Chapter 1. Preliminaries

Numerical Analysis, a branch of Applied Mathematics situated at the border between Mathematics and Computer Science, produces methods and procedures for finding numerical solutions of various problems, with a given precision. Standard topics in a Numerical Analysis course are the approximation of problems by simpler problems, the construction of algorithms, iteration methods, error analysis, stability, asymptotic error formulas, the effects of machine arithmetic, etc.

The study will focus on issues such as:

- Problems modeled by functions that do not have an analytical expression, whose values are known only at a discrete set of points. Based on these, we want to approximate values of the function at new points, values of the derivatives or integrals of the function, etc. Such problems lead to finite and divided differences, interpolation formulas, numerical differentiation and integration schemes and many others.
- In many practical situations it is necessary to solve various types of equations or systems of equations, such as: algebraic, transcendental, differential, or integral equations, whose exact solutions cannot be found. They need to be approximated numerically.

An approximating procedure must satisfy several properties:

- 1. To be *convergent*, meaning the sequence of iterations (successive approximations) must converge to the exact solution, in order to produce "better and better" approximations.
- 2. To be *stable* (to have stability), meaning that small variations of the input data should lead to small variations in the results (the approximating solutions).
- 3. From its structure and properties, to be able to estimate the *error* and the *speed of convergence* of the approximating method.

Many times, the conditions that ensure stability of a numerical method coincide with the ones that guarantee its convergence.

1 Taylor Polynomials

We start with a very useful tool from Calculus, Taylor's theorem. This will be needed for both the development and understanding of many of the numerical methods discussed later on.

In fact, Taylor polynomials give the very first example of a numerical method, being used as a way to evaluate other functions approximately.

Theorem 1.1. [Taylor's Theorem] Let $f : [a, b] \to \mathbb{R}$ be a function with n + 1 continuous derivatives on [a, b], for some $n \ge 0$ and let $x, x_0 \in [a, b]$. Then

$$f(x) = T_n(x) + R_n(x),$$
 (1.1)

where

$$T_n(x) = f(x_0) + \frac{(x - x_0)}{1!} f'(x_0) + \dots + \frac{(x - x_0)^n}{n!} f^{(n)}(x_0)$$
(1.2)

is Taylor's polynomial of degree n of f at x_0 and

$$R_n(x) = \frac{1}{n!} \int_{x_0}^x (x-t)^n f^{(n+1)}(t) dt = \frac{(x-x_0)^{n+1}}{(n+1)!} f^{(n+1)}(\xi), \ \xi \text{ between } x \text{ and } x_0, \ (1.3)$$

is the error of the Taylor approximation.

Example 1.2. Using Taylor's theorem with $x_0 = 0$, we obtain the following well-known formulas:

$$e^{x} = 1 + x + \frac{x^{2}}{2!} + \dots + \frac{x^{n}}{n!} + \frac{x^{n+1}}{(n+1)!}e^{\xi_{x}},$$

$$\cos x = 1 - \frac{x^{2}}{2!} + \frac{x^{4}}{4!} - \dots + (-1)^{n}\frac{x^{2n}}{(2n)!} + (-1)^{n+1}\frac{x^{2n+2}}{(2n+2)!}\cos\xi_{x},$$

$$\sin x = x - \frac{x^{3}}{3!} + \frac{x^{5}}{5!} - \dots + (-1)^{n-1}\frac{x^{2n-1}}{(2n-1)!} + (-1)^{n}\frac{x^{2n+1}}{(2n+1)!}\sin\xi_{x},$$

$$1 + x)^{a} = 1 + {\binom{a}{1}}x + {\binom{a}{2}}x^{2} + \dots + {\binom{a}{n}}x^{n} + {\binom{a}{n+1}}\frac{x^{n+1}}{(1+\xi_{x})^{n+1-a}},$$

with

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$$\left(\begin{array}{c}a\\k\end{array}\right) = \frac{a(a-1)\cdots(a-k+1)}{k!}, \ k=1,2,3,\ldots, \ a\in\mathbb{R}$$

An important special case of the last formula is a = -1, with x replaced by -x:

$$\frac{1}{1-x} = 1 + x + x^2 + \dots + x^n + \frac{x^{n+1}}{1-x},$$

where the remainder has a simpler form than before. This is easily proved by multiplying both sides by 1 - x and then simplifying. Rearranging the terms, we obtain the familiar formula for a partial sum of the Geometric series:

$$1 + x + x^{2} + \dots + x^{n} = \frac{1 - x^{n+1}}{1 - x}, \ x \neq 1.$$

Series representations for the functions in Example 1.2 can be obtained by letting $n \to \infty$. Recall that the series for the first three functions converge for all $x \in \mathbb{R}$, while those for the last two converge for |x| < 1. So, by taking Taylor polynomials of higher and higher degree, we can obtain better and better approximations of the functions above.



Fig. 1: Taylor approximations of e^x

Example 1.3. Consider the function $f(x) = e^x$. Figure 1 shows the approximations of f with Taylor polynomials of degree 1, 2, 3 and 4, for $x \in [-1,1]$. The errors of these approximations, $\max_{x \in [-1,1]} \{e^x - p_n(x)\}$, are graphed in Figure 2.



Fig. 2: Errors in Taylor approximations of e^x

Taylor's formula in two dimensions

Theorem 1.4. Let $f : D \subset \mathbb{R}^2 \to \mathbb{R}^2$ be a function which is n + 1 times continuously differentiable on D, for some $n \ge 0$ and let $(x, y), (x_0, y_0) \in D$. Then

$$f(x,y) = T_n(x,y) + R_n(x,y),$$
 (1.4)

where

$$T_{n}(x,y) = f(x_{0},y_{0}) + \frac{x-x_{0}}{1!}f'_{x}(x_{0},y_{0}) + \frac{y-y_{0}}{1!}f'_{y}(x_{0},y_{0}) + \frac{1}{2!}\left[(x-x_{0})^{2}f''_{x^{2}}(x_{0},y_{0}) + 2(x-x_{0})(y-y_{0})f''_{xy}(x_{0},y_{0}) + (y-y_{0})^{2}f''_{y^{2}}(x_{0},y_{0})\right] + \sum_{j=1}^{n}\frac{1}{j!}\left[(x-x_{0})\frac{\partial}{\partial x} + (y-y_{0})\frac{\partial}{\partial y}\right]^{j}f(x,y)\bigg|_{\substack{x=x_{0}\\y=y_{0}}}$$
(1.5)

and

$$R_n(x,y) = \frac{1}{(n+1)!} \left[(x-x_0) \frac{\partial}{\partial x} + (y-y_0) \frac{\partial}{\partial y} \right]^{n+1} f(x,y) \bigg|_{\substack{x=\xi\\y=\eta}},$$
(1.6)

with (ξ, η) a point on the line segment determined by the points (x, y) and (x_0, y_0) .

2 Errors: Sources, Propagation, Analysis

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2.1 Types of Numerical Problems

Let us first consider the following simple examples:

Example 2.1. Compute

$$\int_{1}^{3} \frac{1}{x} \, dx.$$

Solution. Its exact value is

$$\int_{1}^{3} \frac{1}{x} dx = \ln x \Big|_{1}^{3} = \ln 3 - \ln 1 = \ln 3$$

and its approximation is

$$\int_{1}^{3} \frac{1}{x} \, dx = 1.0986...$$

Example 2.2. Solve the equation

$$x^2 = 3, x > 0.$$

Solution. The true solution is

$$x = \sqrt{3},$$

with approximating value

$$x = 1.732...$$

Example 2.3. Consider the data

x_i	1	2	3	4	5	6	7
y_i	5	13	16	23	33	38	40

Discuss the nature of the relationship between x and y.

Solution. We plot the data (see Figure 3).



Fig. 3: Scatterplot and linear function

Notice that a straight line "best" approximates the data. So based on these values, we seek a relationship between x and y of the form

$$f(x) = ax + b.$$

In the next chapters, we will analyze several rigorous procedures for approximating scattered data.

From these examples, we see that a numerical problem is, in general, of the form

$$f(x) = y$$

- If x and f are given, and we seek y, then this is called a **direct problem** or an **evaluation problem** (Example 2.1).
- If y and f are given, and we want to find x, then it is called an **inverse problem** (Example 2.2).
- If x and y are given, and f must be determined, then we have an **identifying problem** (Example 2.3).

For each type of problem, we obtain an *approximating* value, which is affected by *errors*, perturbations from the true value. The study of errors and their propagation is an important task in Numerical Analysis. In practical problems, it is important that the approximations obtained have *small (negligible)* errors, that do not affect the overall precision of the numerical procedure.

2.2 Sources and Propagation of Errors

Let $\mathcal{A} : \mathbb{R} \to \mathcal{P}(\mathbb{R})$ be a mapping that assigns a set $\mathcal{A}(x) \subseteq \mathbb{R}$ to each real number $x \in \mathbb{R}$.

Definition 2.4. Let $x \in \mathbb{R}$. The number x^* is called an **approximation** of $x \in \mathbb{R}$, if $x^* \in \mathcal{A}(x)$ (notation $x \approx x^*$). The mapping \mathcal{A} is called an **approximating procedure (method)**.

In practice, $\mathcal{A}(x)$ consists of numbers that vary "little" from x.

Let $x \in \mathbb{R}$ and $x^* \in \mathcal{A}(x)$ be an approximation of the true (exact) value x.

Definition 2.5. The number

$$\Delta x = x - x^* \tag{2.1}$$

is called the **error** of the approximation x^* .

This quantity is also called the *absolute error*. If $\Delta x > 0$, then x^* is an **under-approximation**, and if $\Delta x < 0$, x^* is an **over-approximation**.

For example, for the number $\pi = 3.141592...$, the number $x^* = 3.14$ is an under-approximation, while $x^* = 3.142$ is an over-approximation.

Definition 2.6. *The quantity*

$$\delta x = \frac{\Delta x}{x}, \ x \neq 0 \tag{2.2}$$

is called relative error.

This gives the size of the error in proportion to the true value being approximated. From (2.2), we get

$$\Delta x = x \delta x \Longrightarrow x^* - x = x \delta x,$$

or

$$x^* = (1 + \delta x)x, \tag{2.3}$$

which is widely used in applications.

Since in practice x is unknown, we use instead

$$\delta x = \frac{\Delta x}{x*}.$$
(2.4)

For the absolute and relative error, one can also use the notations Δx^* and δx^* .

Sources of error

- *Mathematical modeling of a physical problem*. A mathematical model for a physical situation is an attempt to give mathematical relationships between certain quantities of physical interest. Because of the complexity of physical reality, a variety of simplifying assumptions are used to construct a more tractable mathematical model. The resulting model has limitations on its accuracy as a consequence of these assumptions, and these limitations may or may not be troublesome, depending on the uses of the model.
- *Blunders (human errors)*. In pre-computer times, chance arithmetic errors were always a serious problem. With the introduction of digital computers, the type of blunder has changed. Chance arithmetic errors (e.g., computer malfunctioning) are now relatively rare, and programming errors are currently the main difficulty. Often a program error will be repeated many times in the course of executing the program, and its existence becomes obvious because of absurd numerical output (although the source of the error may still be difficult to find). This makes good program debugging very important, even though it may not seem very rewarding immediately.
- *Uncertainty in physical data*. Most data from a physical experiment will contain error or uncertainty within it. This must affect the accuracy of any calculations based on the data, limiting the accuracy of the answers.
- *Machine errors*. This means the errors inherent in using the floating-point representation of numbers. Specifically, we mean the rounding/chopping errors and the underflow/overflow errors. The rounding/chopping errors are due to the finite length of the floating-point mantissa; and these errors occur with all of the computer arithmetic operations.
- *Mathematical truncation error*. This name refers to the error of approximation in numerically solving a mathematical problem, and it is the error generally associated with the subject of Numerical Analysis. It involves the approximation of infinite processes by finite ones, replacing noncomputable problems with computable ones, etc.

2.3 **Propagation of Errors**

We distinguish two types of problems:

Case 1. Given the errors in the approximations of input data, find the errors in the output. We start with the case of a function of two variables $f : D \to \mathbb{R}$, $D = \{(x, y) | x, y \in \mathbb{R}\}$. Let (x^*, y^*) be the approximating values of (x, y). Their absolute errors are then

$$\Delta x = x - x^* \Rightarrow x = x^* + \Delta x,$$

$$\Delta y = y - y^* \Rightarrow y = y^* + \Delta y.$$

We want to compute the absolute error

$$\Delta f = f(x, y) - f(x^*, y^*)$$
(2.5)

and the relative error δf .

We use Taylor's formula in two variables (1.5), at x^* and y^* . We have:

$$f(x^* + \Delta x, y^* + \Delta y) = f(x^*, y^*) + \Delta x f'_x(x^*, y^*) + \Delta y f'_y(x^*, y^*) + \frac{(\Delta x)^2}{2!} f''_{x^2}(x^*, y^*) + 2 \frac{\Delta x \Delta y}{2!} f''_{xy}(x^*, y^*) + \frac{(\Delta y)^2}{2!} f''_{y^2}(x^*, y^*) + \dots,$$
(2.6)

or

$$\Delta f = \Delta x f'_x(x^*, y^*) + \Delta y f'_y(x^*, y^*) + \frac{1}{2!} \Big[(\Delta x)^2 f''_{x^2}(x^*, y^*) + 2\Delta x \Delta y f''_{xy}(x^*, y^*) + (\Delta y)^2 f''_{y^2}(x^*, y^*) \Big] + \dots$$
(2.7)

If Δx and Δy are small, then $(\Delta x)^2$, $\Delta x \Delta y$ and $(\Delta y)^2$ can be neglected. Then we find the **absolute** error of function f or the maximum absolute error:

$$\Delta f \simeq \Delta x f'_x(x^*, y^*) + \Delta y f'_y(x^*, y^*).$$
(2.8)

In general, for a function of n variables, we have

$$\Delta f \simeq \Delta x_1 f'_{x_1}() + \Delta x_2 f'_{x_2}() + \dots + \Delta x_n f'_{x_n}() = \sum_{i=1}^n \Delta x_i f'_{x_i}().$$
(2.9)

This is the **propagated error**. Then the relative error δf is

$$\delta f = \frac{\Delta f}{f} \simeq \sum_{i=1}^{n} \Delta x_i \frac{f'_{x_i}(f)}{f} = \sum_{i=1}^{n} \Delta x_i \frac{d}{dx_i} \ln f(f)$$
$$= \sum_{i=1}^{n} x_i \delta x_i \frac{d}{dx_i} \ln f(f) = \sum_{i=1}^{n} x_i \left(\frac{d}{dx_i} \ln f(f)\right) \delta x_i.$$
(2.10)

Case 2. The inverse problem is to determine the precision needed in the input data that will guarantee a (given, preset) accuracy in the output data.

To do this, we use the so-called *principle of equal effects*. This assumes that all terms $f'_{x_i}\Delta x_i$ in (2.9) have the *same effect*, i. e.

$$f'_{x_1}\Delta x_1 = f'_{x_2}\Delta x_2 = \dots = f'_{x_n}\Delta x_n.$$

Then (2.9) becomes

$$\Delta f \simeq n \Delta x_i f'_{x_i}(),$$

from which it follows that the absolute error is

$$\Delta x_i \simeq \frac{\Delta f}{n f'_{x_i}} \tag{2.11}$$

and the *relative error* is given by

$$\delta x_i \simeq \frac{\delta f}{n x_i \frac{d}{dx_i} \ln f} \tag{2.12}$$

3 Floating-Point Representation of Numbers. Significant Digits

Definition 3.1. A number r written in base b (an even number) has the **floating-point representation** as

$$r = \pm r_0 r_1 \dots r_p \cdot r_{p+1} \dots r_k \times b^e$$

where r_0, r_1, \ldots, r_k form the mantissa (significand), e is the exponent and b is the base.

In order to have uniqueness of the representation, the floating-point numbers are *normalized*, that is, we change the representation, not the value, such that $r_0 \neq 0$. The term *floating-point number* will be used to mean a real number that can be exactly represented in this format.

Definition 3.2. The significant digits of a floating-point number written in base b are any of the digits 0, 1, 2, ..., b - 1 in its representation that are non-zero, or located between non-zero digits, or preceded by at least one non-zero digit.

So 0 can be a significant digit if it does more than just indicate the floating decimal point or fills the places of unknown or omitted digits. The leftmost significant digit is called the *most significant* digit.

For example,

- the number 7063 has all significant digits;
- the number 0.02340 the last 4 digits are significant (the zeros in front of 2 merely indicate the decimal point, and are, therefore, *not* significant);
- $1.230 \times 10^3 = 1230 \text{zero}$ is a significant digit, as it is preceded by a nonzero digit.

Consider the number r > 0 with the following representation in base 10:

$$r = r_0 10^k + r_1 10^{k-1} + \dots + r_{n-1} 10^{k-n+1} + r_n 10^{k-n} + \dots$$

Definition 3.3. *The number*

$$r^* = r_0^* 10^k + r_1^* 10^{k-1} + \dots + r_{n-1}^* 10^{k-n+1}$$

approximates r correctly with n significant digits if

$$|\Delta r^*| \le \frac{1}{2} \times 10^{k-n+1}$$
 (3.1)

Example 3.4. Find the number of significant digits with which $e^* = 2.718282$ approximates correctly the number e = 2.71828182...

Solution. Let $e^* = 2.718282$. This can be written as

$$e^* = 2 \cdot 10^0 + 7 \cdot 10^{-1} + 1 \cdot 10^{-2} + \cdots$$

So k = 0. Then

$$\Delta e^* = e^* - e = 0.00000018 = 0.18 \times 10^{-6} \le \frac{1}{2} \times 10^{-6}.$$

Comparing it to (3.1), we get

$$10^{k-n+1} = 10^{-6} \Rightarrow 0 - n + 1 = -6 \Rightarrow n = 7.$$

Thus, the number $e^* = 2.718282$ approximates correctly e = 2.71828182... with 7 significant digits.

The following relation exists between the number of significant digits and the relative error:

$$\delta r^* \leq \frac{1}{r_0^* \times 10^{n-1}},$$
(3.2)

where r^* is the correct approximation with n significant digits, with the normalized representation

$$r^* = r_0^* 10^k + r_1^* 10^{k-1} + \dots + r_{n-1}^* 10^{k-n+1}.$$
(3.3)

This can be considered the maximum relative error.

4 Stability and Conditioning

A number of mathematical problems have solutions that are quite sensitive to small computational errors, for example rounding errors. To deal with this phenomenon, we use the concepts of *stability* and *condition (conditioning) number*.

Consider a general problem of the type

$$y = f(x), f : \mathbb{R}^m \to \mathbb{R}^n.$$

We are interested in how "sensitive" is f to small perturbations of x, that is, do small perturbations of x lead to small perturbations in y – which means the problem is *stable*, or not – *unstable* problem. In the latter case, approximation is not very helpful. We would like to "measure" the degree of sensitivity (how stable or unstable a problem is) by a single number, called the **condition number** of f at x.

The function f is assumed to be evaluated exactly, with infinite precision, as we perturb x. The condition number of f, therefore, is an inherent property of the map f and does not depend on any algorithmic considerations concerning its implementation.

Recall from the propagation of error formula (2.10) that when $f : \mathbb{R}^m \to \mathbb{R}$, we found

$$\delta f \approx \sum_{i=1}^m x_i \frac{f'_{x_i}}{f} \delta x_i.$$

Now, for the general case $f : \mathbb{R}^m \to \mathbb{R}^n, x = [x_1, \dots, x_m]^T \in \mathbb{R}^m, f(x) = (f_1(x), \dots, f_n(x))^T \in \mathbb{R}^n$, let

$$\gamma_{ij} = (\operatorname{cond}_{ij}f)(x) = \frac{x_i \frac{\partial f_j}{\partial x_i}}{f_j(x)}, \ i = \overline{1, m}, \ j = \overline{1, n}$$

and

$$\Gamma(x) = [\gamma_{ij}] = \begin{bmatrix} \frac{x_1 \frac{\partial f_1}{\partial x_1}}{f_1(x)} & \cdots & \frac{x_m \frac{\partial f_1}{\partial x_m}}{f_1(x)} \\ \vdots & & \vdots \\ \frac{x_1 \frac{\partial f_n}{\partial x_1}}{f_n(x)} & \cdots & \frac{x_m \frac{\partial f_n}{\partial x_m}}{f_n(x)} \end{bmatrix},$$
(4.1)

called the **conditioning matrix**. Then, the **condition number** of f at x is defined by

$$(\text{cond } f)(x) = ||\Gamma(x)||,$$
 (4.2)

for a matrix norm $|| \cdot ||$. If f is a linear function, then

$$(\operatorname{cond} f)(x) = \frac{||x|| \left| \left| \frac{\partial f}{\partial x} \right| \right|}{||f(x)||}.$$
(4.3)

Particular cases for m = n = 1

1. If $x \neq 0, f(x) \neq 0$, then

$$(\operatorname{cond} f)(x) = \left| \frac{xf'(x)}{f(x)} \right|.$$

2. If $x = 0, f(x) \neq 0$, then we take only absolute error for x,

$$(\operatorname{cond} f)(x) = \left| \frac{f'(x)}{f(x)} \right|.$$

3. If $x \neq 0$, f(x) = 0, then we take only absolute error for f(x),

$$(\operatorname{cond} f)(x) = |xf(x)|.$$

4. If x = f(x) = 0, then

$$(\operatorname{cond} f)(x) = |f'(x)|$$

Ill-conditioned and ill-posed problems

If the condition number of a problem is large $((\text{cond } f)(x) \gg 1)$, then even for small (relative) errors in the input data, we can expect large errors in the output data. Such problems are called **ill-conditioned** problems. There is no clear line between ill- and well-conditioned problems, it all depends on precision specifications.

If the result of a mathematical problem depends in a discontinuous way on data that varies continuously, then it is impossible to give an accurate numerical solution in a neighborhood of the discontinuity. In such cases, the result can be significantly perturbed, even if the input data is exact and we use high precision procedures. These problems are called **ill-posed** problems. An ill-posed problem can appear if, for example, an integer result is computed from real data (which varies continuously).

Example 4.1.

1. Number of distinct real roots of a polynomial. Let

$$p(x,c) = x^3 - 2x^2 + x + c, \ c > -\frac{4}{27}.$$

The polynomial p can have one real root (if c > 0), two distinct real roots (for c = 0), or three (if c < 0). Therefore, for values of c close to zero, the number of distinct real zeros of p is an *ill-posed* problem.

2. Recurrence relations. Let

$$I_n = \int_0^1 \frac{t^n}{t+5} dt, \ n \in \mathbb{N}.$$

Individual values of this integral can be computed exactly or numerically, e.g. $I_0 = \ln \frac{6}{5}$, $I_{100} \approx 0.0017$. What about a recurrence relation?

Writing

$$\frac{t^n}{t+5} = t^{n-1} - 5\frac{t^{n-1}}{t+5} \text{ or } \frac{t^n}{t+5} = \frac{1}{5}\left(t^n - 5\frac{t^{n+1}}{t+5}\right),$$

we can find *direct recurrence*

$$I_n = \frac{1}{n} - 5I_{n-1} = f_n(I_0),$$

or inverse recurrence

$$I_n = \frac{1}{5} \left(\frac{1}{n+1} - 5I_{n+1} \right) = g_n(I_{100}).$$

Then it can be shown that

$$(\text{cond } f_n)(I_0) = \mathcal{O}(5^n) \text{ and } (\text{cond } g_n)(I_{100}) = \mathcal{O}(5^{100-n}),$$

which means that direct recurrence is an *ill-conditioned* problem, whereas inverse recurrence is *well-conditioned*.

Remark 4.2. In the next chapter, we will consider in greater detail the conditioning of algebraic linear systems and matrices.

5 Divided and Finite Differences

These are expressions that are helpful in writing, computing and implementing various iterative numerical procedures.

5.1 Divided Differences

Definition 5.1. Let $f : [a, b] \to \mathbb{R}$ be a differentiable function on [a, b], and $x_i \in [a, b]$, $i = \overline{0, n}$, be n + 1 distinct nodes. The quantity

$$f[x_0, x_1] = \frac{f(x_1) - f(x_0)}{x_1 - x_0}.$$
(5.1)

is called the first-order divided difference of f at the nodes x_0 and x_1 .

Remark 5.2.

1. An alternative notation is $[x_0, x_1; f]$.

2. The first-order divided difference of a function can be thought of as a *discrete* version of the derivative.

3. If we consider $f[x_0] = f(x_0)$ the divided difference of order 0 at x_0 , then (5.1) can be written as

$$f[x_0, x_1] = \frac{f[x_1] - f[x_0]}{x_1 - x_0}.$$
(5.2)

We define higher-order divided differences recursively using lower-order ones.

Definition 5.3. The divided difference of order n of f at the distinct nodes x_0, x_1, \ldots, x_n is the quantity

$$f[x_0, x_1, \dots, x_n] = \frac{f[x_1, x_2, \dots, x_n] - f[x_0, x_1, \dots, x_{n-1}]}{x_n - x_0}.$$
(5.3)

Remark 5.4.

1. The denominator in (5.3) is the difference between the nodes that are *not* common to the differences at the numerator.

2. For easy computation (and implementation) of divided differences, we generate the *table of divided differences*, illustrated below for 4 nodes. The divided differences are obtained on the first row.

Example 5.5. Let $f(x) = \sin \pi x$, and the nodes $x_0 = 0$, $x_1 = \frac{1}{6}$, $x_2 = \frac{1}{2}$. Let us construct the divided difference table.

Solution.

Divided differences with multiple nodes

Divided differences with multiple nodes can be expressed in terms of the derivatives of the function f, as follows

$$f[x_0, x_0] = \lim_{x_1 \to x_0} f[x_0, x_1] = \lim_{x_1 \to x_0} \frac{f(x_1) - f(x_0)}{x_1 - x_0} = f'(x_0)$$

In general, the **divided difference of order** n at the node x_0 , of multiplicity n + 1, is defined as

$$f[\underbrace{x_0, x_0, \dots, x_0}_{n+1}] = \frac{f^{(n)}(x_0)}{n!},$$
(5.4)

,

and further, for mixed nodes (some simple, some multiple), we use definition 5.3.

Example 5.6. For a function f, construct the divided differences table for the double node x_0 and the simple node x_1 .

Solution. The new nodes are

$$z_0 = x_0, \ z_1 = x_0, \ \text{and} \ z_2 = x_1.$$

The divided differences table:

$$z_{0} = x_{0} \quad f[z_{0}] = f(x_{0}) \longrightarrow f'(x_{0}) \longrightarrow \frac{f(x_{1}) - f(x_{0})}{x_{1} - x_{0}} - f'(x_{0})$$

$$z_{1} = x_{0} \quad f[z_{1}] = f(x_{0}) \longrightarrow \frac{f(x_{1}) - f(x_{0})}{x_{1} - x_{0}}$$

$$z_{2} = x_{1} \quad f[z_{2}] = f(x_{1})$$

Example 5.7. Let us see the divided differences table for 3 double nodes.

Solution. We have the nodes x_0, x_1, x_2 and the values $f(x_i), f'(x_i), i = 0, 1, 2$. We define the sequence of nodes z_0, z_1, \ldots, z_5 by

$$z_{2i} = z_{2i+1} = x_i, i = 0, 1, 2.$$

We build the divided difference table relative to the nodes $z_i, i = \overline{0, 5}$.

Since $z_{2i} = z_{2i+1} = x_i$ for every $i = 0, 1, 2, f[z_{2i}, z_{2i+1}] = f[x_i, x_i]$ is a divided difference with a double node and it is equal to $f'(x_i)$; therefore we will use $f'(x_0), f'(x_1), f'(x_2)$ instead of first order divided differences $f[z_0, z_1], f[z_2, z_3], f[z_4, z_5]$.

$$z_{0} = x_{0} \quad f[z_{0}] \longrightarrow f[z_{0}, z_{1}] = f'(x_{0}) \longrightarrow f[z_{0}, z_{1}, z_{2}] \dots$$

$$z_{1} = x_{0} \quad f[z_{1}] \longrightarrow f[z_{1}, z_{2}] = \frac{f(x_{1}) - f(x_{0})}{x_{1} - x_{0}} \longrightarrow f[z_{1}, z_{2}, z_{3}] \dots$$

$$z_{2} = x_{1} \quad f[z_{2}] \longrightarrow f[z_{2}, z_{3}] = f'(x_{1}) \longrightarrow f[z_{2}, z_{3}, z_{4}] \dots$$

$$z_{3} = x_{1} \quad f[z_{3}] \longrightarrow f[z_{3}, z_{4}] = \frac{f(x_{2}) - f(x_{1})}{x_{2} - x_{1}} \longrightarrow f[z_{3}, z_{4}, z_{5}] \dots$$

$$z_{4} = x_{2} \quad f[z_{4}] \longrightarrow f[z_{4}, z_{5}] = f'(x_{2})$$

$$z_{5} = x_{2} \quad f[z_{5}]$$