Operators 102, 104, 105
 Order of convergence 86
 Order of equations 410
 Ordinary differential equations 410
 Outer loop 108
 Output devices 22
 Output statement 100
 Over determined 209
 Overflow 55

 Parabolic equation 484, 492
 Parallel computer 9, 20
 Parameter statement 99
 Partial derivatives 483
 Partial differential equation 410, 483
 Partial pivoting 220
 Particular solution 412
 PCAT 19
 PCXT 19
 Pentium 19
 Personal computer 19
 Perturbations 80
 Picard's method 417
 Piecewise polynomial 300
 Pivot element 219
 Pivot equation 214
 Pivoting 219
 Poisson's equation 490
 Polygon method 432
 Polynomial equation 122
 Polynomial fitting 333
 Polynomial method 473
 Positional form 98
 Power form 277
 Power method 477
 Precedence rule 102
 Precision 63
 Predictor 426
 Predictor-corrector method 426
 Printer 20
 Procedural error 65
 Processing block 96
 Program statement 95

- Program unit header 95
 Purification 177
 Quadratic convergence 148
 Quadratic equation 123, 195
 Quadruplets 43
 Radioactive decay 409
 Radix point 42
 Rate of convergence 8, 86
 Real constants 98
 Real numbers 51
 Real root 173
 Real type 97
 Recursive trapezoidal rule 391
 Regression analysis 323
 Regression line 326
 Regression plane 342
 Regula falsi 138
 Relational expression 104
 Relational operators 104
 Relative error 70
 Relaxation parameter 266
 Representation error 64
 Residual vector 243
 Richardson extrapolation 362, 389
 Romberg integration formula 390
 Romberg integration 371, 388
 Root squaring method 516
 Roots 121
 Rounding 65
 Roundoff error 8, 62, 65
 Row vector 506
 Row-sum norm 247
 Runge-Kutta method 436
 Run-time errors 114
 Sampling points 370
 Saturation growth rate 331
 Scaling factor 478
 Scatter diagram 323
 Scientific form 98
 Search bracket 126
 Secant algorithm 153
 Secant formula 152
 Secant method 151
 Second generation 16
 Semi-numeric method
 Sensitivity analysis 80
 Sequence structure 33
 Shifted power form 277
 Shooting method 464, 465
 Sign bit 50
 Significant digits 62
 Simple difference 294
 Simple harmonic motion 409
 Simpson's 1/3 rule 379
 Simpson's 3/8 rule 385
 Simpson's formula 445
 Simulation 13
 Simultaneous equations 207
 Singular systems 209
 Size 99
 Slide rule 14
 Softcopy 20
 Software 24
 Source program 28
 Special purpose computer 18
 Spiral convergence 163
 Spiral divergence 163
 Spline function 300
 Spline interpolation 299
 Spline interpolation polynomial 301
 Splines 300, 302
 Square matrix 505
 Stability 80
 Starting values 444, 499
 Statement functions 112
 Statement label 96
 Statement number 96
 Stopping criterion 127
 Straight-line logic 33
 Structured programming 33
 Subprogram 109
 Subroutine 109, 110
 Subscripted variable 99
 Subscripts 99
 Subtractive cancellation 56, 76
 Successive approximations 161
 Successive over relaxation 266
 Successive relaxation method 266
 Supercomputer 19
 Superlinear convergence 157
 Symbolic computation 9
 Symbolic constant 96
 Symmetric rounding 65, 66
 Synthetic division 175
 Tabulated function 347
 Taylor expansion 278
 Taylor series method 413
 Testing 113

- Third generation 16
Three-point formula 349
Top-down design 31
Trace of a matrix 474, 507
Transcendental equation 122
Transistor 16, 40
Translators 26, 28
Transpose matrix 506
Trapezoidal rule 372
Triangle inequality 73
Triangularisation 213
Tridiagonal system 307
Triplets 45, 49
Truncating 65
Truncation error 8, 67
Two-dimensional array 100
Two-point formula 348
Type declaration 95

Under determined 209
- Underflow 55
Unique solution 208
UNIVAC 15
Upper triangular matrix 231
User manual 114
Utility programs 25

Vacuum tube 16
Validation 3
Variable declaration 99
Variable loop 34
Variable-order approach 388
Variables 98
Vector computers 9

Wave equation 497
Weighting coefficient 370
Winchester disk 24
Workstation 19

Zero matrix 506

