# Revision notes - MA2213

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## 1 Computer Arithmetic and Computational Errors

## 1.1 Number Systems

**Definition 1.1** (Decimal System).

The number system we are used to is the **decimal system** and the number "10" which plays an important role is called the **base** of the decimal system.

In general, we can take other positive integer N > 1 as a base. In particular, N = 2, 8, 16 are used in most digital computers and the systems with N = 2, 8, 16 are known as **binary**, **octal**, **hexadecimal** number systems respectively.

## 1.1.1 Decimal to Binary

The conversion of decimal number to binary is performed in two steps. Take  $(53.7)_{10}$  as an example.

1. **Integer Part**. Convert decimal integers to binary by deviding by 2 successively and recording the remainders.

The remainders, 0 or 1, are recorded by starting at the decimal point and moving away to the left. For  $(53)_{10}$ , we have

$$53/2 = 26 \mathbf{R} 1$$
  
 $26/2 = 13 \mathbf{R} 0$   
 $13/2 = 6 \mathbf{R} 1$   
 $6/2 = 3 \mathbf{R} 0$   
 $3/2 = 1 \mathbf{R} 1$   
 $1/2 = 0 \mathbf{R} 1$ 

Therefore, the base 10 number 53 can be written as binary

$$(110101)_2 = 2^5 + 2^4 + 2^2 + 2^0 = 53$$

2. Fractional Part. Convert  $(0.7)_{10}$  to binary as follows:

Multiply by 2 successfully and record the integer parts, moving away from the decimal point to the right:

$$.7 \times 2 = .4 + 1$$

$$.4 \times 2 = .8 + 0$$

$$.8 \times 2 = .6 + 1$$

$$.6 \times 2 = .2 + 1$$

$$.2 \times 2 = .4 + 0$$

$$.4 \times 2 = .8 + 0$$
.....

Notice that the process repeats after four steps and will repeat infinitely exactly the same way. Therefore,

$$(0.7)_{10} = (.1\overline{0110})_2$$

3. Hence,

$$(53.7)_{10} = (110101.1\overline{0110})_2$$

#### 1.1.2 Binary to Decimal

The conversion of binary number to decimal needs to tackle nonterminating binary numbers. Take  $x = (.10\overline{101})_2$  as an example.

Multiplying by  $2^2$  shifts x to

$$y := 2^2 x = (10.\overline{101})_2 = (10)_2 + (.\overline{101})_2 = (2)_{10} + (.\overline{101})_2$$

The fractional part of y, i.e.,

$$z := (.\overline{101})_2$$

is calculated as follows:

$$2^3z = (101.\overline{101})_2$$

So,

$$(2^3 - 1)z = (101)_2 = (5)_{10}$$

i.e.

$$z = \left(\frac{5}{7}\right)_{10}$$

Thus,

$$y = (2)_{10} + \left(\frac{5}{7}\right)_{10} = \left(\frac{19}{7}\right)_{10}$$

and hence,

$$x = \frac{y}{2^2} = \left(\frac{19}{28}\right)_{10}$$

**Theorem 1.1** (Representation of Numbers).

In general, any **natural numbers**  $N \geq 2$  can be used as base.

Every positive real number a has a unique representation of the form

$$a = a_m N^m + a_{m-1} N^{m-1} + \dots + a_1 N^1 + a_0 N^0 + a_{-1} N^{-1} + a_{-2} N^{-2} + \dots$$

where

$$0 \le a_i \le N - 1, \quad a_m \ne 0$$

Equivalently,

$$a = \sum_{i=0}^{\infty} a_{m-i} N^{m-i}$$

where

$$0 < a_m < N; 0 \le a_{m-i} \le N - 1 \quad \forall i \ge 1$$

The above non-terminating representation of number a is not feasible in practical computation. We onl deal with, in practice, numbers which have terminating representation, i.e. of the following form

$$\alpha = \sum_{i=0}^{n-1} \alpha_{m-i} N^{m-i}$$

where  $0 < \alpha_m < N$  and  $0 \le \alpha_{m-i} < N$  for i = 1, 2, ..., n - 1.

**Definition 1.2** (Significant decimal digits).

If N = 10, then the numbers  $\alpha_{m-i}$ ,  $0 \le i \le n-1$  are called **significant decimal digits** and the number  $\alpha$  is said to have n significant decimal digits.

## 1.2 Chopping and Rounding

If the number a has more significant digits than what we wish to have, then we need to replace a by an approximate number  $a^*$  which contains a smaller number of digits, say n. There are two ways of terminating the number a to a given significant number of digits, namely **chopping** or **rounding**.

**Definition 1.3** (Chopping).

In **chopping**, we retain only the first n significant digits in the number a.

**Definition 1.4** (Rounding).

In **rounding**, the following rules are usually practiced:

- Retain the first n significant digits, and
- if the (n+1)th significant digit is less than 5, leave the nth significant digit unchanged;
- otherwise, if the (n + 1)th significant digit is greater or equal to 5, add unity to the nth significant digit.

## 1.3 Floating Point Representation

**Definition 1.5** (Floating Point Numbers).

Number of the form

$$\pm (0.a_1a_2a_3...a_m) \times N^e$$
, with  $a_1 \neq 0$ 

are called **floating point numbers**. The factor  $0.a_1a_2a_3...a_n$  is called the **mantissa**, e is called the **exponent** and N is the **base**.

The folating-point mode is for storing real numbers.

## 1.4 Basic Concepts in Error Estimation

**Definition 1.6** (Sources of Error).

Numerical results are affected by many types of error.

## • Error in Given Input Data

The input data can be result of measurements which are inexact, or produced by some arithmetic process using round-off process.

## • Round-off Errors During Computation

This is due to working with finite machine precision.

#### • Truncation Error

Consider the Taylor series xpansion for sin(x) about x = 0

$$\sin(x) = x - \frac{x^3}{3!} + \frac{x^5}{5!} - \frac{x^7}{7!} + \dots$$

which is an infinite series.

Truncation error occurs when the *infinite series* is broken off after a finite number of terms.

In general, truncation errors are error committed when a limiting process is truncated before one has come to the limiting value.

- Simplifications in the Mathematical Model
- Human Error

#### 1.5 Absolute and Relative Error

## **Definition 1.7** (Absolute Error).

Let  $x^*$  be an approximation to x, then the **absolute error** in approximating x is given by  $|x-x^*|$ .

From the above definition, the floating point representation of x, fl(x) has absolute error

$$|x - \mathrm{fl}(x)|$$

## Definition 1.8 (Absolute Error Bound).

Any non-negative number  $\delta(x^*)$  satisfying the inequality

$$|x - x^*| < \delta(x^*)$$

is called an absolute error bound.

## **Definition 1.9** (Decimal Places).

The two numbers x is said to agree to k decimal places if k is the largest non-negative integer such that

$$|x - x^*| \le 0.5 \times 10^{-k}$$

## **Definition 1.10** (Relative Error).

If  $x^*$  is an approximation to x, then the **relative error** is defined by

$$\frac{|x - x^*|}{|x|} \quad \text{provided that } x \neq 0$$

For a t-digit machine,

$$\frac{|x - \text{fl}(x)|}{|x|} \le \frac{1}{2} \times 10^{1-t}$$

in rounding mode and

$$\frac{|x - \mathrm{fl}(x)|}{|x|} \le 10^{1-t}$$

in chopping mode.

## 1.6 Computation in Floating Point Arithmetic

Take the example of addition/subtraction of two floating point number  $0.a_1a_2a_3...a_m \times 10^{N_1} \pm 0.b_1b_2b_3...b_m \times 10^{N_2}$ .

$$0.a_1 a_2 a_3 \dots a_m \times 10^{N_1} \pm 0.b_1 b_2 b_3 \dots b_m \times 10^{N_2}$$

$$= \begin{cases} (0.a_1 a_2 a_3 \dots a_m \pm 0.b_1 b_2 b_3 \dots b_m) \times 10^{N_1} & \text{if } N_1 = N_2 \\ (0.a_1 a_2 a_3 \dots a_m 0 \dots 0 \pm 0.0 \dots 0 b_1 b_2 b_3 \dots b_m) \times 10^{N_1} & \text{if } N_1 > N_2 \end{cases}$$

## 1.7 Propagation of Errors in Function Evaluation

#### 1.7.1 Function of one variable

In finding the value of f(x) by approximating  $f(x^*)$ , where  $x^*$  is the known approximating value of x,

let  $\delta(x^*)$  be an absolute error bound for  $|x - x^*|$ , i.e.,

$$|x - x^*| \le \delta(x^*)$$

If the function f(x) is differentiable, then by the mean-value theorem, we have

$$f(x) - f(x^*) = f'(p)(x - x^*)$$

where p is some number between x and  $x^*$ .

Therefore,

$$\delta(f(x^*)) \le \max_{t \in \mathbf{I}} |f^*(t)| \delta(x^*)$$

where **I** is the interval  $(x, x^*)$  if  $x < x^*$  or  $(x^*, x)$ .

Sometimes, upper bound of f' is unavailable. Yet, with the following assumption:

- $\delta(x^*)$  is small
- $f'(x^*) \neq 0$ , and
- f'(t) is nearly constant near  $x^*$  (i.e., f'(t) does not very greatly for t between x and  $x^*$ )

We have  $D \approx |f'(x^*)|$ , and consequently

$$\delta(f(x^*)) \approx |f'(x^*)|\delta(x^*)$$

#### 1.7.2 Function of Several Variables

**Theorem 1.2** (Error bounds of function of several variables).

Generalisation suggests, if  $x_i^*$  is an estimate of  $x_i$ ,  $1 \le i \le n$ , then with  $x^* = (x_1^*, x_2^*, \dots, x_n^*)$ , we have

$$\delta(f(x^* = (x_1^*, x_2^*, \dots, x_n^*)) \approx \left| \frac{\partial f(x^*)}{\partial x_1} \right| \delta(x_1^*) + \left| \frac{\partial f(x^*)}{\partial x_2} \right| \delta(x_2^*) + \dots + \left| \frac{\partial f(x^*)}{\partial x_n} \right| \delta(x_n^*)$$

## 1.8 Catastrophic Cancellation

Calculations involving the **subtraction** of two *nearly equal* numbers can result in considerable loss of accuracy due to cancellation.

Let  $x_1$  and  $x_2$  be two nearly equal numbers and denote the error in  $x_1$  and  $x_2$  by  $\delta(x_1)$  and  $\delta(x_2)$ , respectively, we have

$$y = x_1 - x_2$$

$$\Rightarrow |\delta(y)| \le |\delta(x_1)| + |\delta(x_2)|$$

$$\Rightarrow \left|\frac{\delta(y)}{y}\right| \le \frac{|\delta(x_1)| + |\delta(x_2)|}{|x_1 - x_2|}$$

Specifically, suppose two nearly equal numbers  $x_1$  and  $x_2$ , having the same exponent n and being in the k digi representations, are represented in floating point form:

$$x_1 = 0.a_1 a_2 \cdots a_p \alpha_{p+1} \cdots \alpha_k \times 10^n$$
  
$$x_2 = 0.a_1 a_2 \cdots a_p \beta_{p+1} \cdots \beta_k \times 10^n$$

Then their difference  $x_1 - x_2$  in floating point representation will have some digits in the mantissa to be 0, i.e.,

$$x_1 - x_2 = \text{fl}(\text{fl}(x_1) - \text{fl}(x_2)) = 0.\gamma_{p+1} \cdots \gamma_k \underbrace{0 \cdots 0}_{p \text{ 0's}} \times 10^{n-p}$$

So, in floating point calculation,  $x_1 - x_2$  has at most k - p significant digits; p significant digits have been lost or canceled due to the subtraction.

#### 1.8.1 Ways to Avoid Subtraction

Theorem 1.3  $(\sqrt{x+\varepsilon}-\sqrt{x})$ .

If  $|\varepsilon| \ll x$ , then

$$\sqrt{x+\varepsilon} - \sqrt{x} = \frac{\varepsilon}{\sqrt{x+\varepsilon} + \sqrt{x}}$$

presents a better way of calculation that reduces loss of accuracy.

**Theorem 1.4** (Roots of  $ax^2 + bx + c = 0$ ).

Consider the quadratic equation

$$ax^2 + bx + c = 0 \quad a \neq 0$$

The roots of equations are

$$x_1 = \frac{-b + \sqrt{b^2 - 4ac}}{2a}$$
 and  $x_2 = \frac{-b - \sqrt{b^2 - 4ac}}{2a}$ 

To reduce loss of accuracy, some alternative pathway include:

- $x_1 = \frac{-2c}{b + \sqrt{b^2 4ac}}$  and  $x_2 = \frac{2c}{-b + \sqrt{b^2 4ac}}$
- $x_1 + x_2 = -b$
- $\bullet \ x_1 x_2 = c$

**Theorem 1.5** (Evaluation of polynomial f(x)).

First note that  $ax^n = \text{fl}\left(\text{fl}\left(a\right) \times \text{fl}\left(x^n\right)\right)$  where  $\text{fl}\left(x^n\right) = \text{fl}\left(\text{fl}\left(x^{n-1}\right) \times \text{fl}\left(x\right)\right)$  recursively. Instead of calculating  $f(x) = \text{fl}\left(\sum_{i=0}^n a_i x^i\right)$ , nesting may provide better result:

$$f(x) = f((\cdots (a_n x + a_{n-1})x + \cdots)x + a_0)$$

## 1.9 Numerical Instability

**Definition 1.11** (Numerical Stability).

An algorithm is said to be **stable** if the effect of its errors on the final result is **negligible**.

## 2 Numerical Solution of Linear Systems of Equations

Definitions and method of computation of matrix can be found in MA2101.pdf, and are therefore omitted here.

## 2.1 System of Linear Equations

In this chapter, only linear systems with *unique* solution is concerned. Consider the linear system

$$Ax = b$$

where  $A \in \mathbb{M}_n(\mathbb{R})$  and  $b \in \mathbb{R}_c^n$ . If A is invertible, a unique solution x exists and given by

$$x = A^{-1}b$$

There are two classes of method of solving Ax = b numerically:

- **Direct** methods
- Iterative methods

In this chapter, only direct methods are studied.

Two basic direct method include

- Cramer's rule
- Gaussian Elimination

#### 2.1.1 Direct Method: Cramer's Rule

Definition 2.1 (Cramer's Rule).

By Cramer's rule, the  $n \times n$  linear system

$$Ax = b$$

has solution

$$x_i = \frac{d_i}{d} \quad i = 1, \dots, n$$

where  $d := \det(A) \neq 0, d_i = \det(A_i), A_i$  the matrix obtained by replacing the *i*-th column of A by b.

The determinant d is given by **Laplace Theorem**:

$$d = (-1)^{i+1}a_{i1}D_{i1} + (-1)^{i+2}a_{i2}D_{i2} + \dots + (-1)^{i+n}a_{in}D_{in}$$

where  $D_{ij}$  is the determinant of the submatrix obtained from A by deleting its ith row and jth column.

Theorem 2.1 (Computational Complexity of Cramer's Rule).

Suppose it needs  $m_n$  operations of multiplication to compute the determinant of a  $n \times n$  matrix. Then, we have

$$m_n = n + n m_{n-1}$$
  $m_1 = 1$ 

Here, the first term n is results from multiplication of  $a_{ij}$  and  $D_{ij}$ . Thus,

$$m_n = n + n m_{n-1} = n + n[(n-1) + (n-1)m_{n-2}]$$
  
=  $n + n(n-1) + n(n-1)(n-2) + \dots + n(n-1) \dots 3 \cdot 2$   
>  $n!$ 

Hence, in order to solve an  $n \times n$  linear system by Cramer's rule and Laplace Theorem, we have do at least

$$(n+1)n! = (n+1)!$$

multiplication.

Clearly, Cramer's rule is too computationally expensive.

Also, solving nonlinear linear system by **matrix inversion** is *computationally expensive* and often leads to more *inaccuracies*.

#### 2.2 Gaussian elimination

**Definition 2.2** (Elementary Row Operations).

Let  $E_i$  be the *i*th equation in a linear system Ax = b. The three **elementary row operations** permitted to solve this linear system are:

•  $E_i \leftarrow \lambda E_i$ , where  $\lambda$  is a nonzero constant.

$$\begin{bmatrix} 1 & & & & \\ & \ddots & & & \\ & & \lambda & & \\ & & & \ddots & \\ & & & & 1 \end{bmatrix} \begin{bmatrix} a_{11} & \cdots & a_{1n} \\ \vdots & & \vdots \\ a_{i1} & \cdots & a_{in} \\ \vdots & & \vdots \\ a_{n1} & \cdots & a_{nn} \end{bmatrix} = \begin{bmatrix} a_{11} & \cdots & a_{1n} \\ \vdots & & \vdots \\ \lambda a_{i1} & \cdots & \lambda a_{in} \\ \vdots & & \vdots \\ a_{n1} & \cdots & a_{nn} \end{bmatrix}$$

•  $E_i \leftarrow E_i - lE_j$ , where l is a nonzero constant.

$$\begin{bmatrix} 1 & & & & \\ & \ddots & & & \\ & -l & 1 & & \\ & & & \ddots & \\ & & & 1 \end{bmatrix}^{1} \begin{bmatrix} a_{11} & \cdots & a_{1n} \\ \vdots & & \vdots \\ a_{i1} & \cdots & a_{in} \\ \vdots & & \vdots \\ a_{n1} & \cdots & a_{nn} \end{bmatrix} = \begin{bmatrix} a_{11} & \cdots & a_{1n} \\ \vdots & & \vdots \\ a_{i1} - la_{j1} & \cdots & a_{in} - la_{jn} \\ \vdots & & \vdots \\ a_{n1} & \cdots & a_{nn} \end{bmatrix}$$

 $<sup>^{1}-</sup>l$  is at ith row and jth column of the leftmost matrix.

$$\begin{bmatrix}
1 & & & \\
& \ddots & & 1 \\
& & 1 & \\
& & 1 & \\
& & 1 & \\
& & & 1
\end{bmatrix}
\begin{bmatrix}
a_{11} & \cdots & a_{1n} \\
a_{i1} & \cdots & a_{in} \\
\vdots & \cdots & \vdots \\
a_{j1} & \cdots & a_{jn} \\
\vdots & \cdots & \vdots \\
a_{n1} & \cdots & a_{nn}
\end{bmatrix} =
\begin{bmatrix}
a_{11} & \cdots & a_{1n} \\
a_{j1} & \cdots & a_{jn} \\
\vdots & \cdots & \vdots \\
a_{i1} & \cdots & a_{in} \\
\vdots & \cdots & \vdots \\
a_{n1} & \cdots & a_{nn}
\end{bmatrix}$$

By a sequence of the above operations, a linear system can be transformed to a more easily solved linear system with the **same** set of solutions.

The Gaussian elimination for solving Ax = b involves 2 stages/

- A forward course where a sequence of elementary row operations are applied to A to reduce it to an upper triangular form **U**.
- Backward substitution process

**Theorem 2.2** (General method of Gaussian Elimination on Ax = B). Let the original system be denoted by

$$\mathbf{A}^{(0)}\mathbf{x} = \mathbf{b}^{(0)}$$
 where  $\mathbf{A}^{(0)} = \mathbf{A} = (a_{ij}^{(0)}), \ \mathbf{b}^{(0)} = \mathbf{b} = \begin{bmatrix} b_1^{(0)} \\ \vdots \\ b_i^{(0)} \\ \vdots \\ b_n^{(0)} \end{bmatrix}$ 

Step 1: If  $a_{11}^{(0)} \neq 0$ , let

$$l_{i,1} = \frac{a_{i,1}^{(0)}}{a_{1,1}^{(0)}} \qquad i = 2, 3, \dots, n$$

$$a_{i,j}^{(1)} = a_{i,j}^{(0)} - l_{i,1} \times a_{1,j}^{(0)3} \qquad i, j = 2, 3, \dots, n$$

$$b_i^{(1)} = b_i^{(0)} - l_{i,1} \times b_1^{(0)} \qquad i = 2, \dots, n$$

Then we obtain the equivalent system

$$\mathbf{A}^{(1)}\mathbf{x} = \mathbf{b}^{(1)}$$

where

$$\mathbf{A}^{(1)} = \begin{bmatrix} a_{11}^{(0)} & a_{12}^{(0)} & \cdots & a_{1n}^{(0)} \\ 0 & a_{22}^{(1)} & \cdots & a_{2n}^{(1)} \\ 0 & \vdots & \ddots & \vdots \\ 0 & a_{n2}^{(1)} & \cdots & a_{nn}^{(1)} \end{bmatrix}, \quad \mathbf{b}^{(1)} = \begin{bmatrix} b_1^{(0)} \\ b_2^{(1)} \\ \vdots \\ b_n^{(1)} \end{bmatrix}$$

 $<sup>^{3}</sup>$ In general,  $l_{p,q}$  is used in the operations on qth column, and denotes the ratio between the entry on the pth row(p > q),  $a_{p,q}^{(q-1)}$ , to the qth diagonal entry,  $a_{q,q}^{(q-1)}$ .

3In general,  $a_{i,j}^{(k)}$ , (i < j) is obtained by operations on the kth column, and is the result of zeroing the  $a_{i,k}^{(k)}$ .

entry(i > k), so it equals the previous entry  $a_{i,j}^{(k-1)}$  minus  $l_{i,k}$  times the entry on the kth row  $a_{k,j}^{(k-1)}$ .

Step 2: If  $a_{22}^{(1)} \neq 0$ , repeat Step 1 to eliminate  $\mathbf{x}_2$  from row 3 to n. After k steps, we obtain the equivalent system

$$\mathbf{A}^{(k)}\mathbf{x} = \mathbf{b}^{(k)}$$

where  $\mathbf{A}^{(k)}$  takes the form

$$\mathbf{A}^{(k)} = \begin{bmatrix} a_{11}^{(0)} & \cdots & \cdots & \cdots & a_{1n}^{(0)} \\ & a_{22}^{(1)} & \cdots & \cdots & a_{2n}^{(1)} \\ & & \ddots & \vdots & \cdots & \ddots \\ & & a_{k+1,k+1}^{(k)} & \cdots & a_{k+1,n}^{(k)} \\ & & \vdots & \ddots & \vdots \\ & & a_{n,k+1}^{(k)} & \cdots & a_{n,n}^{(k)} \end{bmatrix}$$

If  $a_{k+1,k+1}^{(k)} \neq 0$ , then let

$$l_{i,k+1} = \frac{a_{i,k+1}^{(k)}}{a_{k+1,k+1}^{(k)}}$$

$$a_{i,j}^{(k+1)} = a_{i,j}^{(k)} - l_{i,k+1} \times a_{k+1,j}^{(k)}$$

$$b_i^{(k+1)} = b_i^{(k)} - l_{i,k+1} b_{k+1}^{(k)}$$

Finally, after at most n-1 steps, the system becomes

$$\mathbf{A}^{(n-1)}\mathbf{x} = \mathbf{b}^{(n-1)}$$

which is an upper triangular system, where

$$\mathbf{U} = \mathbf{A}^{(n-1)} = \begin{bmatrix} a_{11}^{(0)} & a_{12}^{(0)} & \cdots & a_{1n}^{(0)} \\ & a_{22}^{(1)} & \cdots & a_{2n}^{(1)} \\ & & \ddots & \vdots \\ & & & a_{nn}^{(n-1)} \end{bmatrix}$$

and

$$\mathbf{b}^{(n-1)} = \begin{bmatrix} b_1^{(0)} \\ b_2^{(1)} \\ \vdots \\ b_n^{(n)} \end{bmatrix}$$

Step 3: The solution vector  $\mathbf{x}$  is solved by back substitution:

$$x_n = \frac{b_n^{(n-1)}}{a_{nn}^{(n-1)}}$$

and for  $k = n - 1, n - 2, \dots, 1$ ,

$$x_k = \frac{1}{a_{kk}^{(k-1)}} \left( b_k^{(k-1)} - \sum_{j=k+1}^n a_{kj}^{(k-1)} x_j \right)$$

The above algorithm is known as **Gaussian elimination**.

## 2.3 Triangular Factorisation: A = LU

We note, in the (k+1)th step of above Gaussian elimination, all rows below the (k+1)th row are minused by a multiple of the (k+1)th row. This operation can be viewed by a collection of elementary row operation 2 and represented by

$$\mathbf{A}^{(k+1)} = \mathbf{L}^{(k)} \mathbf{A}^{(k)}$$

where

$$\mathbf{L}^{(k)} = \begin{bmatrix} 1 & & & & & \\ & \ddots & & & \\ & & 1 & & \\ & & -l_{k+2,k+1} & \ddots & \\ & & \vdots & & \ddots & \\ & & -l_{n,k+1} & & & 1 \end{bmatrix}$$

This relation gives

$$\mathbf{U} = \mathbf{A}^{(n-1)} = \mathbf{L}^{(n-2)} \mathbf{L}^{(n-3)} \cdots \mathbf{L}^{(0)} \mathbf{A}^{(0)}$$

Also, note the inverse of  $\mathbf{L}^{(k)}$  is

$$(\mathbf{L}^{(k)})^{-1} = \begin{bmatrix} 1 & & & & \\ & \ddots & & & \\ & & 1 & & \\ & & l_{k+2,k+1} & \ddots & \\ & & \vdots & & \ddots & \\ & & l_{n,k+1} & & & 1 \end{bmatrix}$$

a flip of signs of all  $l_{i,k+1}$  from  $\mathbf{L}^{(k)}$ .

Therefore,

$$(\mathbf{L}^{(0)})^{-1}(\mathbf{L}^{(1)})^{-1}\cdots(\mathbf{L}^{(n-2)})^{-1}\mathbf{U}=\mathbf{A}$$

Note that  $(\mathbf{L}^{(0)})^{-1}(\mathbf{L}^{(1)})^{-1}\cdots(\mathbf{L}^{(n-2)})^{-1}$  is upper triangular and is defined as  $\mathbf{L}$ :

$$(\mathbf{L}^{(0)})^{-1}(\mathbf{L}^{(1)})^{-1} \cdots (\mathbf{L}^{(n-2)})^{-1} = \begin{bmatrix} 1 & & & & \\ l_{21} & \ddots & & & \\ \vdots & & 1 & & & \\ l_{k+2,1} & \cdots & l_{k+2,k+1} & \ddots & & \\ \vdots & & \vdots & & \ddots & \\ l_{n,1} & \cdots & l_{n,k+1} & \cdots & l_{n,n-1} & 1 \end{bmatrix} := \mathbf{L}$$

Hence,  $\mathbf{A} = \mathbf{L}\mathbf{U}$ , where  $\mathbf{U} = \mathbf{A}^{(n-1)}$  and  $\mathbf{L}$  is a lower triangular matrix.

**Theorem 2.3.** If **U** and **L** are the upper and lower triangular matrices defined above, then A = LU.

## 2.4 Compact Forms of Gaussian Elimination

If **A** admits an **LU** decomposition, we can solve the system of equations  $\mathbf{LUx} = \mathbf{b}$  in two stages.

- Set z := Ux. Solve Lz = b for z.
- Solve  $\mathbf{U}\mathbf{x} = \mathbf{z}$  for  $\mathbf{x}$

Theorem 2.4 (Existence and Uniqueness of LU decomposition).

Let **A** be an  $n \times n$  matrix and  $\mathbf{A}^{(k)}$  be the  $k \times k$  matrix formed from the first k rows and columns of **A**. If  $\det(\mathbf{A}^{(k)}) \neq 0 \forall k = 1, 2, ..., n-1$ , then there exists a unique lower triangular matrix  $\mathbf{L} = (l_{ij})$  with  $l_{ii} = 1 \forall i = 1, 2, ..., n$  and a unique upper triangular matrix  $\mathbf{U} = (u_{ij})$  such that  $\mathbf{A} = \mathbf{L}\mathbf{U}$ .

## 2.5 Compact Forms of Gaussian Elimination

The **LU** decomposition can be calculated directly by Gaussian Elimination. Given that mathbfA admits a **LU** decomposition  $\mathbf{A} = \mathbf{LU}$ , we have, entry-wise

$$a_{ij} = \sum_{k=1}^{n} l_{ik} u_{kj}$$
  $i, j = 1, 2, \dots, n$ 

Note that the above system has  $n^2$  equations and  $n^2 + n$  unknowns  $l_{ik}$ ,  $i \ge k$  and  $u_{kj}$ ,  $k \le j$ . Thus n unknown may be set arbitrarily.

#### Theorem 2.5 (Doolittle Method).

For Doolittle Method, set diagonal entries of  $\mathbf{L}$ ,  $l_{kk} := 1, k = 1, 2, \dots n$  and assume that  $u_{kk} \neq 0, \forall k \in [1, n] \cap \mathbb{Z}$ .

The sequence of calculation is as follows:

- 1. First row of **U**:  $u_{1k}$ ,  $1 \le k \le n$
- 2. First column of **L**:  $l_{k1}, 2 \le k \le n^4$
- 3. Second row of **U**:  $u_{2k}$ ,  $2 \le k \le n$ .

:

Obviously, when calculating kth row of U, row 1 to k-1 of U and column 1 to k-1 of L will be known.

Also, when calculating kth row of **L**, row 1 to k of **U** and column 1 to k-1 of **L** will be known.

Suppose  $u_{kj}$ ,  $(j \ge k)$  is of concern, we have

$$a_{kj} = \sum_{r=1}^{k} l_{kr} u_{rj}$$

<sup>&</sup>lt;sup>4</sup>if  $k = 1, l_{k1} = l_{11} = 1$  by definition

Substitute in  $l_{kk} = 1$  and rearranging,

$$u_{kj} = a_{kj} - \sum_{r=1}^{k-1} l_{kr} u_{rj}$$

Afterwards, suppose  $l_{ik}$ , (i > k) is of concern, we have

$$a_{ik} = \sum_{r=1}^{k} l_{ir} u_{rk}$$

Rearranging,

$$a_{ik} = \sum_{r=1}^{k-1} l_{ir} u_{rk} + l_{ik} u_{kk}$$

Therefore,

$$l_{ik} = \frac{a_{ik} - \sum_{r=1}^{k-1} l_{ir} u_{rk}}{u_{kk}}$$

The above two equation can be interleaved to obtain L and U.

Theorem 2.6 (Crout Method).

For Crout Method, set diagonal entries of  $\mathbf{U}$ ,  $u_{ii}=1,k=1,2,\ldots,n$  and assume that  $l_{kk} \neq 0 \forall k \in [1,n] \cap \mathbb{Z}$ .

The sequence of calculation is as follows:

1. First column of L:  $l_{k1}$ ,  $1 \le k \le n$ 

2. First row of **U**:  $u_{1k}$ ,  $2 \le k \le n$ 

3. Second column of L:  $l_{k2}, 2 \le k \le n$ 

:

Obviously, when calculating kth column of L, column 1 to k-1 of L and row 1 to k-1 of U will be known.

Also, when calculating kth row of U, column 1 to k of L and row 1 to k-1 of U will be known.

Suppose  $l_{ik}$ ,  $(i \ge k)$  is of concern, we have

$$a_{ik} = \sum_{r=1}^{k} l_{ir} u_{rk}$$

Substitute in  $u_{kk} = 1$  and rearranging,

$$l_{ik} = a_{ik} - \sum_{r=1}^{k-1} l_{ir} u_{rk}$$

Afterwards, suppose  $u_{kj}$ ,  $(j \ge k + 1)$  is of concern, we have

$$a_{kj} = \sum_{r=1}^{k} l_{kr} u_{rj}$$

Rearranging,

$$u_{kj} = \frac{a_{kj} - \sum_{r=1}^{k-1} l_{kr} u_{rj}}{l_{kk}}$$

The above two equation can be interleaved to obtain L and U.

## 2.6 LU Decomposition for Tridiagonal Matrices

#### Theorem 2.7.

Let A be a tridiagonal matrices

$$\mathbf{A} = \begin{bmatrix} a_1 & c_1 \\ b_2 & a_2 & c_2 \\ & \ddots & \ddots & \ddots \\ & & \ddots & \ddots & \ddots \\ & & & b_{n-1} & a_{n-1} & c_{n-1} \\ & & & b_n & a_n \end{bmatrix}$$

If an LU decomposition exists for A, then

$$\mathbf{A} = \mathbf{L}\mathbf{U}$$

where

By applying Doolittle's method, it can be shown that

$$\alpha_1 = a_1$$

and for k = 2, 3, ..., n,

$$\beta_k = \frac{b_k}{\alpha_{k-1}}$$

$$\alpha_k = a_k - \beta_k c_{k-1}$$

Theorem 2.8 (Thomas's Algorithm).

If  $\mathbf{A}\mathbf{x} = \mathbf{g}$  and define  $\mathbf{U}\mathbf{x} = h$  where  $\mathbf{x} := (x_1, x_2, \dots, x_n)^T$ ,  $\mathbf{g} = (g_1, g_2, \dots, g_n)^T$  and  $\mathbf{h} = (h_1, h_2, \dots, h_n)^T$ .

$$h_1 = g_1$$
  
 $h_i = g_i - \beta_i h_{i-1}, \quad i = 2, 3, \dots, n$ 

and

$$x_n = \frac{h_n}{\alpha_n}$$
$$x_i = \frac{h_i - c_i x_{x+1}}{\alpha_i}$$

This method is known as **Thomas algorithm**.

## 2.7 Pivoting Strategies

#### 2.7.1 Partial Pivoting

**Definition 2.3** (Partial Pivoting).

At the (k+1)th step of the Gaussian Elimination process,  $k=0,1,\ldots,n-2$ , choose the element having maximum absolute value in the (k+1)th column of  $\mathbf{A}^{(k)}$  that lies on or below the diagonal so that

 $|a_{s,k+1}^{(k)}| = \max_{i} |a_{i,k+1}^{(k)}| \quad k+1 \le i \le n$ 

and interchange row k+1 with row s.

#### 2.7.2 Scaled Partial Pivoting

**Definition 2.4** (Scaled Partial Pivoting).

In the beginning, calculate  $s_i$  for all row i, where

$$s_i = \max_{1 \le j \le n} \{|a_{ij}|\}, i = 1, \dots, n$$

At the (k+1)th step of the Gaussian Elimination process,  $k=0,1,\ldots,n-2$ , choose rth row of  $\mathbf{A}^{(k)}$  that lies on or below the diagonal where r is determined by

$$\frac{|a_{r,k+1}^{(k)}|}{s_r} = \max_i \frac{|a_{i,k+1}^{(k)}|}{s_i} \quad k+1 \le i \le n$$

and exchange row k+1 with row r and  $s_{k+1}$  with  $s_r$ .

## 3 Error of Approximation

**Definition 3.1** (Error).

Given n+1 data values  $(x_0, f_0), (x_1, f_1), \ldots, (x_n, f_n)$ , we may define the error of approximation, E, by

$$E = \sum_{i=0}^{n} |p(x_i) - f_i|$$

If we take p(x) to be a polynomial of degree n then by choosing

$$p(x_i) = f_i, i = 0, 1, \dots, n$$

we can make E = 0. The function p(x) is the **Lagrange interpolating polynomial**. Alternatively, when n is large, we may fit a polynomial of degree k where  $k \ll n$ . We use

$$E = \sum_{i=0}^{n} [p(x_i) - f_i]^2$$

By minimising E we achieve a **least square fit**.

If the function f(x) is continuous specified on the interval [a, b], we may define the error E in approximating f(x) by p(x) as

$$E = \max_{a \le x \le b} |p(x) - f(x)|$$

It is common in practice to take p(x) as a polynomial and minimise E with respect to variations in the coefficients. This generate the **minimax** polynomial approximation.

The following theorem justifies the choice of polynomial for approximation

**Theorem 3.1** (Weierstrass Approximation Theorem).

Suppose f is defined and continuous on [a, b]. For each  $\varepsilon > 0$  there exists a polynomial P(x), defined on [a, b], with property that

$$|f(x) - P(x)| < \varepsilon \quad \forall x \in [a, b]$$

## 3.1 Least Square Approximation

#### 3.1.1 Discrete Data

Given a set of m+1 discrete data points  $(x_0, f(x_0)), (x_1, f(x_0)), \dots (x_m, f(x_m))$ , each carrying weight  $w_0, w_1, \dots, w_m$  respectively.

The task is to find an approximating polynomial function<sup>5</sup>

$$p_n(x; a_0, a_1, \dots, a_n) := a_0 + a_1 x + \dots + a_n x^n$$

<sup>&</sup>lt;sup>5</sup>The subscript n denotes the power of the polynomial; thus the polynomial  $p_n$  has n+1 unknowns

(m > n) such that

$$E(a_0, a_1, \dots, a_n) = \sum_{i=0}^{m} w_i [f(x_i) - p_n(x_i; a_0, a_1, \dots, a_n)]^2$$

is minimised with respect to the parameters  $a_0, a_1, \dots, a_n$ . We require

$$\frac{\partial E}{\partial a_i} = 9 \forall j = 0, 1, \dots, n$$

Therefore, by differentiating with  $a_i$ ,

$$\sum_{i=0}^{m} w_i [f(x_i) - p_n(x_i; a_0, a_1, \dots, a_n)] \frac{\partial}{\partial a_j} (-p_n(x_i; a_0, a_1, \dots, a_n)) \quad \forall j = 0, 1, \dots, n$$

Rearranging, we have

$$\sum_{i=0}^{m} w_{i} \underbrace{(a_{0} + a_{1}x_{i} + \dots + a_{n}x_{i}^{n})}_{p_{n}(x_{i}; a_{0}, a_{1}, \dots, a_{n})} x_{i}^{j} = \sum_{i=0}^{m} w_{i} f(x_{i}) x_{i}^{j} \quad \forall j = 0, 1, \dots, n$$

There is n+1 unknowns, and n+1 equations, so the linear system is<sup>7</sup>

$$\begin{bmatrix} \sum_{i=0}^{m} w_{i} x_{i}^{0} & \sum_{i=0}^{m} w_{i} x_{i}^{1} & \cdots \sum_{i=0}^{m} w_{i} x_{i}^{n} \\ \sum_{i=0}^{m} w_{i} x_{i}^{1} & \sum_{i=0}^{m} w_{i} x_{i}^{2} & \cdots \sum_{i=0}^{m} w_{i} x_{i}^{n+1} \\ \vdots & \vdots & \ddots & \vdots \\ \sum_{i=0}^{m} w_{i} x_{i}^{n} & \sum_{i=0}^{m} w_{i} x_{i}^{n+1} & \cdots \sum_{i=0}^{m} w_{i} x_{i}^{2n} \end{bmatrix} \begin{bmatrix} a_{0} \\ a_{1} \\ \vdots \\ a_{n} \end{bmatrix} = \begin{bmatrix} \sum_{i=0}^{m} w_{i} f(x_{i}) x_{i}^{0} \\ \sum_{i=0}^{m} w_{i} f(x_{i}) x_{i}^{1} \\ \vdots \\ \sum_{i=0}^{m} w_{i} f(x_{i}) x_{i}^{n} \end{bmatrix}$$

#### 3.1.2 Continuous Function

Given a continuous function f(x) defined on an interval [a, b] with each value  $x \in [a, b]$  associated with a weight w(x).

The task is to find an approximating polynomial function  $p_n(x; a_0, a_1, \dots, a_n)$  such that

$$E(a_0, a_1, \dots, a_n) = \int_a^b w(x) [f(x) - p_n(x; a_0, a_1, \dots, a_n)]^2 dx$$

is minimised with respect to the parameters  $a_0, a_1, \ldots, a_n$ .

The necessary condition sfor  $E(a_0, a_1, \ldots, a_n)$  to be minimum are

$$\frac{\partial E}{\partial a_j} = 0 \forall j = 0, 1, \dots, n$$

Differentiating with respect to  $a_j$ ,

$$\frac{dE}{da_j} = -2 \int_a^b w(x) [f(x) - p(x; a_0, a_1, \dots, a_n)] \frac{\partial}{\partial a_j} (p_n(x; a_0, a_1, \dots, a_n)) dx = 0$$

 $<sup>^6</sup>E$  can be understood as a weighted sum of squares of residue for each data.

<sup>&</sup>lt;sup>7</sup>the jth column corresponding to the equation of specific j

Rearranging,<sup>8</sup>

$$\int_{a}^{b} w(x)(a_0 + a_1 x + \dots + a_n x^n) x^j dx = \int_{a}^{b} w(x) f(x) x^j dx \quad \forall j = 0, 1, \dots, n$$

$$\sum_{k=0}^{n} \int_{a}^{b} w(x) x^{k+j} dx = \int_{a}^{b} w(x) f(x) x^j dx \quad \forall j = 0, 1, \dots, n$$

Similarly, a matrix equation can be formed from this linear system of n + 1 variables and n + 1 equations. However, this matrix is **ill-conditioned**.

## 3.2 Lagrange Interpolation

#### Definition 3.2.

Given a set of n+1 data points  $(x_0, f_0), (x_1, f_1), \ldots, (x_n, f_n)$ , we seek to find a polynomial  $p_n(x)$  of degree n which passes through each of the given points, i.e.

$$p_n(x_i) = f_i \quad i = 0, 1, 2, \dots, n$$

If we write  $p_n(x) = a_0 + a_1x + a_2x^2 + \cdots + a_nx^n$ , we have

$$\begin{bmatrix} 1 & x_0 & \cdots & x_0^n \\ 1 & x_1 & \cdots & x_1^n \\ \vdots & \vdots & \ddots & \vdots \\ 1 & x_n & \cdots & x_n^n \end{bmatrix} \begin{bmatrix} a_0 \\ a_1 \\ \vdots \\ a_n \end{bmatrix} = \begin{bmatrix} f_0 \\ f_1 \\ \vdots \\ f_n \end{bmatrix}$$

The determinant of the  $(n+1) \times (n+1)$  Vandermonde matrix is  $\prod_{0 \le i < j \le n} (x_j - x_i)$ , so the system has a unique solution as long as  $x_i, i = 1, 2, \ldots, n$  are all distinct. The solved polynomial is the unique Lagrange interpolating polynomial.

**Theorem 3.2** (Computation of Lagrange Interpolating Polynomial). Write

$$p_n(x) = \alpha_0(x - x_1)(x - x_2) \cdots (x - x_n) + \alpha_1(x - x_0)(x - x_2) \cdots (x - x_n) + \cdots + \alpha_n(x - x_0)(x - x_1) \cdots (x - x_{n-1})$$

Then, we have

$$\alpha_i = \frac{f_i}{\prod_{\substack{j=0\\j\neq i}}^n (x_i - x_j)}$$

 $<sup>\</sup>frac{\partial}{\partial a_i}(p_n(x; a_0, a_1, \dots, a_n)) = x^j$  is obvious.

So,

$$p_n(x) = \sum_{i=0}^n l_i(x) f_i$$

where

$$l_i(x) = \prod_{\substack{j=0\\j\neq i}}^n \frac{x - x_j}{x_i - x_j}$$

Note that  $l_i(x)$  admits the property that

$$l_i(x_i) = 1$$
  
 $l_i(x_j) \neq 1 \text{ for } j \neq i$ 

**Theorem 3.3** ( $\psi(x)$  and alternative expression of Lagrange Interpolating Polynomial). Let  $\psi(x)$  denote the polynomial

$$\psi(x) = (x - x_0)(x - x_1) \cdots (x - x_n) = \prod_{i=0}^{n} (x - x_i)$$

then, by differentiating on each product term once at a time, we have

$$\psi'(x) = \sum_{i=0}^{n} \prod_{\substack{j=0\\j\neq i}}^{n} (x - x_j)$$

and hence

$$\psi'(x_i) = (x_i - x_0) \cdots (x_i - x_{i-1})(x_i - x_{i+1}) \cdots (x_i - x_n)$$

as all other terms containing  $(x - x_i)$  will vanish. Thus,<sup>9</sup>

$$l_i(x) = \frac{\psi(x)}{(x - x_i)\psi'(x_i)}$$

So  $p_n(x)$  may be written as

$$p_n(x) = \sum_{i=0}^{n} \frac{\psi(x)}{(x - x_i)\psi'(x_i)} f_i$$

## 3.3 Truncation Error of Interpolating Polynomial

**Theorem 3.4** (Extended Mean Value Theorem). Suppose that

$$a \le x_0 < x_1 < \dots < x_k \le b$$

$$f(x_0) = f(x_1) = \dots = f(x_k) = 0$$

and  $f(x), f'(x), \ldots, f^{(k)}(x)$  are all continuous on [a, b]. There is a  $\xi \in (x_0, x_n) \subset (a, b)$  such that

$$f^k(\xi) = 0$$

<sup>&</sup>lt;sup>9</sup>Do NOT forget multiply  $(x - x_i)$  while using this formula!

Theorem 3.5 (Lagrange Interpolating Polynomial Error Formula).

Let  $f(x) \in C^{(n+1)}[a,b]$ . Further, let  $p_n(x)$  interpolates f(x) at (n+1) distinct points  $x+0, x_1, \ldots, x_n \in [a,b]$ . Then, for  $x \in [a,b], x \neq x_i, i=0,1,\ldots,n$ ,

$$f(x) - p_n(x) = \psi(x) \frac{f^{(n+1)}(\xi)}{(n+1)!}$$

for some  $xi \in \text{Spr}\{x, x_0, x_1, \dots, x_n\}$ .

Here,  $\operatorname{Spr}\{x, x_0, x_1, \dots, x_n\}$  denotes the smallest interval containing  $x, x_0, x_1, \dots, x_n$ .

It follows that an upper bound for the error is

$$\delta(f(x)) := |f(x) - p_n(x)| \le \frac{M}{(n+1)!} |\psi(x)|$$

where

$$M = \max_{a \le \xi \le b} |f^{(n+1)}(\xi)|$$

#### 3.4 Divided Differences

**Definition 3.3** (Divided Differences).

Let  $(x_0, f(x_0)), \ldots, (x_n, f(x_n))$  be n+1 given data points. The **first divided difference** of f(x) between  $x_0$  and  $x_1$ , denoted by  $f[x_0, x_1]$  is defined by

$$f[x_0, x_1] := \frac{f(x_1) - f(x_0)}{x_1 - x_0}$$

Analogously, the **second divided difference** for the triplet  $(x_0, x_1, x_2)$  is defined by

$$f[x_0, x_1, x_2] := \frac{f[x_1, x_2] - f[x_0, x_1]}{x_2 - x_0}$$

And, the *n*th divided difference of f(x) at (n+1)-tuple  $(x_0, \ldots, x_n)$  is defined recursively in terms of the (n-1)th divided difference as follows:

$$f[x_0, x_1, \dots, x_n] := \frac{f[x_1, \dots, x_n] - f[x_0, \dots, x_{n-1}]}{x_n - x_0}$$

**Theorem 3.6** (Divided difference of self).

We define  $f[x_i, x_i]$  by

$$f[x_i, x_i] = \lim_{x_{i+1} \to x_i} f[x_i, x_{i+1}] = f'(x_i)$$

provided  $f'(x_i)$  exists.

Similarly, we define rth divided difference of self is

$$f[\underbrace{x_i,\ldots,x_i}_{rx_i's}] = \frac{1}{r!}f^{(r)}(x_i)$$

## 3.5 Newton Interpolation Formula

It is easy to see, by definition of first order divided difference, that

$$f(x) = f(x_0) + (x - x_0)f[x, x_0]$$

Therefore, we can obtain

$$f(x) = f(x_0) + (x - x_0)f[x_0, x_1] + \dots + (x - x_0) + \dots + (x - x_{n-1})f[x_0, \dots, x_n] + R_n(x)$$

$$= f(x_0) + \sum_{i=0}^{n-1} f[x_0, \dots, x_{i+1}] \prod_{i=0}^{i} (x - x_i) + R_n(x)$$

where

$$R_n(x) = \psi(x) f[x, x_0, x_1, \dots, x_n]$$

The Newton interpolating formula of these n+1 data points is

$$p_n(x) = f(x_0) + (x - x_0)f[x_0, x_1] + \dots + (x - x_0) \cdot \dots \cdot (x - x_{n-1})f[x_0, \dots, x_n]$$

To see this, we note that at  $x_0, \ldots, x_n$ ,  $R_n = 0$ , so  $f(x) = p_n(x)$ ; also  $p_n(x)$  is at most of degree n.

When expanding the formula  $p_n$  to  $p_{n+1}$ , we note that  $f = p_{n+1} = p_n + (x - x_0) \cdots (x - x_n) f[x_0, \dots, x_{n+1}]$ , so the divided difference can be easily acquired.

**Remark**:  $f[x_0, \ldots, x_n] = f[x_{i_0}, \ldots, x_{i_n}]$ , where  $\{i_k\}_{k=1}^n$  is a permutation of  $\{k\}_{k=1}^n$ .

## 3.6 Cubic Spline Interpolation

Cubic Spline interpolation is a piecewise approximation used to replace high order polynomial interpolation, latter of which has the problem of wild oscillation.

## Definition 3.4 (Cubic Spline Approximation).

Given a function f defined on [a,b] and a set of data points  $x_i$  with  $a=x_0 < x_1 < \cdots < x_{n-1} < x_n = b$ , a **cubic spline interpolant** to f is a *piecewise* function S satisfying the following conditions:

- 1. S is a cubic polynomial, denoted  $S_j$  on the subinterval  $[x_j, x_{j+1}]$  for  $j = 0, 1, \ldots, n-1$ .
- 2. Interpolation condition:  $S(x_j) = f(x_j)$ .
- 3.  $S \in C^1(a,b): S'_j(x_{j+1}) = S'_{j+1}(x_{j+1}) \quad j = 0, 1, \dots, n-2.$
- 4.  $S \in C^2(a,b): S_j''(x_{j+1}) = S_{j+1}''(x_{j+1})$   $j = 0, 1, \dots, n-2$ .
- 5. One of the following sets of boundary condition is satisfied
  - $S_0''(x_0) = S_{n-1}''(x_n) = 0$  (Natural boundary condition)
  - Or,  $S'_0(x_0) = f'(x_0)$  and  $S'_{n-1}(x_n) = f'(x_n)$  (Clamped boundary condition)

**Theorem 3.7** (Solving cubic spline interpolation efficiently).

The following method will produce the cubic spline interpolation of n+1 data points by solving a linear system of maximum size  $(n+1) \times (n+1)$ .

Let  $s_i(x)$  be a cubic polynomial on  $[x_i, x_{i+1}]$ . Then  $s_i''(x)$  is the linear polynomial on  $[x_i, x_{i+1}]$ . Let  $M_i = s_i''(x_i)$ .  $M_i$  will be calculated later using (3) and (5).

Using lagrange interpolation polynomial, we have

$$s_i''(x) = \frac{x_{i+1} - x}{h_i} M_i + \frac{x - x_i}{h_i} M_{i+1}$$

where  $h_i = x_{i+1} - x_i$  the step size. Integrating twice, we have

$$s_i(x) = \frac{(x_{i+1} - x)^3}{6h_i} M_i + \frac{(x - x_i)^3}{6h_i} M_{i+1} + Ax + B$$

for some constant A and B. Using (2), i.e.,

$$s_i(x_i) = f_i$$
  $s_i(x_{i+1}) = f_{i+1}$ 

we have, nicely<sup>10</sup>

$$\frac{h_i^2}{6}M_i + Ax_i + B = f_i$$

$$\frac{h_i^2}{6}M_{i+1} + Ax_{i+1} + B = f_{i+1}$$

Solving which we have

$$\begin{cases} A &= \frac{f_{i+1} - f_i}{h_i} + \frac{h_i}{6} (M_i - M_{i+1}) \\ B &= \frac{x_{i+1} f_i - x_i f_{i+1}}{h_i} + \frac{h_i}{6} (x_i M_{i+1} - x_{i+1} M_i) \end{cases}$$

Therefore,

$$s_i(x) = \frac{6}{h_i} [(x_{i+1} - x)^3 M_i + (x - x_i)^3 M_{i+1}]$$
$$- \frac{h_i}{6} [(x_{i+1} - x) M + (x - x_i) M_{i+1}]$$
$$+ \frac{1}{h_i} [(x_{i+1} - x) f_i + (x - x_i) f_{i+1}]$$

From here, differentiate  $s_i(x)$ , we have expression for  $s'_i(x)$ :

$$s_i'(x) = \frac{1}{2h_i} \left[ -(x_{i+1} - x)^2 M_i + (x - x_i)^2 M_{i+1} \right] - \frac{h_i}{6} \left( -M_i + M_{i+1} \right) + \frac{1}{h_i} \left( -f_i + f_{i+1} \right)$$

 $<sup>^{10}</sup>$ At each substitution, one cubic power vanishes and the other degenerates to  $h_i$ , which simplifies the calculation.

Replacing the subscript i by i-1, we have

$$s'_{i-1}(x) = \frac{1}{2h_{i-1}} \left[ -(x_i - x)^2 M_{i-1} + (x - x_{i-1})^2 M_i \right] - \frac{h_{i-1}}{6} \left( -M_{i-1} + M_i \right) + \frac{1}{h_{i-1}} \left( -f_{i-1} + f_i \right)$$

Imposing (3), i.e.  $s'_{i-1}(x_i) = s'_i(x_i)$ , we will have

$$h_{i-1}M_{i-1} + 2(h_{i-1} + h_i)M_i + h_iM_{i+1} = 6\frac{f_{i+1} - f_i}{h_i} - 6\frac{f_i - f_{i-1}}{h_{i-1}}$$

for i = 1, ..., n-1. This equation represents a system of n-1 linear equations with n+1 unkwowns; the other 2 equation is obtained from the boundary condition.

#### Natural boundary condition:

Under natural boundary condition,  $M_0 = M_n = 0$  by definition. Then we have

$$Am = b$$

where

$$\mathbf{A} = \begin{pmatrix} 2(h_0 + h_1) & h_1 \\ h_1 & 2(h_1 + h_2) & h_2 \\ & \ddots & \ddots & \ddots \\ & & h_{n-3} & 2(h_{n-3} + h_{n-2}) & h_{n-2} \\ & & & h_{n-2} & 2(h_{n-2} + h_{n-1}) \end{pmatrix}$$

$$\mathbf{m} = \begin{pmatrix} M_1 & M_2 & \cdots & M_{n-1} \end{pmatrix}^t$$

and

$$\mathbf{b} = 6 \times \begin{pmatrix} \frac{f_0}{h_0} - \frac{f_1}{h_0} - \frac{f_1}{h_1} + \frac{f_2}{h_1} \\ \vdots \\ \frac{f_{i-1}}{h_{i-1}} - \frac{f_i}{h_{i-1}} - \frac{f_i}{h_i} + \frac{f_{i+1}}{h_i} \\ \vdots \\ \frac{f_{n-2}}{h_{n-2}} - \frac{f_{n-1}}{h_{n-2}} - \frac{f_{n-1}}{h_{n-1}} + \frac{f_n}{h_{n-1}} \end{pmatrix}$$

#### Clamped boundary condition:

Under clamped boundary condition,  $s'_0(x_0) = f'(x_0)$  and  $s'_{n-1}(x_n) = f'(x_n)$ , which we will have, from  $s'_i(x)$  equation,

$$2h_0 M_0 + h_0 M_1 = 6\left(-\frac{f_0}{h_0} + \frac{f_1}{h_0} - f_0'\right)$$
$$h_{n-1} M_{n-1} + 2h_{n-1} M_n = 6\left(\frac{f_{n-1}}{h_{n-1}} - \frac{f_n}{h_{n-1}} + f_n'\right)$$

Therefore, we can form  $\mathbf{Am} = \mathbf{b}$  accordingly, where

$$\mathbf{A} = \begin{pmatrix} 2h_0 & h_0 \\ h_0 & 2(h_0 + h_1) & h_1 \\ & \ddots & \ddots & \ddots \\ & & h_{n-2} & 2(h_{n-2} + h_{n-1}) & h_{n-1} \\ & & & h_{n-1} & 2h_{n-1} \end{pmatrix}$$

$$\mathbf{m} = \begin{pmatrix} M_0 & M_1 & \cdots & M_n \end{pmatrix}^t$$

and

$$\mathbf{b} = 6 \times \begin{pmatrix} -\frac{f_0}{h_0} + \frac{f_1}{h_0} - f'_0 \\ \frac{f_0}{h_0} - \frac{f_1}{h_0} - \frac{f_1}{h_1} + \frac{f_2}{h_1} \\ \vdots \\ \frac{f_{i-1}}{h_{i-1}} - \frac{f_i}{h_{i-1}} - \frac{f_i}{h_i} + \frac{f_{i+1}}{h_i} \\ \vdots \\ \frac{f_{n-2}}{h_{n-2}} - \frac{f_{n-1}}{h_{n-2}} - \frac{f_{n-1}}{h_{n-1}} + \frac{f_n}{h_{n-1}} \\ \frac{f_{n-1}}{h_{n-1}} - \frac{f_n}{h_{n-1}} + f'_n \end{pmatrix}$$