

# Métodos Numéricos

Mercedes Román Ruiz

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# Capítulo 1

## Introducción

### 1. Vector Spaces

**Definition 1.1.** A vector space over the numeric field  $K$  ( $K = \mathbb{R}$  or  $K = \mathbb{C}$ ) is a nonempty set  $V$ , whose elements are called vector and in which two operations are defined, called addition and scalar multiplication, that enjoy the following properties:

1. addition is commutative and associative;
2. there exists an element  $0 \in V$  (the zero vector or null vector) such that  $v + 0 = v$  for each  $v \in V$
3.  $0 \cdot v = 0$ ,  $1 \cdot v = v$ , for each  $v \in V$ , where  $0$  and  $1$  are respectively the zero and the unity of  $K$ ;
4. for each element  $v \in V$  there exists its opposite,  $-v$ , in  $V$  such that  $v + (-v) = 0$ ;
5. the following distributive properties hold

$$\forall \alpha \in K, \forall v, w \in V, \alpha(v + w) = \alpha v + \alpha w$$

$$\forall \alpha, \beta \in K, \forall v \in V, (\alpha + \beta)v = \alpha v + \beta v$$

6. the following associative property holds

$$\forall \alpha, \beta \in K, \forall v \in V, (\alpha\beta)v = \alpha(\beta v)$$

**Definition 1.2.** We say that a nonempty part  $W$  of  $V$  is a vector subspace of  $V$  if  $W$  is a vector space over  $K$ .

**Definition 1.3.** A system of vector  $v_1, \dots, v_n$  of a vector space  $V$  is called linearly independent if the relation

$$\alpha_1 v_1 + \alpha_2 v_2 + \dots + \alpha_n v_n = 0$$

with  $\alpha_1, \alpha_2, \dots, \alpha_n \in K$  implies that  $\alpha_1 = \alpha_2 = \dots = \alpha_n = 0$ . Otherwise, the system will be called linearly dependent.

We call a basis of  $V$  any system of linearly independent generator of  $V$ . If  $u_1, \dots, u_n$  is a basis of  $V$ , the expression  $v = v_1 u_1 + \dots + v_n u_n$  is called the decomposition of  $v$  with respect to the basis and the scalars  $v_1, \dots, v_n \in K$  are the components of  $v$  with respect to the given basis. Moreover, the following pro holds.

**Property 1.1.** Let  $V$  be a vector space which admits a basis of  $n$  vectors. Then every system of linearly independent vector of  $V$  has at most  $n$  elements and any other basis of  $V$  has  $n$  elements. The number  $n$  is called the dimension of  $V$  and we write  $\dim(V) = n$ . If, instead, for any  $n$  there always exist  $n$  linearly independent vectors of  $V$ , the vector space is called infinite dimensional.

### 2. Matrices

Let  $m$  and  $n$  be two positive integers. We call a matrix having  $m$  rows and  $n$  columns, or a matrix  $m \times n$ , or a matrix  $(m, n)$ , with elements in  $K$ , a set of  $mn$  scalars  $a_{ij} \in K$ , with  $i = 1, \dots, m$  and  $j = 1, \dots, n$ , represented in the following rectangular array

$$A = \begin{bmatrix} a_{11} & a_{12} & \cdots & a_{1n} \\ a_{21} & a_{22} & \cdots & a_{2n} \\ \vdots & \vdots & \ddots & \vdots \\ a_{m1} & a_{m2} & \cdots & a_{mn} \end{bmatrix} \quad (1.1)$$

When  $K = \mathbb{R}$  or  $K = \mathbb{C}$  we shall respectively write  $A \in \mathbb{R}^{m \times n}$  or  $A \in \mathbb{C}^{m \times n}$ , to explicitly outline the numerical fields which the elements of  $A$  belong to. Capital letters will be used to denote the matrices, while the lower case letters corresponding to those upper case letters will denote the matrix entries.

We shall abbreviate (1.1) as  $A = (a_{ij})$  with  $i = 1, \dots, m$  and  $j = 1, \dots, n$ . The index  $i$  is called row index, while  $j$  is the column index. The set  $(a_{i1}, a_{i2}, \dots, a_{in})$  is called the  $i$ -th row of  $A$ ; likewise,  $(a_{1j}, a_{2j}, \dots, a_{mj})$  is the  $j$ -th column of  $A$ .

If  $n = m$  the matrix is called squared or having order  $n$  and the set of the entries  $(a_{11}, a_{22}, \dots, a_{nn})$  is called its main diagonal.

A matrix having one row or one column is called a row vector or column vector respectively. Unless otherwise specified, we shall always assume that a vector is a column vector. In the case  $n = m = 1$ , the matrix will simply denote a scalar of  $K$ .

**Definition 2.1.** Let  $A$  be a matrix  $m \times n$ . Let  $1 \leq i_1 < i_2 < \dots < i_k \leq m$  and  $1 \leq j_1 < j_2 < \dots < j_l \leq n$  two sets of contiguous indexes. The matrix  $S(k \times l)$  of entries  $s_{pq} = a_{i_p j_q}$  with  $p = 1, \dots, k, q = 1, \dots, l$  is called a submatrix of  $A$ . If  $k = l$  and  $i_r = j_r$  for  $r = 1, \dots, k$ ,  $S$  is called a principal submatrix of  $A$ .

**Definition 2.2.** A matrix  $A(m \times n)$  is called block partitioned or said to be partitioned into submatrices if

$$A = \begin{bmatrix} A_{11} & A_{12} & \cdots & A_{1l} \\ A_{21} & A_{22} & \cdots & A_{2l} \\ \vdots & \vdots & \ddots & \vdots \\ A_{k1} & A_{k2} & \cdots & A_{kl} \end{bmatrix}$$

where  $A_{ij}$  are submatrices of  $A$ .

### 3. Operations with Matrices

#### 3.1. Matrices and Linear Mappings

**Definition 3.1.** A linear map for  $\mathbb{C}^n$  into  $\mathbb{C}^m$  is a function  $f : \mathbb{C}^n \rightarrow \mathbb{C}^m$  such that  $f(\alpha x + \beta y) = \alpha f(x) + \beta f(y)$ ,  $\forall \alpha, \beta \in K$  and  $\forall x, y \in \mathbb{C}^n$ .

The following result links matrices and linear maps.

**Property 3.1.** Let  $f : \mathbb{C}^n \rightarrow \mathbb{C}^m$  be a linear map. Then, there exists a unique matrix  $A_f \in \mathbb{C}^{m \times n}$  such that

$$f(x) = A_f x \quad \forall x \in \mathbb{C}^n \quad (1.2)$$

Conversely, if  $A_f \in \mathbb{C}^{m \times n}$  then the function defined in (1.2) is a linear map from  $\mathbb{C}^n$  into  $\mathbb{C}^m$ .

### 4. Well-posedness and Condition Number of a Problem

Consider the following problem: find  $x$  such that

$$F(x, d) = 0 \quad (1.3)$$

where  $d$  is the set of data which the solution depends on and  $F$  is the functional relation between  $x$  and  $d$ . According to the kind of problem that is represented in (1.3), the variables  $x$  and  $d$  may be real numbers, vectors or functions. Typically, (1.3) is called a direct problem if  $F$  and  $d$  are given and  $x$  is unknown, inverse problem if  $F$  and  $x$  are known and  $d$  is the unknown, identification problem when  $x$  and  $d$  are given while the functional relation  $F$  is the unknown.

Problem (1.3) is well posed if it admits a unique solution  $x$  which depends with continuity on the data. We shall use the terms well posed and stable in an interchanging manner and we shall deal henceforth only with well-posed problems.

A problem which does not enjoy the property above is called ill posed or unstable and before undertaking its numerical solution it has to be regularized, that is, it must be suitably transformed into a well-posed problem. Indeed, it is not appropriate to pretend the numerical method can cure the pathologies of an intrinsically ill-posed problem.

Let  $D$  be the set of admissible data, i.e. the set of the values of  $d$  in correspondance of which problem (1.3) admits a unique solution. Continuous dependence on the data means that small perturbations on the data  $d$  of  $D$  yield "small" changes in the solution  $x$ . Precisely, let  $d \in D$  and denote by  $\delta d$  a perturbation admissible in the sense that  $d + \delta d \in D$  and by  $\delta x$  the corresponding change in the solution, in such a way that

$$F(x + \delta x, d + \delta d) = 0 \quad (1.4)$$

Then, we require that

$$\exists \eta_0 = \eta_0(d) > 0, \quad \exists K_0 = K_0(d) \text{ such that if } \|\delta d\| \leq \eta_0 \text{ then } \|\delta x\| \leq K_0 \|\delta d\| \quad (1.5)$$

The norms used for the data and for the solution may not coincide, whenever  $d$  and  $x$  represent variables of different kinds.

**Remark 4.1.** The property of continuous dependence on the data could have been stated in the following alternative way, which is more akin to the classical form of Analysis  $\forall \epsilon > 0 \exists \delta = \delta(\epsilon)$  such that if  $\|\delta d\| \leq \delta$  then  $\|\delta x\| \leq \epsilon$ .

The form (1.5) is however more suitable to express in the following the concept of numerical stability, that is, the property that small perturbations on the data yield perturbations of the same order on the solution.

With the aim of making the stability analysis more quantitative, we introduce the following definition.

**Definition 4.1.** For problem (1.3) we define the relative condition number to be

$$K(d) = \sup\left\{\frac{\|\delta x\|/\|x\|}{\|\delta d\|/\|d\|}, \delta d \neq 0, d + \delta d \in D\right\} \quad (1.6)$$

Whenever  $d = 0$  or  $x = 0$ , it is necessary to introduce the absolute condition number, given by

$$K_{abs}(d) = \sup\left\{\frac{\|\delta x\|}{\|\delta d\|}, \delta d \neq 0, d + \delta d \in D\right\} \quad (1.7)$$

Problem (1.3) is called ill-conditioned if  $K(d)$  is “big” for any admissible datum  $d$ .

The property of a problem of being well-conditioned is independent of the numerical method that is being used to solve it. In fact, it is possible to generate stable as well as unstable numerical schemes for solving well-conditioned problems. The concept of stability for an algorithm or for a numerical method is analogous to that used for problem (1.3) and will be made precise in the next section.

**Remark 4.2.** (Ill-posed problems) Even in the case in which the condition number does not exist (formally, it is infinite), it is not necessarily true that the problem is ill-posed. In fact there exist well posed problems for which the condition number is infinite, but such that they can be reformulated in equivalent problems with a finite condition number.

If problem (1.3) admits a unique solution, then there necessarily exists a mapping  $G$ , that we call resolvent, between the sets of the data and of the solutions, such that

$$x = G(d), \text{ that is } F(G(d), d) = 0 \quad (1.8)$$

According to this definition, (1.4) yields  $x + \delta x = G(d + \delta d)$ . Assuming that  $G$  is differentiable in  $d$  and denoting formally by  $G'(d)$  its derivative with respect to  $d$  (if  $G : \mathbb{R}^n \rightarrow \mathbb{R}^m$ ,  $G'(d)$  will be the Jacobian matrix of  $G$  evaluated at the vector  $d$ ), a Taylor’s expansion of  $G$  truncated at first order ensures that

$$G(d + \delta d) - G(d) = G'(d)\delta d + o(\|\delta d\|) \quad \text{for } \delta d \rightarrow 0$$

where  $\|\cdot\|$  is a suitable vector norm and  $o(\cdot)$  is the classical infinitesimal symbol denoting an infinitesimal term of higher order with respect to its argument. Neglecting the infinitesimal of higher order with respect to  $\|\delta d\|$ , from (1.6) and (1.7) we respectively deduce that

$$K(d) \approx \|G'(d)\| \frac{\|d\|}{\|G(d)\|}, \quad K_{abs}(d) \approx \|G'(d)\| \quad (1.9)$$

where the symbol  $\|\cdot\|$ , when applied to a matrix, denotes the induced matrix norm (1.10) associated with the vector norm introduced above. The estimates in (1.9) are of great practical usefulness in the analysis of problems in the form (1.8).

**Theorem 4.1.** Let  $\|\cdot\|$  be a vector norm. The function

$$\|A\| = \sup_{x \neq 0} \frac{\|Ax\|}{\|x\|} \quad (1.10)$$

is a matrix norm called induced matrix norm or natural matrix norm.

In view of (1.9), the quantity  $\|G'(d)\|$  is an approximation of  $K_{abs}(d)$  and is sometimes called first order absolute condition number. This latter represents the limit of the Lipschitz constant of  $G$  as the perturbation on the data tends to zero.

Such a number does not always provide a sound estimate of the condition number  $K_{abs}(d)$ . This happens, for instance, when  $G'$  vanishes at a point whilst  $G$  is nonnull in a neighborhood of the same point. For example, take  $x = G(d) = \cos(d) - 1$  for  $d \in (-\pi/2, \pi/2)$ , we have  $G'(0) = 0$  while  $K_{abs}(0) = 2/\pi$ .

## 5. Stability of Numerical Methods

We shall henceforth suppose the problem (1.3) to be well posed. A numerical method for the approximate solution of (1.3) will consist, in general, of a sequence of approximate problems

$$F_n(x_n, d_n) = 0 \quad n \geq \quad (1.11)$$

depending on a certain parameter  $n$  (to be defined case by case). The understood expectation is that  $x_n \rightarrow x$  as  $n \rightarrow \infty$ , i.e. that the numerical solution converges to the exact solution. For that, it is necessary that  $d_n \rightarrow d$  and that  $F_n$  approximates  $F$ , as  $n \rightarrow \infty$ . Precisely, if the datum  $d$  of problem (1.3) is admissible for  $F_n$ , we say that (1.11) is consistent if

$$F_n(x, d) = F_n(x, d) - F(x, d) \rightarrow 0 \text{ for } n \rightarrow \infty \quad (1.12)$$

where  $x$  is the solution to problem (1.3) corresponding to the datum  $d$ .

A method is said to be strongly consistent if  $F_n(x, d) = 0$  for any value of  $n$  and not only for  $n \rightarrow \infty$ .

In some cases problem (1.11) could take the following form

$$F_n(x_n, x_{n-1}, \dots, x_{n-q}, d_n) = 0 \quad n \geq q \quad (1.13)$$

where  $x_0, x_1, \dots, x_{q-1}$  are given. In such case, the property of strong consistency becomes  $F_n(x, x, \dots, x, d) = 0$  for all  $n \geq q$ .

Recalling what has been previously state about problem (1.3), in order for the numerical method to be well posed (or stable) we require that for any fixed  $n$ , there exists a unique solution  $x_n$  corresponding to the datum  $d_n$ , that  $x_n$  depends continuously on the data. More precisely, let  $d_n$  be an arbitrary element of  $D_n$ , wehre  $D_n$  is the set of all admissible data for (1.11). Let  $\delta d_n$  be a perturbation admissible in the sense that  $d_n + \delta d_n \in D_n$ , and let  $\delta x_n$  denote the corresponding perturbation on the solution, that is

$$F_n(x_n + \delta x_n, d_n + \delta d_n) = 0$$

Then we require that

$$\begin{aligned} \exists \eta_0 = \eta_0(d_n) < 0, \exists K_0 = K_0(d_n) \text{ such that} \\ \text{if } \|\delta d_n\| \leq \eta_0 \text{ then } \|\delta x_n\| \leq K_0 \|\delta d_n\| \end{aligned} \quad (1.14)$$

As done in (1.6), we introduce for each problem in the sequence (1.11) the quantities

$$\begin{aligned} K_n(d_n) &= \sup \left\{ \frac{\|\delta x_n\| / \|x_n\|}{\|\delta d_n\| / \|d_n\|}, \delta d_n \neq 0, d_n + \delta d_n \in D_n \right\}, \\ K_{abs,n}(d_n) &= \sup \left\{ \frac{\|\delta x_n\|}{\|\delta d_n\|}, \delta d_n \neq 0, d_n + \delta d_n \in D_n \right\} \end{aligned} \quad (1.15)$$

The numerical method is said to be well condition if  $K_n(d_n)$  is “small” for any admissible datum  $d_n$ , ill conditioned otherwise. As in (1.8), let us consider the case where, for each  $n$ , the functional relation (1.11) defines a mapping  $G_n$  between the sets of the numerical data and the solutions

$$x_n = G_n(d_n), \text{ that is } F_n(G_n(d_n), d_n) = 0 \quad (1.16)$$

Assuing that  $G_n$  is differentiable, we can obtain from (1.15)

$$K_n(d_n) \approx \|G'_n(d_n)\| \frac{\|d_n\|}{\|G_n(d_n)\|}, \quad K_{abs,n} \approx \|G'_n(d_n)\| \quad (1.17)$$

We observe that, in the case where the sets of admissible data in problems (1.3) and (1.11) coincide, we can use in (1.14) and (1.15) the quantity  $d$  instead of  $d_n$ . In such case, we can define the relative and absolute asymptotic condition number corresponding to the datum  $d$  as follows

$$K^{num}(d) = \lim_{k \rightarrow \infty} \sup_{n \geq k} K_n(d)$$

$$K_{abs}^{num}(d) = \lim_{k \rightarrow \infty} \sup_{n \geq k} K_{abs,n}(d)$$

The final foal of numerical approximation is, of course, to build, through numerical problems of type (1.11), solutions  $x_n$  that “get closer” to the solution of problem (1.3) as much as  $n$  gets larger. This concept is made precise in the next definition.

**Definition 5.1.** The numerical method (1.11) is convergent iff

$$\begin{aligned} \forall \epsilon > 0 \exists n_0 = n_0(\epsilon), \exists \delta = \delta(n_0, \epsilon) > 0 \text{ such that} \\ \forall n > n_0(\epsilon), \forall \delta d_n : \|\delta d_n\| \leq \delta \rightarrow \|x(d) - x_n(d + \delta d_n)\| \leq \epsilon \end{aligned} \quad (1.18)$$

where  $d$  is na admissible datum for the problem (1.3),  $x(d)$  is the corresponding solution and  $x_n(d + \delta d_n)$  is the solution of the numerical problem (1.11) with datum  $d + \delta d_n$ .

To verify the implication (1.18) it suffices to check that under the same assumptions

$$\|x(d + \delta d_n) - x_n(d + \delta d_n)\| \leq \frac{\epsilon}{2} \quad (1.19)$$

Indeed, thanks to (1.5) we have

$$\begin{aligned} \|x(d) - x_n(d + \delta d_n)\| &\leq \|x(d) - x(d + \delta d_n)\| \\ &\quad + \|x(d + \delta d_n) - x_n(d + \delta d_n)\| \leq K_0 \|\delta d_n\| + \frac{\epsilon}{2} \end{aligned}$$

Choosing  $\delta = \min\{\eta_0, \epsilon/(2K_0)\}$  one obtains (1.18).

Measures of the convergence of  $x_n$  to  $x$  are given by the absolute error or the relative error, respectively defined as

$$E(x_n) = |x - x_n|, \quad E_{rel}(x_n) = \frac{|x - x_n|}{|x|} \quad (\text{if } x \neq 0) \quad (1.20)$$

In the cases where  $x$  and  $x_n$  are matrix or vector quantities, in addition to the definitions in (1.20) it is sometimes useful to introduce the relative error by component defined as

$$E_{rel}^c(x_n) = \max_{i,j} \frac{|x - x_n|_{ij}}{|x_{ij}|} \quad (1.21)$$

## 5.1. Relations between Stability and Convergence

The concepts of stability and convergence are strongly connected.

First of all, if problem (1.3) is well posed, a necessary condition in order for the numerical problem (1.11) to be convergent is that it is stable.



Let us thus assume that the method is convergent, that is, (1.18) holds for an arbitrary  $\epsilon > 0$ . We have

$$\begin{aligned} \|\delta x_n\| &= \|x_n(d + \delta d_n) - x_n(d)\| \leq \|x_n(d) - x(d)\| \\ &+ \|x(d) - x(d + \delta d_n)\| + \|x(d + \delta d_n) - x_n(d + \delta d_n)\| \\ &\leq K(\delta(n_0, \epsilon), d) \|\delta d_n\| + \epsilon \quad (1.22) \end{aligned}$$

having used (1.5) and (1.19) twice. Choosing now  $\delta d_n$ , suchh that  $\|\delta d_n\| \leq \eta_0$ , we deduce that  $\|\delta d_n\| \leq \eta_0$ , we deduce that  $\|\delta x_n\|/\|\delta d_n\|$  can be bounded by  $K_0 = K(\delta(n_0, \epsilon), d) + 1$ , provided that  $\epsilon \leq \|\delta d_n\|$ , so that the method is stable. Thus, we are interested in stable numerical methods since only these can be convergent.

The stability of a numerical method becomes a sufficient condition for the numerical problem (1.11) to converge if this latter is also consistent with problem (1.3). Indeed, under these assumptions we have

$$\begin{aligned} \|x(d + \delta d_n) - x_n(d + \delta d_n)\| &\leq \|x(d + \delta d_n) - x(d)\| \\ &+ \|x(d) - x_n(d)\| + \|x_n(d) - x_n(d + \delta d_n)\| \end{aligned}$$

Thanks to (1.5), the first term at right-hand side can be bounded by  $\|\delta d_n\|$ . A similar bound holds for the third term, due to the stability property (1.14). Finally, concerning the remaining term, if  $F_n$  is differentiable with respect to the variable  $x$ , an expansion in a Taylor series gives

$$F_n(x(d), d) - F_n(x_n(d), d) = \frac{\partial F_n}{\partial x} \Big|_{(x,d)} (x(d) - x_n(d))$$

for a suitable  $x$  “between”  $x(d)$  and  $x_n(d)$ . Assuming also that  $\partial f_n/\partial x$  is invertible, we get

$$x(d) - x_n(d) = \left( \frac{\partial F_n}{\partial x} \Big|_{(x,d)} \right)^{-1} [F_n(x(d), d) - F_n(x_n(d), d)] \quad (1.23)$$

On the other hand, replacing  $F_n(x_n(d), d)$  with  $F_n(x(d), d)$  and passing to the norms, we find

$$\|x(d) - x_n(d)\| \leq \left\| \left( \frac{\partial F_n}{\partial x} \Big|_{(x,d)} \right)^{-1} \right\| \|F_n(x(d), d) - F(x(d), d)\|$$

Thanks to (1.12) we can thus conclude that  $\|x(d) - x_n(d)\| \rightarrow 0$  for  $n \rightarrow \infty$ . The result that has just been proved, although stated in qualitative terms, is a milestone in numerical analysis, known as equivalence theorem (or Lax-Richtmyer theorem): “for a consistent numerical method, stability is equivalent to convergence”.

## 6. Sources of Error in Computational Models

Whenever the numerical problem (1.11) is an approximation to the mathematical problem (1.3) and

this latter is in turn a model of a physical problem, we shall say that (1.11) is a computational model for PP.

In this process the global error, denoted by  $e$ , is expressed by the difference between the actually computed solution,  $\hat{x}_n$ , and the physical solution,  $x_{ph}$ , of which  $x$  provides a model. The global error  $e$  of the mathematical model, given by  $x - x_{ph}$ , and the error  $e_c$  of the computational model,  $\hat{x}_n - x$ , that is  $e = e_m + e_c$ .

The error  $e_m$  will in turn take into account the error of the mathematical model in strict sense and the error on the data. In the same way,  $e_c$  turns out to be the combination of the numerical discretization error  $e_n = x_n - x$ , the error  $e_a$  introduced by the numerical algorithm and the roundoff error introduced by the computer during the actual solution of problem (1.11).

In general, we can thus outline the following sources of error:

1. error due to the model, that can be controlled by a proper choice of the mathematical model;
2. errors in the data, that can be reduced by enhancing the accuracy in the measurement of the data themselves;
3. truncation error, arising from having replaced in the numerical model limits by operations that involve a finite number of steps;
4. rounding errors.

The error at the items 3. and 4. give rise to the computational error. A numerical method will thus be convergent if this error can be made arbitrarily small by increasing the computational effort. Of course, convergence is the primary, albeit not unique, goal of a numerical method, the others being accuracy, reliability and efficiency.

Accuracy means that the errors are small with respect to a fixed tolerance. It is usually quantified by the order of infinitesimal of the error  $e_n$  with respect to the discretization characteristic parameter. By the way, we notice that machine precision does not limit, on theoretical grounds, the accuracy.

Reliability means it is likely that the global error can be guaranteed to be below a certain tolerance. Of course, a numerical model can be considered to be reliable only if suitably tested, that is, successfully applied to several test cases.

Efficiency means that the computational complexity that is needed to control the error is as small as possible.

By algorithm we mean a directive that indicates, through elementary operations, all the passages that are needed to solve a specific problem. An algorithm can in turn contain sub-algorithms and must have the feature of terminating after a finite number of

elementary operations. As a consequence, the executor of the algorithm must find within the algorithm itself all the instructions to completely solve the problem at hand.

Finally, the complexity of an algorithm is a measure of its executing time. Calculating the complexity of an algorithm is therefore a part of the analysis of the efficiency of a numerical method. Since several algorithms, with different complexities, can be employed to solve the same problem  $P$ , it is useful to introduce the concept of complexity of a problem, this latter meaning the complexity of the algorithm that has a minimum complexity among those solving  $P$ . The complexity of a problem is typically measured by a parameter directly associated with  $P$ .

## 7. Machine Representation of Numbers

### 8. Ejercicios

(1) Se considera la siguiente sucesión definida por recursión

$$x_0 = 1 \quad x_1 = \frac{1}{5} \quad x_{n+1} = \frac{36}{5}x_n - \frac{7}{5}x_{n-1}$$

Esta sucesión tiene como solución  $x_n = \frac{1}{5^n}$ . Utilizar Matlab para calcular  $\frac{1}{5^n}$  utilizando la sucesión recursiva del principio, para  $n \leq 32$ . Hacer el estudio del error absoluto y relativo.

(2) Repetir el ejercicio anterior tomando  $x_0 = 2$  y  $x_1 = \frac{36}{5}$ , teniendo en cuenta que la solución es ahora  $x_n = \frac{1}{5^n} + 7^n$ .

(3) Compara los resultados de los ejercicios 1 y 2 y dar una explicación formal de lo que está sucediendo.

# Bibliografía

- [1] A. Quarteroni; R. Sacco; F. Sale. *Numerical Mathematics*. Springer, 2007.