

# گزارش همایش ماهانهٔ انجمن ریاضی ایران

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#### PREFACE

The aim of this foreword is to give a short presentation of the Monthly Colloquium of the Iranian Mathematical Society (IMS). It will be best understood through a brief history of the IMS activities and the role of conferences therein.

The IMS originated its effective life from a meeting in 1970 and has organized several conferences thereafter. The Annual Mathematics Conference in Iran (AMCI) has undoubtedly the largest evaluating number among them. This is reflected from both scientific and social aspects. Its scientific importance is due to the general talks given by the main invited speakers, the number and the length of its refereed papers, and the fact that the proceedings of the past 34 AMCI have been almost regularly published and are now reviewed in both Mathematical Reviews and MathSci. Its social importance is credited to the number of participants, the enthusiastic panels concerning some local or universal problems in mathematical education or research, the distribution of prizes to winners in its opening ceremonies, the traditional annual session of the General Assembly of the IMS held simultaneously, the announcement of some decisions made by the General Assembly in the closing ceremonies, cultural activities containing visits of a few tourist sites in the country offered to the foreign speakers, folkloric or classical concerts, etc.

These annual conferences remain one of the main concerns of the IMS. However, from the early days of the IMS's life, other needs have motivated the organization of a lot of different seminars and colloquia as well as the publication of periodical and non periodical issues.

During the first decade (up to 1979), a Weekly Seminar has also been

held in Tehran at different universities following the speaker's affiliation. From the end of that period, many specialized seminars have been organized by the IMS and held at different universities not necessarily in Tehran. For instance, the first Algebra Seminar was held in Ahwaz (December 31st 1977 - January 1st 1978). Due to a lot of circumstances, the proceedings of this seminar has not been published. There are lots of other annual or biannual well-established seminars in algebra, analysis, geometry and topology, differential equations and dynamical systems, etc. They are mostly publishing their proceedings. A detailed analysis of these conferences and seminars requires a serious study, which is quite interesting, but goes beyond the scope of this preface.

Let me now return back to the philosophy of the Monthly Colloquium as I feel its spirit. The general topics addressed in AMCI presented by main invited speakers to a large audience are of great importance. However, their number are limited by the restrictions naturally imposed by such a program with a big variety in a short period of time (4 days only). The main speeches presented at specialized seminars have also the same shortage of time (2 days only) and, in plus, their participants are almost exclusively scholars involved in the field. By the way, the participants of the AMCI and the specialized seminars are practically the only persons who are receiving the proceedings. But a general talk or a good survey article is of interest to a larger public. Anyone of our graduate students and us needs to obtain information about what is done by others even outside his specialties. So, the IMS has decided to organize a Monthly Colloquium presenting mainly survey articles by research leaders in the country or from abroad. The articles presented for the moment at Tehran are addressed to a reasonable audience, around one hundred graduate students and professors. Ideally, a first draft of

the papers will be distributed to the audience the day of the colloquium. Moreover, they will potentially be recorded in videotapes, for the use of universities and research centers outside the capital, etc. Finally, they will be published, in a final form, annually or biannually, when a lot of papers are ready for printing.

During the period 2000-2003, the organizing committee of the Monthly Colloquium that I had the honor to manage has published the first volume containing 8 articles written in Persian, and is publishing now the second volume containing 11 articles in English. These 2 volumes are available from the IMS head office and can be obtained by the members of the society as well as the interested general public.

The third and next volumes will be published by our successors. I am confident that they will continue to fulfill the requirements of the mathematical community more perfectly. The monthly colloquium of the Iranian Mathematical Society is an important step in the propagation of mathematics in our country and it is worth all the efforts done for it.

For the final word, I present my deepest gratitude to the authors who have accepted our invitation and submitted their priceless manuscripts for publication in volume one and this volume. Personally, I have learned a lot from them. The publication of this issue would not have been possible without the endeavors of the secretaries of the IMS whom I thank hereby. Specifically, I would like to thank Mrs. F. Samadian who adapted the format of the Farhang-va-Andishe-ye-Riazi for the first volume and that of the Bulletin of the Iranian Mathematical Society for the present volume.

Arsalan Chademan Tehran, May 10, 2004

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Proceedings of the Monthly Colloquium of the Iranian Mathematical Society Vol. 2, pp 1-20 (2004)

# Numerical Solution of Integral Equations

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Abstract: Mathematical models of a wide class of problems are integral equations. Most of integral equations have no analytic solution. In this talk I introduce different kinds of integral and integro-differential equations, their origins and applications. I also explain, in some detail, the numerical methods for solving integral equations.

# 1. Introduction

In some sense integral equations must be felt to be either more advanced or of less practical interest than differential equations. We turn more

MSC (2000): 45A05, 45B05, 45D05, 45E05, 45J05

Keywords: Integral equations, Fredholm integral equations, Volterra integral equations, Integro-differential equations, Nystrom method, Galerkin method, Expansion method.

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readily to a differential formulation of a problem than to an integral formalism. Yet the theory of linear nonsingular integral equations is at least as well developed as that of differential equations and it is in many respects rather simpler. The corresponding operators are bounded rather than unbounded, leading to a very straightforward existence theory (the Fredholm theory). There is a much tighter link between the theory and practice of integral equations than is the case for with differential equations. Most of the convergence proofs are constructive in nature and all or nearly all of the constructions have been used as the basis of algorithms for the numerical solution of the underlying equations (although not always with any great success).

For those who require detailed mathematical theory of Integral Equations (I. E.), a number of books are available:

- 1. S. Smithies, Integral Equations, Combridge University Press (1958).
- 2. R. P. Kanwal, Linear Integral Equations, Academic Press (1971).
- 3. C. D. Green, Integral Equation Methods, Nelson (1969).
- 4. J. A. Cochran, The Analysis of Linear Integral Equations, McGraw-Hill (1972).

The books which you may find useful in connection with the numerical solution of integral equations are listed below:

- 1. S. G. Mikhlin and J. L. Smolitsky, Approximate Methods for the Solution of Differential and Integral Equations, Elsevier (1967).
- L. M. Delves and J. Walsh (eds.), Numerical Solution of Integral Equations, Oxford University Press (1974).
- C. T. H. Baker, The Numerical Treatment of Integral Equations, Oxford University Press (1977).
- 4. C. T. H. Baker and G. F. Miller (eds.), Treatment of Integral Equations by Numerical Methods, Academic Press (1982).

 L. M. Delves and J. L. Mohamed, Computational Methods for Integral Equations, Cambridge University Press (1985).

## 2. Classification

 Linear Fredholm Integral Equations of the First and Second Kind, respectively

$$y(s) = \lambda \int_a^b K(s,t)x(t)dt,$$
  
$$x(s) = y(s) + \lambda \int_a^b K(s,t)x(t)dt,$$

where x(t) is the unknown function and K(s,t) is the kernel of the equation.

ii) Linear Volterra Integral Equations of the First and Second Kind

$$y(s) = \lambda \int_a^s K(s,t)x(t)dt,$$
  
$$x(s) = y(s) + \lambda \int_a^s K(s,t)x(t)dt.$$

iii) Homogeneous equation, eigenvalue equation, or a Fredholm equation of the third kind

$$x(s) = \lambda \int_a^b K(s, t) x(t) dt$$

those values of  $\lambda$  for which nontrivial solutions exist are called characteristic values of the equation.

iv) Nonlinear Integral Equations

In (i) to (iii) if you replace K(s,t) with K(s,t,x(s),x(t)) you obtain a nonlinear integral equation.

- v) Integral equation involving unknown functions  $x(s_1, s_2, ..., s_m)$  also occur regularly in practice but lie outside the scope of this talk!
- vi) Integro-differential equations

$$Q(s)x'(s) + R(s)x(s) = y(s) + \lambda \int_a^b K(s,t)x(t)dt, \quad x(a) = x_0.$$

# 3. Origin of Integral Equations

Consider the following initial value problem

$$\frac{dx(s)}{ds} = f(s,x), \quad x(0) = x_0,$$

under suitable continuity conditions on f(s, x), we have:

$$x(s) = x_0 + \int_0^s f(t, x(t))dt,$$

which is an integral equation for x(s).

ii) Consider the second order differential equation

$$\frac{d^2x}{ds^2} = f(s, x)$$

$$x(0) = x_0, \frac{d(x(0))}{ds} = x_1,$$

this equation converts to

$$x(s) = x_0 + x_1 s + \int_0^s (s - t) f(t, x(t)) dt$$

which is an integral equation for x(s).

iii) Two-point boundary value problems may be converted to an integral equation. Consider the problem

$$\begin{cases} \frac{d^2x}{ds^2} = f(s, x), \\ x(0) = \alpha, x(l) = \beta, \end{cases}$$

this problem converts to

$$x(s) = y(s) - \int_0^t K(s,t)f(t,x(t))dt,$$

where

$$y(s) = \alpha + \frac{\beta - \alpha}{l}s,$$

$$K(s,t) = \begin{cases} \frac{t(l-s)}{l}, & 0 \le t \le s, \\ \frac{s(l-t)}{l}, & s \le t \le l. \end{cases}$$

# 4. Applications

We take a random selection of research papers in a variety of fields.

 Currents in a superconducting strip (Rhoderick and Wilson, 1962 [18])

$$y(s) = \frac{1}{\pi} \int_0^1 \frac{t-s}{(t-s)^2 + a^2} x(t) dt.$$

ii) Flow round a hydrofil (Kershaw, 1974 [15])

$$x(s) = y(s) + \lambda \int_C K(s,t)x(t)dt,$$

subject to:  $\int_C x(t)dt = 0$ , C a closed contour.

iii) Population competition (Dounham and Shah, 1976 [12])

$$\mathbf{x}(s) = \mathbf{y}(s) \int_a^b \mathbf{k}(s, t, \mathbf{x}(s, t)) \mathbf{x}(t) dt, \mathbf{x}^T = (x_1, x_2, x_3).$$

iv) Quantum scattering: closed coupled calculations (Horn and Frazer, 1975 [14])

$$x(s) = y(s) - \lambda \int_0^\infty K_1(s,z) \int_0^\infty K_2(z,t)x(t)dtdz.$$

# 5. Existence of Solution to Integral Equations

Smithies [19], defined the concept of relatively uniformly convergence of sequences of functions and kernels.

**Definition.** Suppose  $\{f_n(s)\}$  be a sequence of  $\mathcal{L}^2$  functions. We say that this sequence converges relatively uniformly (c.r.u.) to f(s) if there exists a non-negative  $\mathcal{L}^2$  function P(s) such that

$$\forall \epsilon \exists N(\epsilon) \forall n \forall s \ (n \ge N(\epsilon) \Longrightarrow |f_n(s) - f(s)| \le \epsilon P(s)).$$

It can be shown that if  $\{f_n(s)\}$  (c.r.u.) to f(s) then it converges pointwise, but not necessarily uniformly. Absolutely relatively uniformly convergence of sequences and series of functions and kernels can be defined similarly.

It is proved that if for a fixed  $\lambda$  there exists kernel  $H_{\lambda}(s,t) \in \mathcal{L}^{2}[a,b] \times [a,b]$  such that

$$K(s,t) - H_{\lambda}(s,t) = \lambda H_k K = \lambda K H_{\lambda},$$

then  $x = y + \lambda K H_{\lambda}$  is a solution of the equation

$$x = y + \lambda K x.$$

It is also shown that if  $||\lambda K|| < 1$  then the series

$$K(s,t) + \lambda K^2(s,t) + \lambda^2 K^3(s,t) + \dots$$

known as Neumann series, is absolutely relatively uniformly convergent to  $H_{\lambda}(s,t)$ . (For a detailed discussion of the above concepts see Smithies [19]).

Values of  $\lambda$  for which  $H_{\lambda}(s,t)$  exist are called regular values of the kernel K. It is shown that the set of regular values of K is open.

# 6. Numerical Solution of Fredholm Integral Equations

### 6.1 Iterative method and the Neumann series

The Fredholm I. E. of the second kind

$$x(s) = y(s) + \lambda \int_{a}^{b} K(s, t)x(t)dt$$
 (1)

can be written in operator form as

$$x = y + \lambda Kx$$
.

The Neumann series is

$$x = \sum_{i=0}^{\infty} \lambda^i K^i y, \tag{2}$$

and in truncated form

$$x \simeq x_n = \sum_{i=0}^n \lambda^i K^i y.$$

If the series (2) converges, then  $\lim_{n\to\infty} ||x-x_n|| = 0$ . The approximate solution  $x_n$  is most easily produced iteratively via the obvious recurrence relation:

$$x_{n+1} = y + \lambda K x_n$$
$$x_0 = y,$$

provided that the function  $Kx_n = \int_a^b K(s,t)x_n(t)dt$  can be computed.

# 6.2 The Nyström (or quadrature) method

In equation (1) if you use a quadrature rule you find

$$\int_a^b K(s,t)x(t)dt \simeq \sum_{j=1}^n w_j K(s,t_j)x(t_j),$$

where  $w_j$  and  $t_j$  are weights and nodes of the quadrature rule. So (1) converts to (approximately)

$$x(s) = y(s) + \lambda \sum_{j=1}^{n} w_j K(s, t_j) x(t_j),$$
 (3)

if you put  $s = t_i$ , i = 1, ..., n, in (3) you get the following system of linear equations for the unknowns  $x(t_1), ..., x(t_n)$ ,

$$x(t_i) = y(t_i) + \lambda \sum_{j=1}^{n} w_j K(t_i, t_j) x(t_j), \quad i = 1, 2, \dots, n.$$
 (4)

If you define  $\mathbf{x}^i = (x(t_1), \dots, x(t_n)), \mathbf{y}^i = (y(t_1), \dots, y(t_n))$  and  $\mathbf{k}_{ij} = w_j K(t_i, t_j)$  the system (4) becomes

$$(I - \lambda \mathbf{k})\mathbf{x} = \mathbf{y},\tag{5}$$

this equation has a solution if  $I - \lambda k$  is not singular, and this may happen even if  $||\lambda k|| > 1$ . In general the matrix  $I - \lambda k$  is well-conditioned and the system (5) can be solved easily by direct methods.

# 6.3 Expansion method for Fredholm Equations

# **6.3.1** Methods based on an expansion of the solution of Fredholm I. E. of the second kind

Suppose that the sequence of functions  $\{h_n(s)\}$  is orthogonal and complete in  $\mathcal{L}^2(a,b)$  and

$$x(s) = \sum_{j=1}^{\infty} b_j h_j(s),$$

then for sufficiently large N, we may approximate x as closely as we please by  $x_N$ :

$$x(s) \simeq x_N(s) = \sum_{j=1}^N a_j h_j(s).$$

An expansion method is then an algorithm for determining the  $a_i$  for either an arbitrary or a specified choice of the sequence  $\{h_n(s)\}$ . There are many possible algorithms and we consider only the most common.

1. Residual minimisation methods Assuming  $L = I - \lambda K$ , the equation (1) can be written as

$$Lx = y$$
.

Now if  $\epsilon_N = x - x_N$ ,  $r_N = y - Lx_N$  then  $r_N = L\epsilon_N$  and computing  $r_N$  requires no knowledge of x. We now choose the vector  $\mathbf{a} = (a_1, \ldots, a_n)^t$  from the minimisation criterion:  $\min_{\mathbf{a}} ||r_N||$ .

It can be shown that

$$\frac{||r_N||}{1+||K||} \le ||\epsilon_N|| \le \frac{||r_N||}{1-||K||},$$
 provided that  $||K|| < 1.$ 

There are many norms avialable, the choice of norm is influenced by the analytical properties of the kernel K.

Chebyshev norm. For continuous kernels a possible choice is the Chebyshev  $(L_{\infty})$  norm:

$$||x|| = \max_{a \le x \le b} |x(s)|, \quad ||K|| = \max_{a \le s \le b} \int_a^b |K(s,t)| dt.$$

We seek to compute

$$||r_N|| = \min_{a} \max_{a \le s \le b} |y(s) - x_N(s) + \int_a^b K(s,t)x_N(t)dt|,$$

with  $x_N = \sum_{i=1}^N a_i h_i(s)$ , setting  $k_i(s) = \int_a^b K(s,t) h_i(t) dt$ ,

$$||r_N|| = \min_{s} \max_{s} |y(s) - \sum_{i=1}^n a_i (h_i(s) - k_i(s))|,$$

usually  $k_i(s)$  can be estimated by numerical quadrature (for fixed s) and to search, not over [a, b], but over a discrete point set  $\{\xi_i, i = 1, \ldots, q\}$ , i.e.

$$||r_N|| \simeq \min_{\mathbf{a}} \max_{k=1}^q |y(\xi_k) - \sum_{i=1}^n (h_i(\xi_k) - k_i(\xi_k))|.$$

## $L_2$ norm: Least squares approximation

Alternatively, we may choose to minimise  $||y - Lx||_2$ :

$$\min_{\mathbf{a}} \int_a^b \left| \left[ x_N(s) - y(s) - \int_a^b K(s,t) x_N(t) dt \right] \right|^2 ds,$$

whence on inserting  $x_N$  we find the system

$$L_{ls}a = y_{ls}$$

where

$$(\mathbf{L}_{ls})_{ij} = \int_a^b (Lh_i)^*(s)h_j(s)ds, \quad i, j = 1, ..., N$$

$$(\mathbf{y}_{ls})_i = \int_a^b y(s)(Lh_i)^*(s)ds, \quad i = 1, ..., N$$

$$Lh_i(s) = h_i(s) - \int_a^b K(s,t)h_i(t)dt.$$

In this norm we have to evaluate  $N^2$  apparently triple integrals of the form

$$\int_a^b \left[ \int_a^b K^*(s,t) h_i^*(t) dt \right] \left[ \int_a^b K(s,t') h_j(t) dt' \right] ds.$$

2. Galerkin method. Our aim is to make r(s) zero. Now if the set  $\{h_i\}$  is complete and orthonormal in  $\mathcal{L}^2(a,b)$ , the statement r(s)=0 is equivalent to

$$(r(s), h_i(s)) = 0, \quad i = 1, 2, \dots$$

Now with only N parameters  $a_i$ , i = 1, ..., N, at our disposal the best we can do is to make the residual  $r_N(s)$  orthogonal to the first N functions  $h_1(s), ..., h_N(s)$ , i.e.

$$\int_a^b h_i^*(s)[y(s) - (Lx_N)(s)]ds = 0, \quad i = 1, \dots, N.$$

this leads to the system  $L_G a = y_G$  where now

$$(\mathbf{L}_G)_{ij} = \int_a^b h_i^*(s)(Lh_j(s))ds, \quad i, j = 1, \dots, N$$
$$(\mathbf{y}_G)_i = \int_a^b h_i^*(s)y(s)ds, \quad i = 1, \dots, N.$$

3. The Fast Galerkin algorithm Delves [10,11] and his Ph.D. students used Chebyshev sequence of polynomials  $\{T_i(s)\}_{i=0}$  and set  $x(s) \simeq \sum_{i=0}^{N} a_i T_i(s)$  and for

$$x(s) = y(s) + \int_{-1}^{1} K(s,t)x(t)dt$$

obtained

$$(\bar{\mathbf{D}} - \bar{\mathbf{B}})\mathbf{a} = \bar{\mathbf{y}}$$

where

$$\bar{B}_{ij} = \int_{-1}^{1} \frac{T_i(s)}{(1-s^2)^{\frac{1}{2}}} \left( \int_{-1}^{1} K(s,t) T_j(t) dt \right) ds, \quad i, j = 0, 1, \dots, N$$

$$\bar{y}_i = \int_{-1}^{1} \frac{T_i(s)}{(1-s^2)^{\frac{1}{2}}} y(s) ds, \quad i = 0, 1, \dots, N.$$

They used Fast Fourier Transformation (FFT) and evaluated approximate values  $B_{ij}$  and  $y_i$  for  $\bar{B}_{ij}$  and  $\bar{y}_i$  in  $\mathcal{O}(N^2 \log N)$  operations. Another advantage of this method is that

$$|\bar{y}_i| \le c\hat{i}^{-p}, \hat{i} = \max\{1, i\}, \quad i = 0, 1, \dots$$

where p depends on the smoothness of the function y(s). Similarly they showed that  $|\bar{B}_{ij}| \leq c''\hat{i}^{-(r+1)}\hat{j}^{-2}$  if  $|k_{ij}| \leq c'\hat{i}^{-(r+1)}\hat{j}^{-(q+1)}$  where  $K(s,t) = \sum_{i=0}^{\infty} \sum_{j=0}^{\infty} k_{ij}T_i(s)T_j(t)$ , again r,q depend on the countinuity properties of K(s,t) as functions of s and t, respectively. They introduced asymptotically lower diagonal matrix M of type A and degree p

$$\frac{|M_{ij}|}{|M_{ii}||M_{jj}|} \le ci^{-p}, \quad i > j, c > 0, p \ge 0,$$

and showed that the matrix  $I - \lambda \bar{B}$  is asymptotically lower diagonal of type A and degree r [13]. This was the basis of an iterative method for solving  $\mathbf{L}_G \mathbf{a} = \mathbf{y}_G$  with  $\mathcal{O}(N^2)$  operations. Some authors have used splines or wavelet functions to solve Fredholm I. Es. of the second kind (Razzaghi and Ordokhani [17]).

# 7. Fredholm Integral Equations of the First Kind

All of the numerical methods discussed for solving second kind equations apply formally also to first kind equation

$$y(s) = \int_a^b K(s, t)x(t)dt, \tag{6}$$

with the single exeption of the Neumann iterations.

Fredholm integral equations of the first kind are among ill-posed problems, i.e.

- i) They may have no solution;
- ii) They may have not a unique solution;
- iii) The solution does not depend on y(s) continuously, i.e. small changes on y(s) may cause dramatic changes on the solution x(s).

For example the equation  $\int_0^{2\pi} \cos s \sin tx(t)dt = e^s$  has no solution. If x is a solution of Kx = y and  $\phi \neq 0$  is such that  $K\phi = 0$  then  $x + \alpha\phi$  is also a solution of Kx = y for arbitrary  $\alpha \in \mathbb{C}$ .

# 7.1 Eigenfunction expansion

The first method we consider (Baker et al., 1964; Turchin, Kozlov and Malkevich, 1971; Hanson, 1971 [9]) is a direct extension of the numerical method for finite rank kernels.

**Theorem.** Let K be of finite rank n and have the representation:

$$K=\sum_{\nu=1}^n a_\nu\oplus b_\nu$$

then (6) has no solution unless  $y \in \text{span}\{a_1, \ldots, a_n\}$ , i.e.

$$\exists a_1,\ldots,a_n:y(s)=\sum_{\nu=1}^n\alpha_{\nu}a_{\nu}.$$

In this case (6) has the solution

$$x=\sum_{\nu=1}^n\beta_\nu a_\nu,$$

if and only if the following system of linear equations is nonsingular

$$\sum_{\nu=1}^{n} (b_{\mu}, a_{\nu}) \beta_{\nu} = a_{\mu}, \quad \mu = 1, \dots, n.$$

#### Hermitian kernels

If K is Hermitian, i.e.  $\overline{K(t,s)} = K^*(s,t) = K(s,t)$ , it has real eigenvalues  $\lambda_i$ , with corresponding eigenfunctions  $v_i$ , satisfying the equation  $\int_a^b K(s,t)v_i(t)dt = \lambda_i v_i(s)$ . If the  $\{v_i\}$  form a complete set, any function, and in particular the solution x and driving term y, have expansions of the form

$$x = \sum_{i=1}^{\infty} \beta_i v_i, \quad y = \sum_{i=1}^{\infty} \alpha_i v_i$$

giving  $Kx = \sum_{i=1}^{\infty} \beta_i K v_i = \sum_{i=1}^{\infty} \beta_i \lambda_i v_i = y = \sum_{i=1}^{\infty} \alpha_i v_i$ . So,  $\beta_i = \alpha_i/\lambda_i$  and  $x = \sum_{i=1}^{\infty} (v_i, y) \lambda_i^{-1} v_i$ .  $[v_i, i = 1, 2, \dots]$  are assumed to be orthonormal]. The eigenvalues  $\lambda_i$  tend to zero as i increases, giving the sum

$$\sum_{i=1}^{\infty} (v_i, y) \lambda_i^{-1} v_i, \tag{7}$$

a dangerous look. Now if  $\lim_{i\to\infty}(v_i,y)\lambda_i^{-1}\neq 0$  the sum (7) does not represent an  $\mathcal{L}^2$  function; that is the equation has no solution, a numerical test for consistency. However, even if  $(v_i,y)\lambda_i^{-1}\underset{i\to\infty}{\longrightarrow} 0$ , the series (7) is very unstable. Suppose we replace y by  $y+\epsilon v_k$ . Then the solution x changes to  $x+\epsilon\lambda_k^{-1}v_k$  and the "response ratio", defined as  $\|\delta x\|/\|\delta y\|$  is  $|\lambda_k^{-1}|$ . Since  $\lim_{k\to\infty}|\lambda_k|=0$ , this response ratio can be made arbitrary large, showing that the solution x does not depend continuously on y.

#### Non-Hermitian kernels

If K is not Hermitian  $K^*K$  and  $KK^*$  are Hermitian. Singular values  $\mu_i$  and singular functions  $\{u_i, v_i\}$  satisfy the relations

$$u_i = \mu_i K v_i, \quad v_i = \mu_i K^* u_i, \quad (u_i, u_j) = (v_i, v_j) = \delta_{ij}.$$

Further,  $\lim_{i\to\infty}\mu_i^{-1}=0$ .

Now suppose we make the expansions

$$x = \sum_{i=1}^{\infty} \beta_i v_i = \sum_{i=1}^{\infty} (v_i, x) v_i,$$
$$y = \sum_{i=1}^{\infty} \alpha_i u_i = \sum_{i=1}^{\infty} (u_i, y) u_i.$$

Then it follows that  $\beta_i = \mu_i(u_i, y)$  and

$$x = \sum_{i=1}^{\infty} (u_i, y) \mu_i v_i.$$

### The method of regularisation

In many practical applications the driving term y(s) represents some measured function, the integral operator K represents a model of the instrument used to do the measuring and x(s) represents the "true" measured quantity, y(s) being a "smeared" version of x(s) as seen through the instrument being used. Now y(s) will be known to some finite accuracy

 $\epsilon$ ; we should therefore only expect to find  $||Kx - y|| \leq \epsilon$ . Of all the functions x satisfying this relation, we seek that which is the smoothest, in the sence that for some linear operator L, ||Lx|| has the minimum value. This yields the constrained minimisation problem

minimise 
$$||Lx||$$
 subject to  $||Kx - y|| \le \epsilon$ . (8)

This problem can be solved directly, in any norm. It is however relatively complicated to do so. It is therefore usual to solve, not (8), but a related problem, which we develop as we note that the minimum value of ||Lx|| in (8) will decrease as  $\epsilon$  increases, that is as the constrains weaken. Therefore, at the minimum of (8), the constrains will be binding, that is  $||Kx - y|| = \epsilon$ . Now, if we solve the unconstrained problem (for fixed  $\alpha$ ).

$$minimise_x ||Kx - y||^2 + \alpha ||Lx||^2,$$
 (9)

we will find the minimum some value  $\eta$  for ||Kx - y||. As  $\alpha \longrightarrow 0$ ,  $\eta \longrightarrow 0$  provided that a solution of (6) exists, and for some value of  $\alpha, \eta = \epsilon$ . This makes plausible the following results, due to Tikhonov [20,21].

For some  $\alpha$  (which depends on  $\epsilon$ ) the solution of problem (9) is identical to that of problem (8). Problem (9) is easier to solve and it is referred to as the *regularised problem*. The choice of operator L must be made on qualitative grounds, the usual choices are: L = I or  $\frac{d}{as}$  or  $\frac{d^2}{ds^2}$ . When L = I it can be shown that the solution to (9) is the solution of the following Fredholm I. E. of the second kind,

$$\int_{a}^{b} \hat{K}(s,t)x(t)dt + \alpha x(s) = \hat{y}(s)$$
 (10)

where

$$\hat{K}(s,t) = \int_a^b K^*(\xi,s)K(\xi,t)d\xi, \quad \hat{y}(s) = \int_a^b K^*(t,s)y(t)dt.$$

Unfortunately the value of  $\alpha$  for which  $\eta = \epsilon$  is very small and the solution of (10) is very sensitive to  $\alpha$ . This makes this kind of regularisation very expensive and impractical.

Regularisation method was also used by Phillips [16] and Twomey [22], they used  $L = \frac{d^2}{ds^2}$ .

## The augmented Galerkin method

We now turn to the use of expansion methods for (6), and recall that (6) converts to

$$\bar{B}\mathbf{a} = \bar{\mathbf{y}} \tag{11}$$

where  $\bar{B}$  is very ill-conditioned, and even singular when K is of finite rank. Recall that we assumed

$$x(s) = \sum_{i=0}^{\infty} b_i h_i(s)$$

and for  $x(s) \in \mathcal{L}^2(a,b)$  we should have

$$\sum_{i=0}^{\infty} |b_i|^2 < \infty.$$

Hence we may assume that there exist constants  $c>0,\,r\geq\frac{1}{2}$  such that

$$|b_i| \leq C_b \hat{i}^{-r} = \delta_i, \quad i = 0, 1, \ldots$$

again r depends on the smoothness of the solution x(s).

So, instead of solving the ill-conditioned system (11) we solve the well-conditioned problem:

minimise 
$$\|\bar{B}\mathbf{a} - \bar{\mathbf{y}}\|$$
  
subject to:  $|a_i| \leq \delta_i, i = 0, 1, ..., N,$  (12)

where  $\delta_i = C_f \hat{i}^{-r}$ .

The problem is now to estimate  $C_f$ , r. There are a number of heuristic and practical ways to set  $C_f$ , r (see Babolian & Delves (1979) [8] and Babolian (1980) [7]). The solution of (12) is not sensitive to the values of  $C_f$ , r, but r should not be set very large unless x is very smooth.

Problem (12) has been solved in  $\|\cdot\|_{\infty}$  and  $\|\cdot\|_{2}$  (Belward (1982)). In 1997, Abbasbandy & Babolian [1,2] proposed some fully automatic algorithms to set  $C_f$  and r. Then some papers were published using automatic augmented Galerkin algorithms [3,4]. Some authors have used direct regularisation, by multiplying both sides of Kx = y by a function Q(s,t) (Maleknejad & Rostami) or used preconditioning to reduce the condition number of the resulting system of equations. But none of them have the general success of automatic augmented Galerkin algorithms.

# 8. Integro-differential Equations

Integro-differential equations are also considered using expansion methods with complete error analysis (see Babolian & Delves (1981) [5]).

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Proceedings of the Monthly Colloquium of the Iranian Mathematical Society Vol. 2, pp 21-38 (2004)

# Factorizations of Finite Groups

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Abstract: Let G be a finite group and A, B proper subgroups of G. If G = AB, then we say that G is a factorizable group and A, B are called factors of this factorization. In this case G is also called the product of two subgroups A and B. The problem of which finite groups are factorizable is still an open probelm. Certain conditions on the structure of G or the factors of the factorization yields some information about G. In this paper we will give a survey of the results obtained so far on the product of finite groups.

# 1. Introduction

Let G be a group and A, B subgroups of G. If  $G = AB = \{ab | a \in A, b \in B\}$ , then we say that G is a factorizable group and A, B are called

MSC (2000): 20D40

Keywords: Factorization, Simple group, Product of groups

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factors of the factorization. We also say that G is the product of its subgroups A and B. We always have G = AG which is called the trivial factorization of G. Therefore we call a factorization G = AB proper or non-trivial if both factors are proper subgroups of G. If  $G \cong A \times B$  is the external direct product of two non-trivial groups A and B, then  $G \cong \bar{A}\bar{B}$ , where  $\bar{A} \cong A$  and  $\bar{B} \cong B$  and so G has a proper factorization. Also factorizations of groups as product of 3 or more subgroups may be of interest, but it is not considered here. In the book [AFD], page 13, the authors ask the following question:

Question: Describe all groups that have a proper factorization.

The above question is an open problem in group theory and it seems to be a difficult problem in general. Not every group has a factorization, for example if G is an infinite group with all proper subgroups finite, then G is not factorizable. And if G is the cyclic group of prime order or one of the sporadic finite simple groups:

 $M_{22}, J_1, J_3, J_4, M^{\circ}L, L_y, O'N, Co_2, Co_3, Fi_{23}, Fi'_{24}, HN, TH, BM$  or M then G does not have a proper factorization. The other 11 sporadic groups have proper factorization, for example:

 $M_{11} = L_2(11)M_{10}, \quad M_{12} = M_{11}M_{11}, \quad M_{23} = F_{23}^{11}M_{22}, \quad M_{24} = M_{23}(M_{12}.2), \quad J_2 = U_3(3)(A_5 \times D_{10}), \quad HS = M_{22}(U_3(5).2), \quad He = (SP_4(2).2)(7^2.SL_2(7)), \quad Ru = L_2(29).^2F_4(2), \quad Suz = G_2(4).U_5(2), \quad Fi_{22} = F_4(2)'.(2.U_6(2)), \quad Co_1 = Co_2.(3.Suz.2).$ 

In the above  $F_{23}^{11}$  denotes a Frobenius group with kernel isomorphic to  $Z_{23}$  and complement isomorphic to  $Z_{11}$ . If L is an exceptional group of Lie type except  $G_2(q), q = 3^n, F_4(q), q = 2^n$  or  $G_2(4)$ , then L does not have a factorization. The simple unitary groups  $U_{2m+1}(q)$  don't have a factorization except when (m,q)=(1,3),(1,5) or (4,2).

To see why the sporadic simple group  $J_1$  is not a factorizable group we recall from [CCNPW] that  $J_1$  has order  $2^3.3.5.7.11.19$  with maximal

subgroups of the form:

$$L_2(11), 2^3: F_7^3, Z_2 \times A_5, F_{19}^6, F_{11}^{10}, D_6 \times D_{10}, F_7^6$$

If  $J_1 = AB$  is a proper factorization of  $J_1$ , then A is contained in a maximal subgroup M of  $J_1$ , therefore  $J_1 = MB$ . We will consider the 7 possible cases for M. If  $M \cong L_2(11), 2^3 : F_7^3, Z_2 \times A_5, F_{11}^{10}, D_6 \times D_{10}$  or  $F_7^6$ , then we have  $19 \times 7 ||B|$ . But then B must be contained in a maximal subgroup of  $J_1$  with order divisible by  $19 \times 7$ , which is not the case. If  $M \cong F_{19}^6$ , then  $7 \times 11 ||B|$  and from the list of maximal subgroups of  $J_1$  we see that  $J_1$  does not have a maximal subgroup with order divisible by  $7 \times 11$ . This final contradiction shows that  $J_1$  is not a factorizable group.

Looking at that factorization of finite groups, the problem goes back to Burnside [Bu] who proved any finite group G of order  $p^aq^b$  is solvable. In the case if p and q are distirct primes and P and Q are p-Sylow and q-Sylow subgroups of G respectively, then G = PQ. N.Ito [It1] proved that if G = AB and if A and B are abelian subgroups of G, then G is metabelian. H. Wielandt [Wi] generalized the result of Ito and proved that if G = AB is a finite group, where A and B are nilpotent subgroups of G, and (|A|, |B|) = 1, then G is solvable. O. Kegel in [Ke] generalized the result obtained by Wielandt without the assumption (|A|, |B|) = 1. Chernikov in [Ch] without the condition of finiteness on G proved that if G = AB, A and B nilpotent subgroups of G satisfying the minimum condition on groups, then G is a solvable group. Ore in [Or] proved that if G is a solvable group and if A and B are two non-conjugate maximal subgroups of G, then G = AB and conversely any maximal factorization of G arises in this way. We recall that a factorization G = AB of a group G is called maximal if both A and B are maximal subgroups of G.

Although there are interesting problems concerning factorizations of infinite groups, but here we are concerned with finite groups. Therefore from now on all groups are assumed to be finite although the general results stated may be true for infinite groups as well. Our aim in this paper is to review results on factorizable finite groups and to state our recent results concerning factorizations G = AB, where one of the factors is the alternating group.

# 2. Preliminary Results.

To start with, first we will mention some elementary results.

**Theorem 2.1.** Let G be a finite group with subgroups A and B, then the following statements are equivalent:

- (i) G = AB,
- (ii) A acts transitively on the set of right cosets of B in G,
- (iii) B acts transitively on the set of right cosets of A in G,
- (iv) If  $\pi_1$  and  $\pi_2$  are the permutation characters of G obtained from (ii) and (iii) respectively, then  $(\pi_1, \pi_2) = 1$ .

**Theorem 2.2.** If G acts transitively on a set  $\Omega$  and H is a transitive subgroup of G, then for any  $\alpha \in \Omega$  we have  $G = G_{\alpha}H$ . In general if G acts k-homogeneously on a set  $\Omega$  and H is a k-homogeneous subgroup of G, then  $G = HG_{(\Delta)}$  where  $\Delta$  is a subset of size k in  $\Omega$ .

To see how theorem 2.2 can be used to obtain factorization of some groups, we mention that any finite group G is a factor of some symmetric group. Because, if G is a group of order m, then we can assume G as a transitive subgroup of  $S_n$  and therefore by Theorem 2.2 we get  $S_n = S_{n-1}G$ .

If we consider the simple group  $L_2(7)$  of order  $2^3.3.7$ , then it is easy to see that  $L_2(7)$  has subgroups of orders 1, 2, 4, 8, 3, 6, 12, 24, 7, 21, 168. Therefore  $L_2(7)$  has faithful transitive actions on sets of cardinalities: 168, 84, 42, 21, 56, 28, 14, 7, 24, 8. Now consider the transitive action of

 $L_2(7)$  on 24 letters. Since the sporadic group  $M_{24}$  has a subgroup isomorphic to  $L_2(7)$ , hence by Theorem 2.2 we will get  $M_{24} = M_{23}.L_2(7)$ . Now we consider the action of  $L_2(7)$  on 8 points. Since  $A_8$  has a subgroup isomorphic to  $L_2(7)$  which is 2-transitive on 8 points, therefore by the first and the second parts of Theorem 2.2 we have the following factorizations of the alternating group  $A_8$ ,  $A_8 = A_7L_2(7) = A_6L_2(7) = S_6L_2(7)$ .

Using part (iv) of Theorem 2.1 and information given in [CCNPW] about permutation characters of certain groups on maximal factorizations of sporadic groups, for instance:  $J_2 = U_3(3).(A_5 \times D_{10}), HS = M_{22}(U_3(5)2).$ 

M.W.Liebeck, C.E.Praeger and J.Saxl in [LPS] determined completely the maximal factorizations of all the finite simple groups and their antomorphism groups. As we mentioned earlier the maximal factorizations of solvable groups were obtained by Ore in [Or]. The main Theorem of [LPS] is the following.

**Theorem 2.3.** ([LPS]) Let L be a finite simple group and let G be a group such that  $L \subseteq G \subseteq Aut(L)$ . Suppose G = AB, where A and B are maximal subgroups of G not containing L. Then the triple (G, A, B) is explicitly known.

The following two results are also obtained about the factorizations of the alternating and symmetric groups.

Theorem 2.4. ([LPS]). Let  $L = A_n (n \ge 5)$  acting naturally on a set  $\Omega$  of n points, and let  $L \le G \le Aut(L)$ . Suppose that G = AB where A and B are arbitrary subgroups of G not containing L. Then one of the following holds:

- (i)  $A_{n-k} \leq A \leq S_{n-k} \times S_k$  for some k with  $1 \leq k \leq 5$ , and B is k-homogeneous on  $\Omega$ .
- (ii) n = 6, 8, 10. If n = 6 the groups A and B have the following property that  $A \cap L = L_2(5)$ ,  $B \cap L = S_3 \wr S_2$ . If n = 8, then A =

 $2^3.L_3(2)$  and  $B \geq Z_5 \times Z_3$ . If n = 10, then  $A = L_2(8)$  or  $L_2(8).3$  and  $A_5 \times A_5 \leq B \leq S_5 \wr S_2$  and B is transitive on  $\Omega$ .

Of course in the above theorem the roles of A and B may be interchanged.

Corollary 2.5. ([LPS]) Let G be  $A_n$  or  $S_n(n > 5)$  and suppose that G = AB with A and B maximal subgroups of G not containing  $A_n$ . Then one of the following holds.

- (i)  $A = (S_{n-k} \times S_k) \cap G, 1 \le k \le 5$ , and B is k-homogeneous.
- (ii)  $n = 6, A = PGL_2(5) \cap B, B = (S_3 \wr S_2) \cap G$ .

If  $k \geq 2$ , then combining the results of W.M.Kantor [Ka] and P.J. Cameron [Ca] it is possible to find the group B mentioned in Theorem 2.4 and Corollary 2.5. If k = 1, then B is a transitive subgroup of  $S_n$  or  $A_n$  and in general it is not difficult to find B in this case.

Factorizations of the symmetric and alternating groups are considered in [WW] as well. We call a factorization G = AB of G exact if  $A \cap B = 1$ . In [WW] all the exact factorizations of the alternating and symmetric groups are found.

# 3. General result on factorizations of group.

Much of recent research is concerned with factorization of finite simple groups, or factorizations involving simple groups. The so called Szep conjecture was proved by E.Fisman and Z.Arad in [FA]. The conjecture states: No simple group has a factorization G = AB with  $Z(A) \neq 1, Z(B) \neq 1$ .

Z.Arad and E.Fisman in [AF] considered factorizations of simple groups G = AB such that (|A|, |B|) = 1. They proved:

**Theorem 3.1.** ([AF]) Let G be a simple group and G = AB where A and B are subgroups of G with (|A|, |B|) = 1. Then one of the followings

holds:

- (i)  $G = A_r$  with  $r \ge 5$  a prime, and  $A \cong A_{r-1}$ , |B| = r.
- (ii)  $G = M_{11}$  and either A is solvable or  $A \cong M_{10}$ .
- (iii)  $G = M_{23}$  and either B is a Frobenius group of order 11.23 or B is cyclic of order 23 and  $A \cong M_{22}$ .
- (iv)  $G = PSL_2(q)$  where either  $q \in \{11, 29, 59\}$  and  $A \cong A_5$  or  $q \not\equiv 1 \pmod{4}, q > 3$  and A is solvable.
- (v)  $G = PSL_r(q)$ , r an odd prime such that (r, q 1) = 1 and either  $G \cong PSL_5(2)$  and |B| = 5.31 or A is maximal parabolic subgroup such that  $PSL_{r-1}(q)$  is involved in A. In particular B is either cyclic or Frobenius.

As a cosequence of the above theorem we obtain the following corollary.

Corollary 3.2. ([AF]) Let G be a group such that G = AB, (|A|, |B|) = 1. Let D be a composition factor of G. Then either D is of type (i) - (v) in Theorem 3.1 or  $\pi(D) \subset \pi(A)$  or  $\pi(D) \subset \pi(B)$ . In particular in the latter case D is either a section of A or of B, respectively. Here  $\pi(G)$  denotes the set of primes involved in the prime factorization of the order of the finite group G.

Before stating the next result we introduce some definitions. Let p be a prime. An elementary abelian p-group  $E_p^n$  is the direct product of n copies of  $Z_p$  and in this case n is called the rank of  $E_p^n$ . If G is a finite group, then the p-rank of G is the maximum p-rank of an elementary abelian p-subgroup of G and is denoted by  $m_p(G)$ .

Now considering the factorizations G = AB and putting restrictions on the structures of A and B, especially on the Sylow 2-subgroups of these groups leads to the following results due to U.Preiser [Pr].

Theorem 3.3. ([Pr]) Let G be a finite group that contains subgroups

A and B with G = AB. Assume that a Sylow 2-subgroup of A has rank 2 and a Sylow 2-subgroup of B is elementary abelian. Then one of the followings holds:

- (i)  $G \cong A$  or B,
- (ii)  $G \cong A \times B$ ,
- (iii)  $G \cong A_6, A \cong A_5 \cong B$ ,
- (iv)  $G \cong A_7$ ,  $A \cong PSL_2(7)$ ,  $B \cong A_5$ ,
- (v)  $G \cong A_9$ ,  $A \cong A_6$  or  $A_7$ ,  $B \cong PSL_2(8)$ ,
- (vi)  $G \cong M_{12}$ ,  $A \cong M_{11}$ ,  $B \cong A_5$  or  $PSL_2(11)$ ,
- (vii)  $G \cong G_2(3^m)$ ,  $A \cong PSL_3(3^m)$ ,  $B \cong^2 G_2(3^m)$ ,  $m \text{ odd, } m \neq 1$ .

Also in the paper [Pr] by U. Prieser the following result is proved which asserts certain finite groups are not factorizable.

**Theorem 3.4.** ([Pr]) (i) The groups  $PSP_4(q)$ , q odd, don't admit a factorization with proper simple subgroups,

- (ii)  $^3D_4(q^3)$ , q odd, don't admit a factorization with two proper simple subgroups,
- (iii)  $PSL_5(q), q \equiv 3 \pmod{4}$  and  $PSU_5(q^2), q \equiv 1 \pmod{4}$ , don't admits factorizations with proper simple subgroups,
- (iv)  $G_2(q), q = p^f, p > 3$ , don't admit a factorization with proper simple subgroups.

In [HOS] finite groups G with factorization G = AB, where A and B have small Sylow 2-subgroups and found.

**Theorem 3.5.** ([HOS]) Let G = AB be a factorization of G and  $G \not\cong A \times B$ . If A and B are finite simple groups such that a Sylow 2-subgroup of B has order 4 or 8, then  $G \cong A_6$  or  $A_7$ .

There are several results by G.Walls in [Wa1] concerning Sylow 2subgroups in a factorization of a finite group and we mention some of them here. **Theorem 3.6.** ([Wa1]) Let G = AB where A and B are simple subgroups of G such that a Sylow 2-subgroup of G is dihedral. Then

- (i)  $G \cong A$  or B,
- (ii)  $G \cong A_6$ ,  $A, B \cong A_5$ ,
- (iii)  $G \cong A_7$ ,  $A \cong A_6$ ,  $B \cong L_2(7)$  or  $A \cong A_5$ ,  $B \cong L_2(7)$ .
- (iv)  $G \cong A \times B$ .

**Theorem 3.7.** ([Wa1]) Let G = AB with A and B simple subgroups of G. If a Sylow 2-subgroup of G is abelian or quasi-dihedral, then  $G \cong A$ , B or  $A \times B$ .

**Theorem 3.8.** ([Wa1]) If G = AB, A and B simple, and if the order of a Sylow 2-subgroup of G is at most 32, then the case (i) - (iv) in Theorem 3.6 occur.

Factorizations of simple groups as the product of two simple subgroups first appeared in the following paper of Ito.

**Theorem 3.9.** ([It2]) If  $G = L_2(q) = AB$  with A and B simple subgroups of G, then

- (i) G = A or B,
- (ii) q = 9 and  $G \cong L_2(9) \cong A_6 \cong A_5 A_5$ .

Extending the above result we have the following results of Gentchev.

**Theorem 3.10.** ([Ge1]) Let G be a simple group of Lie type of Lie rank 1 or 2 over GF(q). Let G = AB with A and B simple non-abelian subgroups of G. Then one of the followings occurs.

- (i)  $G \cong L_2(9), A, B \cong A_5$ ,
- (ii)  $G \cong G_2(4)$ ,  $A \cong J_2$ ,  $B \cong U_3(4)$ ,
- (iii)  $G \cong G_2(q), q = 3^{2n+1} > 3, A \cong L_3(q), B \cong^2 G_2(q),$
- (iv)  $G \cong G_2(q), q = 3^n, A \cong L_3(q), B \cong B_3(q),$
- (v)  $G \cong U_4(3), A \cong L_3(4), B \cong PSP_4(3),$
- (vi)  $G \cong U_4(q), q \not\equiv -1 \pmod{3}, A \cong U_3(q), B \cong PSP_4(q).$

Theorem 3.11. ([Ge2]) Let G be a sporadic simple group such that G = AB where A and B are simple non-abelian subgroups of G. Then

- (i)  $G \cong M_{24}$ ,  $A \cong L_2(23)$ ,  $B \cong M_{22}$  or  $M_{23}$  or  $A \cong L_2(7)$ ,  $B \cong M_{23}$ ,
- (ii)  $G \cong M_{12}, A \cong A_5, L_2(11)$  or  $M_{11}$  and  $B \cong M_{11}$ ,
- (iii)  $G \cong Suz$ ,  $A \cong U_5(2)$ ,  $B \cong G_2(4)$ ,
- (iv)  $G \cong Co_1$ ,  $A \cong G_2(4)$ ,  $B \cong Co_2$ .

# 4. Factorizations with one factor being an altrnating or a symmetric group.

Now we turn to factorizations where one of the factors is isomorphic to an alternating or a symmetric group. O.Kegel and H.Luneberg [KL] proved the following.

Theorem 4.1. ([KL]) Let G = AB be a factorizable group with  $A, B \cong A_5$ . Then  $G \cong A, B, A \times B$  or  $A_6$ .

W.R. Scott in [So] found all finite groups which are equal to the product of  $A_5$  and a finite non-abelian simple group.

**Theorem 4.2.** ([So]) Let G = AB be a finite group, A and B subgrops of G with A a finite non-abelian simple group and  $B \cong A_5$  such that  $Out(A) \not\geq A_5$ . Then one of the following holds:

- (i)  $G \cong A$  or B,
- (ii)  $G \cong A \times B$ ,
- (iii)  $G \cong A_n, A \cong A_{n-1}, n = 6, 10, 12, 15, 20, 30, 60,$
- (iv)  $G \cong A_7, A \cong L_2(7)$ ,
- (v)  $G \cong M_{12}, A \cong M_{11}$ .
- G.Walls in [Wa2] extended the main results of [So] by taking B to be a non-simple group.

Theorem 4.3. ([Wa2]) Suppose that G is a simple group and that

G = AB, where A is a non-abelian simple group and  $B \cong S_5$ . Then one of the following occurs:

- (i)  $G \cong A_n, A \cong A_{n-1}, n = 10, 12, 15, 20, 24, 30, 40, 60, 120,$
- (ii)  $G \cong M_{12}, A \cong M_{11}$
- (iii)  $G \cong A_7$  or  $A_8, A \cong L_2(7)$ ,
- (iv)  $G = A, A \geq A_5$ .

**Theorem 4.4.** ([Wa2]) Suppose that G = AB is not a simple but A is a simple group and  $B \cong S_5$ . Then one of the following cases occurs:

- (i)  $G \cong A \times B$ ,
- (ii)  $G \cong (A \times A_5)\langle \tau \rangle, \tau$  acting as an outer automorphism on both factors,
- (iii)  $G \cong (A_5 \times A_5)\langle \tau \rangle$ ,  $A \cong A_5$  and  $A_5 \times A_5$  is a minimal normal subgroup of G,
  - (iv)  $G \cong S_6, A \cong A_5$ ,
  - (v)  $G \cong A_n \times Z_2$ ,  $A = A_{n-1}$ , n = 10, 12, 15, 20, 24, 30, 40, 60, 120,
- (vi)  $G \cong M_{12}\langle \tau \rangle$ ,  $A \cong M_{11}$ ,  $\tau$  acts as an outer automorphism of order 2 on  $M_{12}$ ,
- (vii)  $G \cong A\langle \tau \rangle$  where  $A \geq A_5$  and  $\tau$  is an outer automorphism of the simple group A which acts as an outer automorphism on the copy of  $A_5$  in A.

In Theorem 4.4 we observe that one of the factors of the factorizable group G is a non-simple group. Walls in [Wa2] proved the following general result.

**Theorem 4.5.** ([Wa2]) Suppose that G = AB is a finite group with subgroups A and B such that  $A \cong S_n, n \geq 5$  and  $B \cong A_5$ . Then one of the following occurs.

- (i)  $G \cong A \times B$ ,
- (ii)  $G \cong S_n = A, n \geq 5$ ,

- (iii)  $G \cong S_{n+1}, n = 5, 9, 11, 14, 19, 29, 59,$
- (iv)  $G \cong (A_n \times A_5)\langle \tau \rangle, \tau$  acting as an automorphism of order 2 on both factors,
- (v)  $n = 5, G \cong (A_5 \times A_5)\langle \tau \rangle, \tau$  an automorphism of order 2,  $A_5 \times A_5$  a minimal normal subgroup of G.

To prove Theorem 4.5 one needs a special treatment of factorizable groups with a factor being non-simple. Lemma 3 of [Wa2] is adjusted to deal with the situations as above.

**Lemma 4.6.** Let G = AB be a factorization of a finite group and A, B subgroup of G such that A is a simple group and B has a unique proper normal subgroup N which is simple non-abelian. Let  $G \not\cong A \times B$  and let M be a minimal normal subgroup of G. Then one of the following holds:

- (i) G = AB = M is simple.
- (ii)  $G = MB, M \cong A \times N, N \cong A$ ,
- (iii)  $G = MB, M \cong NA$  is simple,
- (iv)  $M = A \text{ or } N, [G:AN] = [B:N], AN \cong A \times N,$
- (v)  $M \cap X = 1$ ,  $|X| | [X : A \cap B]$  for  $X \in \{A, B\}$  and  $|M| |A \cap B| = |AM/M \cap BM/M|$ .

Garry Walls in [Wa3] considers factorizable non-simple groups. Some examples of non-simple groups which can be factored as product of two proper subgroups are:

$$2.PSU_4(3^2) = PSP_4(3)PSL_3(4),$$
  
 $3.PSU_4(3^2) = PSP_4(3)PSU_3(3^2),$   
 $3.Suz = PSU_5(2^2)G_2(4).$ 

Motivated by recent results of Walls, We were interested in finding the structure of finite factorizable groups with one factor isomorphic to an alternating group. In [DR] we proved the following results: Theorem 4.7. ([DR]) Let G be a finite group such that G = AB where  $A \cong A_6$  and  $B \cong A_n, n \geq 5$ , then either  $G \cong A \times B$  or G is a simple alternating group as follows:

- (i)  $G \cong A_{n+1}$ , n = 5, 9, 29, 35, 39, 44, 59, 71, 89, 119, 179, 359,
- (ii)  $G \cong A_n, n \geq 6$ ,
- (iii)  $G \cong A_{10}, n = 8$ .

**Theorem 4.8.** ([DR]) Let G be a finite group such that G = AB,  $A \cong A_6$  and  $B \cong S_n$ ,  $n \geq 5$ . Then one of the following occurs:

- (i)  $G \cong A_6 \times S_n$ ,
- (ii)  $G \cong A_{10} \cong A_6 S_8, n = 8$ ,
- (iii)  $G \cong (A_6 \times A_6)\langle \tau \rangle, \tau$  an automorphism of order 2 and  $A_6 \times A_6$  is the minimal normal subgroup of G, n = 6,
  - (iv)  $G \cong A_{n+1}$ , n = 5, 9, 14, 19, 35, 39, 44, 59, 71, 89, 119, 179, 359,
  - (v)  $G \cong S_n, n \geq 6$ ,
  - (vi)  $G \cong A_{10} \times Z_2$ , n = 8,
- (vii)  $G \cong (A_6 \times A_n)\langle \tau \rangle, n \geq 5$ , where  $\tau$  acts as an automorphism of order 2 on both factors.

In [DRW] we extended the result obtained in [So] by considering factorizable groups with one factor isomorphic to the alternating group  $A_6$  or  $S_6$ . These results are as follows:

**Theorem 4.9.** ([DRW]) Let G be a simple group and G = AB be a proper factorization of G with A simple and  $B \cong A_6$ . Then one of the following occurs:

- (i)  $G \cong A_n$ ,  $A \cong A_{n-1}$ , n = 10, 15, 20, 30, 36, 40, 45, 60, 72, 90, 120, 180, 360,
- (ii)  $G \cong A_{10} \cong A_8 A_6$ ,
- (iii)  $G \cong A_8 \cong L_2(7)A_6$ ,
- (iv)  $G \cong A_9 \cong L_2(8)A_6$ ,

- (v)  $G \cong A_7 \cong L_2(7)A_6$
- (vi)  $G \cong O_8^+(2) \cong S_6(2)A_6$ .

Theorem 4.10. ([DRW]) Let G be a simple group with non-trivial factorization G = AB, where A and B are subgroups of G with A simple and  $B \cong S_6$ . Then we have one of the following cases:

- (i)  $G \cong A_n \cong A_{n-1}S_6$ , n = 10, 12, 15, 20, 30, 36, 40, 45, 60, 72, 80, 90, 120, 144, 180, 240, 360, <math>g720,
- (ii)  $G \cong A_{10} \cong A_8 S_6$ ,
- (iii)  $G \cong A_8 \cong L_2(7)S_6$ ,
- (iv)  $G \cong A_9 \cong L_2(8)S_6$ ,
- (v)  $G \cong S_6(2) \cong U_3(3)S_6$ ,
- $(vi) G \cong S_4(4) \cong L_2(16)S_6,$
- (vii)  $G \cong S_8(2) \cong O_8^-(2)S_6$ ,
- (viii)  $G \cong O_8^+(2) \cong S_6(2)S_6$ .

Theorem 4.11. ([DRW]) If G is not a simple group and has a factorization G = AB with A simple and  $B \cong S_6$ , then one of the following occurs:

- (i)  $G \cong A \times B$ ,
- (ii)  $G \cong A(\tau)$  where A is a simple group containing  $A_6$  and  $\tau$  is an outer automorphism of both A and  $A_6$ .
  - (iii)  $G \cong S_n$  or  $A_n \times Z_2$ ,  $A \cong A_{n-1}$ , n = 10, 15, 20, 30, 36, 40, 45, 60, 72, 90, 120, 180, 360,
  - (iv)  $G \cong S_{10}$  or  $A_{10} \times Z_2$ ,  $A \cong A_8$ ,
  - (v)  $G \cong S_p$  or  $A_8 \times Z_2$ ,  $A \cong L_2(7)$ ,
  - (vi)  $G \cong S_9$  or  $A_9 \times Z_2$ ,  $A \cong L_2(8)$ ,
  - (vii)  $G \cong S_7, A \cong L_2(7)$ ,
- (viii)  $G \cong O_8^+(2)\langle \tau \rangle$  or  $O_8^+(2) \times Z_2$ ,  $A \cong S_6(2)$ ,  $\tau$  an outer automorphism of  $O_8^+(2)$ ,

- (ix)  $G \cong (A_6 \times A_6) \langle \tau \rangle$ ,  $A \cong A_6$ , where  $A_6 \times A_6$  is the minimal normal subgroup of G,
- (x)  $G \cong (A \times A_6)\langle \tau \rangle, \tau$  acting as an outer automorphism on both factors.

In general, the structure of groups which are the product of an alternating and a symmetric group is given in [Da]. This paper is a generalization of [DR].

**Theorem 4.12.** ([Da]) Let G be a finite group with two subgroups A and B such that G = AB, where A is isomorphic to some alternating group  $A_r$  and B is isomorphic to some symmetric group  $S_n, r, n \geq 5$ . Then one of the following occurs:

- (i)  $G \cong A_r$  or  $S_n$ , the trivial factorization,
- (ii)  $G \cong A_r \times S_n$ ,
- (iii)  $G \cong A_{10}, S_{10}$  or  $A_{10} \times Z_2$ , where  $A \cong A_6$  and  $B \cong S_8$ ;
- (iv)  $G \cong A_{15}, S_{15}$  or  $A_{15} \times Z_2$ , where  $A \cong A_7$  or  $A_8$  and  $B \cong S_{13}$ ,
- (v)  $G \cong A_{r+1}$  (or  $S_{n+1}$ ),  $A \cong A_r$ ,  $B \cong S_n$ , where  $A_{r+1}$  and  $A_{r+1} \times Z_2$  (or  $S_{n+1}$ ) have a transitive subgroup isomorphic to  $S_n$  (or  $A_r$ ),
- (vi)  $G \cong (A_r \times A_r)\langle \tau \rangle, \tau$ , an outer automorphism of order 2 of  $A_r$  interchanging the two  $A_r$ 's and  $A_r \times A_r$  is the minimal normal subgroup of G,
- (vii)  $G \cong (A_r \times A_n)\langle \tau \rangle, \tau \neq n$ , where  $\tau$  acts as an automorphism of order 2 on both factors.

(viii) 
$$G \cong S_r$$
 or  $S_{r+1}$ .

At the end of this paper we would like to mention some interesting problems concerning factorizable finite groups.

Q1: Find the structure of all finite groups  $G = L_2(7)B$  where B is a non-abetian simple group.

Q2: Find all finite groups G such that  $G = L_2(7)B$  where B is any

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finite group with the property that B' = N is a simple group which is the unique normal subgroup of B and [B:N] = 2.

Q3: Let G be a simple factorizable group such that G = AB, where A and B are proper subgroups of G. If A and B are perfect groups is it true that one of A or B must be a simple group? (A is called a perfect group if A' = A). The answer to this question in the case of alternating G is yes.

Q4: Obtain general results about factorizations of infinite alternating or symmetric groups.

Q5: Find the structure of all finite groups  $G = L_2(q)B$ , where B isomorphic to the symmeteric group  $S_n, n \geq 5$ .

**Q6**: In general find the structure of all finite groups G such that G = AB where A is a simple group and B is isomorphic to a symmetric group.

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Proceedings of the Monthly Colloquium of the Iranian Mathematical Society Vol. 2, pp 39-57 (2004)

# Algebra in a Category: Injectivity in Equational Classes

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Abstract: Much of classical set based universal algebra can be and, if we may say so, should be, studied in a general topos rather than in the topos Set of sets.

Each topos carries its own logic, which is not necessarily the same as the classical logic of the topos Set. Thus, the classical notions of mathematics may behave differently in different topoi.

The notion from classical universal algebra we have been chosen for this talk is injectivity, which is one of the most central and useful notion in many mathematical disciplines.

The main purpose of this talk is to describe the relationship between the class  $mod\Sigma$  of models of a set  $\Sigma$  of equations in the category Set and the corresponding class  $mod(\Sigma, E)$  of models of  $\Sigma$  in a suitable category E with respect to injectivity and some related notions.

MSC (2000): 08B30, 08C05, 18C05

Keywords: Universal algebra, Grothendieck topos, essential boundedness, injectivity.

The basic nature of our results is that, for any given  $\Sigma$ , whatever holds in Set, concerning this notion, also holds in E, provided E has some special properties, in particular when E is a Grothendieck topos.

The talk is general and the results have already been published. For the details of the proof, see [10].

#### 1. Introduction

1.1 Universal Algebra: "Much of the beauty of mathematics is derived from the fact that it affords abstraction. Not only does it allow one to see the forest rather than the individual trees, but it offers the possibility for the study of the structure of the entire forest, in preparation for the next stage of abstraction-comparing forests."

Mathematicians at the beginning of the twentieth century were confronted with a large number of algebraic systems such as groups, rings, quaternions, Lie algebras, number rings, algebraic number fields, vector spaces, Boolean algebras, lattices, etc. A. N. Whitehead [1898] tried to place these diverse algebraic systems within a common frame work. Universal algebra as understood today goes back to the 1930's and it emerged as a natural development of the abstract approach to algebra initiated by Emmy Noether.

Universal Algebra studies features common to familiar algebraic systems mentioned above. Although, "one can become a very good mathematician without being a professional logician even though logical thought is central to mathematics". This goes for algebraists with regard to Universal Algebra. But, such a study places the algebraic notions in their proper setting. It often reveals connections between seemingly different concepts and helps to systematize one's thoughts. However, it does not usually solve the whole problem for us, tidies up a mass of rather trivial

details and thus allows us to concentrate our powers on the hard core of the problem.

Universal algebra has grown very rapidly in the last thirty or forty years. Not only the literature expanded rapidly, but also the problems have become more sophisticated and the results deeper. Young mathematicians entering this field today are indeed fortunate, for there are hard and interesting problems to be attacked and sophisticated tools to be used.

In the last two decades universal algebra has become useful and important in combinatorics and theoretical computer sciences. In particular, structural aspects such as syntax and semantics, data abstraction, etc., are mainly investigated by methods of universal algebras [20, 21, 23, 24, 25, 34].

One of the fundamental ideas of universal algebra is the representation of logical notions in nonlogical terms. The famous Birkhoff Variety Theorem

states that a class of algebras is equationally definable iff it is closed under subalgebras, homomorphic images and products; such a class is called a variety. This characterization result was the starting point of universal algebra.

The general theory of algebras borrows techniques and ideas from lattice theory, logic, and category theory and derives inspirations from older, more specialized branches of algebras such as the theories of groups, rings, and modules.

1.2 Category Categorical methods of speaking and thinking has become widespread in mathematics because they achieve a unification of parts of different mathematical fields; frequently they bring simplifications and provide the impetus for new developments.

Categories, initially a convenient way of formulating exact sequences, and axiomatic homology theory, obtained independent life in the works of Ehresman, Kan, Maclane, Eilenberg, Barr, Freyd, Gray, Lawvere, Linton, Tierney, and others.

A category may be thought of in the first instance as a universe for a particular kind of mathematical discourses. Such a universe is determined by specifying a certain kind of "objects" and a certain kind of "arrows" that links different objects. The most general universe of current mathematical discourse is the category Set of sets with functions between them. Many basic properties of sets and set theoretic operations can be described by reference to the arrows in Set, and these descriptions can be interpreted in any category by means of its arrows. So the question that arises is "when does a category look and behave like Set?" A vague answer is "when it is (at least) a topos".

1.3 Topos In 1963, Lawvere tried to find a purely categorical foundation for all mathematics, beginning with an appropriate axiomatization of the category of sets. When Lawvere heard of the properties of Grothendieck topoi, he soon observed that such a topos admits basic operations of set theory such as the formation of sets  $Y^X$  of all functions from X to Y and of power sets  $\mathcal{P}(X)$ . Subsequently, Lawvere and Tierney, working together at Dalhousie University, defined in an elementary way, free of all set-theoretic assumptions, the notion of an "elementary topos".

A topos is formally a category which has finite limits, exponentiations (abstracting the function set  $B^A$ ) and subobject classifier (abstracting the truth set  $2 = \{0, 1\}$ ).

Recall that, for a category C with finite limits, we say that C has exponentiations (exponentials) if for every objects A and B,

there is an object  $B^A$  together with an arrow  $ev: B^A \times A \to B$  (called evaluation) such that for every arrow  $g: C \times A \to B$  there is a unique arrow  $\hat{g}: C \to B^A$  with  $ev \circ (\hat{g} \times id_A) = g$ .

We also say that C has subobject classifier if there exists an object  $\Omega$  with an arrow  $t: 1 \to \Omega$  (called the truth arrow) such that for every monomorphism  $f: B \to A$  there is a unique arrow  $\chi_f: A \to \Omega$  (called the classifying arrow) making the square

$$\begin{array}{cccc}
B & \xrightarrow{f} & A \\
! & \downarrow & \downarrow & \chi_f \\
1 & \xrightarrow{t} & \Omega
\end{array}$$

a pullback square.

In fact, a topos is informally a category which looks and behaves very much like the category of sets. One of the first examples of a topos is  $Set^{C^{\circ r}}$ .

A Grothendieck topos  $I\!E$  is a reflective subcatgory of  $Set^{C^{op}} = \hat{\mathcal{C}}$   $(i:I\!E \rightleftarrows \hat{\mathcal{C}}:R,\ R\dashv i)$ , for some small category  $\mathcal{C}$ , whose reflection functor R preserves finite limits. For example, for a topological space X, the category PreshX, of presheaves on X, and the category ShX, of sheaves on X, are Grothendieck topoi.

Again, one can become a good mathematician without being a professional category theorist. But

"How can you do "new maths" problems with an "old math" mind?"- Charlie Brown.

"Virtually all algebraic notions in category theory are parodies of their parents in the most "classical" of categories · · · the category of left A-modules." H. Bass.

Here we briefly study injectivity, the most central notion in classical universal algebra, modelled in a topos, rather than in the category Set of sets. It is intended to provide a deeper understanding of the real features of this algebraic notion and to show that how a classical set-based Universal Algebra can be, as we may say so, it should be, studied in a topos (or category) theoretic setting. For more information about this approach see [3, 5, 6, 7, 10, 11, 12, 13, 14, 18, 30].

# 2. Algebra in a category

Let IK be a finitely complete category (that is, a category which has finite products and equalizers, in particular, it has a terminal object 1).

**2.1 Definition** Given a family  $\tau = (n_{\lambda})_{\lambda \in \Lambda}$  of finite cardinal number  $n_{\lambda}$ , indexed by a set  $\Lambda$ , an algebra in K is an entity  $(A, (\lambda_A)_{\lambda \in \Lambda})$ , where A is an object of K and, for each  $\lambda \in \Lambda$ , the  $\lambda$  th operation  $\lambda_A : A^{n_{\lambda}} \to A$  is a morphism in K. The family  $\tau = (n_{\lambda})_{\lambda \in \Lambda}$  is called the type of this algebra. The algebra  $(A, (\lambda_A)_{\lambda \in \Lambda})$  is simply denoted by A.

In the sequel, all algebras are of the same type  $\tau$ .

**2.2 Definition** A homomorphism from an algebra A to an algebra B in  $I\!K$  is a morphism  $h:A\to B$  in  $I\!K$  such that, for each  $\lambda\in\Lambda$ ,  $\lambda_B\circ h^{n_\lambda}=h\circ\lambda_A$ . That is, the following diagram commutes, for each  $\lambda\in\Lambda$ :

$$\begin{array}{cccc} A^{n_{\lambda}} & \xrightarrow{h^{n_{\lambda}}} & B^{n_{\lambda}} \\ \lambda_{A} & \downarrow & & \downarrow & \lambda_{B} \\ & A & \xrightarrow{h} & B \end{array}$$

Clearly the identity morphism on the algebra A in  $I\!\!K$  is a homomorphism  $1_A:A\to A$  and for composable homomorphisms  $A\stackrel{g}{\to} B\stackrel{f}{\to} C,$   $f\circ g$  is a homomorphism.

As a result, the collection of all algebras (of the type  $\tau$ ) in  $I\!K$  and homomorphisms between them forms a category denoted by  $Alg(\tau)I\!K$  (or by  $Alg(\tau)$  if  $I\!K = Set$ ).

**2.3 Remark** For  $A \in Alg(\tau)K$  and any natural number n, the set  $Hom_K(A^n, A)$  of all morphisms in K from  $A^n$  to A can easily be made into an algebra of the type  $\tau$  in Set, by defining the  $\lambda$ th operation as

$$\lambda(\phi_1,\cdots,\phi_{n_\lambda})=\lambda_A\circ\prod_{i=1}^{n_\lambda}\phi_i$$

for any  $\phi: A^n \to A$   $(i = 1, \dots, n_{\lambda})$  in IK, where  $\prod_{i=1}^{n_{\lambda}} \phi_i: A^n \to A^{n_{\lambda}}$  is the morphism determined by  $\phi_i$ 's. That is,  $\lambda_{Hom_K(A^n,A)}$  takes  $(\phi_1, \dots, \phi_{n_{\lambda}})$  to

$$\lambda \circ \prod_{i=1}^{n_{\lambda}} \phi_i : A^n \to A^{n_{\lambda}} \to A$$

Let FX be the absolutely free algebra in Set of the type  $\tau$  on a set  $X = \{x_1, \dots, x_n\}$  of n elements. Extend the map

$$X \to Hom_K(A^n, A)$$
  
 $x_i \mapsto p_i : A^n \to A$ 

 $(p_i \text{ the } i \text{ th projection})$ , freely to

$$\phi: F \to Hom_K(A^n, A)$$

and denote  $\phi(p)$  by  $p_A$  for any  $p \in F$ .

**2.4 Definition** A law (identity or equation) over  $\Lambda$  in the set X of variables is any pair  $\sigma = (p, q) \in F \times F$ , denoted by p = q.

We say A satisfies the equation p = q, written  $A \models (p = q)$ , if  $p_A = q_A$ .

The full subcategory of  $Alg(\tau)I\!\!K$  given by the class of all algebras in  $I\!\!K$  satisfying every equation of a set  $\Sigma$  of equations is denoted by  $mod(\Sigma,I\!\!K)$ , or by  $mod\Sigma$ , if  $I\!\!K=Set$ , and is called an equational category of algebras.

#### 2.5 Examples

1) A group in the category IK = ShX, the category of all sheaves on a topological space X, is an entity  $(G; *, ()^{-1}, e)$ , where G is a sheaf on X and

$$*: G \times G \rightarrow G$$
 ,  $()^{-1}: G \rightarrow G$  ,  $e: 1 \rightarrow G$ 

are morthpisms in ShX (that is, natural transformations) such that the following diagrams are commutative:

i) (associativity of \*)

ii) (the identity condition)

iii) (the inverse condition)

$$(1,()^{-1}) \quad \downarrow \qquad e \downarrow \qquad \downarrow \qquad (()^{-1},1)$$

$$G \times G \stackrel{*}{\longrightarrow} G \stackrel{*}{\longleftarrow} G \times G$$

Notice that the commutativity of the above diagrams in ShX means that the diagrams are commutative at each  $U \in O(X)$  in Set. In fact, a sheaf G is a group in ShX iff GU is a group in Set (see also 2.8).

A homomorphism between groups in ShX, is a sheaf morphism (natural transformation) which is a group homomorphism at each  $U \in O(X)$ .

2) A ring in the category MSet, of all M-sets (that is, sets with an action of the monoid on it), is an entity (R; +, .., -, 0), where R is an M-set and

$$+: R \times R \to R$$
,  $: R \times R \to R$ ,  $-: R \to R$ ,  $0: \{\bullet\} \to R$ 

are action preserving maps satisfying the ring axioms.

In fact, a ring in MSet is an M-set which is also a ring in Set whose ring operations are action preserving maps.

A homomorphism between rings in MSet is an action preserving ring homomorphism.

3) A monoid in the category R - Mod, of all left R-modules, is an entity (M; ., e), where M is an R-module and

$$.: M \times M \to M \ , \ e: \{ \bullet \} \to M$$

are R-module homomorphisms satisfying the following identities:

- i) (m.n).k = m.(n.k)
- ii) e.m = m = m.e

So, an R-module which is also a monoid whose monoid operations are R-module homomorphisms is a monoid in R-Mod.

A homomorphism between monoids in R-Mod is an R-module homomorphism which is also a monoid homomorphism.

**2.6 Remark** Let  $k : IK \to IL$  be a functor, preserving finite limits. Then k induces another functor

$$\overline{k}: Alg(\tau) I\!\! K \to Alg(\tau) I\!\! L$$

defined on objects by

$$\overline{k}A = (A, (k\lambda_A)_{\lambda \in \Lambda})$$

and on homomorphisms  $f: A \to B$ , by

$$\overline{k}(f) = k(f)$$

If  $\sigma=(p,q)$  is an equation, then  $A\models(p=q)$  implies that  $p_A=q_A$ , and hence  $kp_A=kq_A$  which implies that  $p_{\overline{k}A}=q_{\overline{k}A}$ ; and thus  $\overline{k}A\models(p=q)$ . We thus get a functor

$$\overline{k}: mod(\Sigma, I\!\!K) 
ightarrow mod(\Sigma, I\!\!L)$$

for any given set  $\Sigma$  of equations.

Recall that a set  $\Phi$  of objects of a category  $I\!K$  is said to be a set of generators if for every pair of morphisms  $f,g:A\to B$  with  $f\neq g$  there exists  $G\in \Phi$  and a morphism  $h:G\to A$  such that  $fh\neq gh$ .

**2.7 Lemma** Let IK have a set  $\Phi$  of generators. Then, for any  $A \in Alg(\tau)IK$  and a set  $\Sigma$  of equations,  $A \in mod(\Sigma, IK)$  iff  $\overline{h}_G(A) \in mod\Sigma$  for each  $G \in \Phi$ , where  $h_G = Hom_K(G, -)$ .

**Proof** Applying remark 2.6 to the functor  $h_G$ , we get that for  $A \in mod(\Sigma, I\!\!K)$ ,  $\overline{h}_G(A) \in mod\Sigma$ . Conversely, let  $A \in Alg(\tau)I\!\!K$  and  $\overline{h}_G(A) \in mod\Sigma$  for each  $G \in \Phi$ . Let  $\sigma = (p,q)$  be an equation in  $\Sigma$ . By hypothesis, for all  $G \in \Phi$ ,  $\overline{h}_G(A) \models \sigma$ , that is,  $p_{\overline{h}_G(A)} = q_{\overline{h}_G(A)}$ . So,  $h_G(p_A) = h_G(q_A)$  for all  $G \in \Phi$ . Since  $\Phi$  is a set of generators, the preceding equalities yield that  $p_A = q_A$ . Thus  $A \models \sigma$ , and hence  $A \in mod(\Sigma, I\!\!K)$ .

**2.8 Corollary** The category  $Alg(\tau)\hat{C}$  is isomorphic to the category of all  $Alg(\tau)$ -valued presheaves on C. And, for a Grothendieck topos  $I\!\!E$ , and a set  $\Sigma$  of equations,  $A \in mod(\Sigma, I\!\!E)$  iff  $AU \in mod\Sigma$  for all  $U \in C$ .

## 3. Injectivity in equational classes

A natural question to ask would be, what is the relationship between the behaviour of a certain classical algebraic notion in  $mod\Sigma$  and in  $mod(\Sigma, I\!\!K)$ . In the following, we briefly consider the notion of injectivity, and show that the properties of  $mod\Sigma$ , regarding injectivity, survive the passage to  $mod(\Sigma, I\!\!E)$ , for a set  $\Sigma$  of equations and an arbitrary Grothendieck topos  $I\!\!E$  (fixed from now on); for the details and some particular cases see for example [10], [13], [30]. For the case of equational classes of algebras in Set see for example [2], [32].

**3.1 Definition** An object E in a category K is called *injective* if, for any monomorphism  $B \xrightarrow{\overline{f}} C$  and any morphism  $B \xrightarrow{\overline{f}} E$  there exists a morphism  $C \xrightarrow{\overline{f}} E$  with  $\overline{f} \circ i = f$ ; that is the following diagram commutes:

$$\begin{array}{cccc} & B & \xrightarrow{i} & C \\ f & \downarrow & \swarrow & & \overline{f} \\ & E & & \end{array}$$

Replacing B, in the above definition, by E and f by the identity morphism on E, we get the definition of an absolute retract object E in a category.

**3.2 Definition** A monomorphism  $h:A\to B$  in a category K is called **essential** if, for any morphism  $g:B\to C$ , wherever  $g\circ h$  is a monomorphism, then so is g.

- 3.3 Lemma In  $mod(\Sigma, E)$ , for a Grothendieck topos E, we have
- i) any composite of essential monomorphisms is an essential monomorphism, and
- ii) any direct limit of essential monomorphisms is an essential monomorphism.
- **3.4** Lemma In  $mod(\Sigma, I\!\!E)$ , for any monomorphism  $h: A \to B$  there exists a homomorphism  $g: B \to C$  with  $g \circ h$  an essential monomorphism.

**Proof** Take  $\Theta_0$  to be the maximal congruence on B such that  $B/\Theta_0 \in mod(\Sigma, I\!\!E)$  and  $A \to B \to B/\Theta_0$  is a monomorphism. This composition is then essential.

- 3.5 Corollary In  $mod(\Sigma, I\!\!E)$ , an algebra A is an absolute retract iff it has no proper essential extension.
- **3.6 Definition** A category IK is called *essentially bounded* if, for each  $A \in IK$  there exists, up to isomorphism, only a set of essential extensions in IK.
- **3.7** Proposition  $mod(\Sigma, I\!\!E)$  is essentially bounded iff  $mod\Sigma$  is essentially bounded.
- 3.8 Definition In any category IK, pushouts transfer monomorphisms

if, for any pushout diagram

$$\begin{array}{cccc} & A & \xrightarrow{f} & B \\ u & \downarrow & & \downarrow & v \\ & C & \xrightarrow{g} & D \end{array}$$

whenever f is a monomorphisms, then g is also a monomorphism.

If IK has pushouts, the above is equivalent to say that, any diagram

$$\begin{array}{ccc}
A & \xrightarrow{f} & E \\
u & \downarrow & \\
C
\end{array}$$

with f a monomorphism can be completed to a commutative diagram

$$\begin{array}{ccc}
A & \xrightarrow{f} & B \\
u & \downarrow & & \downarrow & v \\
C & \xrightarrow{g} & D
\end{array}$$

with g a monomorphism.

- **3.9 Proposition** Pushout transfer monomorphisms in  $mod(\Sigma, \mathbb{E})$  iff they do in  $mod\Sigma$ .
- **3.10 Lemma** The category  $mod(\Sigma, I\!\!E)$  has enough injectives iff it is essentially bounded and pushouts transfer monomorphisms.
- **3.11 Proposition** The category  $mod(\Sigma, I\!\!E)$ ) has enough injectives iff  $mod\Sigma$  has enough injectives.

# 4. Behaviour of injectivity in $mod(\Sigma, I\!\!E)$

Banaschewski in [2] calls the notion of injectivity in a category IK properly behaved iff the following three propositions hold:

- (I) For any  $A \in IK$ , the following conditions are equivalent:
  - (I1) A is injective.
  - (I2) A is an absolute retract.
  - (I3) A has no proper essential extensions.
- (E) Every  $A \in IK$  has an injective hull, unique up to isomorphisms.
- (H) For any monomorphism  $f: A \rightarrow B$ , the following conditions are equivalent:
  - (H1)  $f: A \to B$  is an injective hull of A.
  - (H2)  $f: A \to B$  is a maximal essential extension.
  - (H3)  $f: A \rightarrow B$  is a minimal injective extension.

For  $IK = mod(\Sigma, IE)$ , we now have the following.

- 4.1 Proposition For  $mod(\Sigma, E)$ , the following are equivalent:
  - (i) Injectivity is properly behaved.
  - (ii)  $mod(\Sigma, I\!\!E)$  has enough injectives.
- (iii)  $mod(\Sigma, I\!\!E)$  is essentially bounded and pushout transfer monomorphisms.

In particular, one has, by proposition 3.11:

**4.2** Corollary Injectivity is properly behaved in  $mod(\Sigma, \mathbb{E})$  iff it is properly behaved in  $mod(\Sigma)$ .

#### 4.3 Examples

1) Recall that, the category Set has enough injectives. So, if we take  $\Sigma$  to be the empty set, then  $mod\Sigma$  as the full subcategory of all algebras

of the type  $\tau = \emptyset$ , is Set and hence has enough injectives. Using 3.11, this implies that for any Grothendieck topos  $I\!\!E$ ,  $mod(\Sigma, I\!\!E) = I\!\!E$  has enough injectives. In particular, the category MSet, of M-sets have enough injectives. Thus injectivity is properly behaved in such categories.

- 2) The category *Boo*, of Boolean algebras has enough injectives (the power set of each set is injective in *Boo*). So, the category of Boolean algebras in any Grothendieck topos has enough injectives.
- 3) The category Ab, of abelian groups has enough injectives (recall that here injectivity means divisibility). So, the category of abelian groups in any Grothendieck topos, in particular in ShX, has enough injectives.
- 4.4 Note For certain  $\Sigma$ , one has characterization of the injective  $A \in mod\Sigma$  by properties of A in terms of its elements and subsets. For example: divisibility for abelian groups, completeness for Boolean algebras (the Sikorski Theorem), and completeness and Booleanness for distributive lattices. A good question to ask is to what extent, that is for what E, such characterizations remain valid in  $mod(\Sigma, E)$ . Banaschewski [3] shows that divisibility = injectivity for abelian groups in the category ShL of sheaves on a frame L iff L is Boolean. Also Ebrahimi [13] shows that for Boolean algebras in the topos of M-sets, injectivity implies "internal" completeness, but the converse is not true. Also, Mahmoudi [30] defines internal injectivity for the category MBoo of Boolean algebras in the topos MSet and finds some equivalent conditions in which the internal version of the Sikorski Theorem holds.

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Proceedings of the Monthly Colloquium of the Iranian Mathematical Society Vol. 2, pp 59-104 (2004)

# Approximation in Complex and Real Lipschitz Algebras

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Abstract: For a compact metric space (X,d) and  $0 < \alpha \le 1$ , the algebra of all complex valued functions f on X for which  $p_{\alpha}(f) = \sup \left\{ \frac{|f(x) - f(y)|}{d^{\alpha}(x,y)} : x,y \in X, x \neq y \right\} < \infty$  is denoted by  $\operatorname{Lip}(X,\alpha)$  and the subalgebra of those  $f \in \operatorname{Lip}(X,\alpha)$  for which  $\frac{|f(x) - f(y)|}{d^{\alpha}(x,y)} \longrightarrow 0$  as  $d(x,y) \longrightarrow 0$  is denoted by  $\operatorname{Lip}(X,\alpha)$ . These algebras are known as the Lipschitz algebras and were first studied by D.R. Sherbert in 1964. The Lipschitz algebras are Banach algebras under the norm  $||f|| = ||f||_X + p_{\alpha}(f)$ , where  $||f||_X$  is the uniform norm on X.

The Hedberg's Theorem concerning the density of certain subalgebras of  $lip(X, \alpha)$  in  $lip(X, \alpha)$  will be discussed and, as a consequence, the density of Lip(X, 1) in  $lip(X, \alpha)$  for  $0 < \alpha < 1$  will be shown.

When X is a compact subset of  $\mathbb{C}^n$  there are interesting subalgebras of Lipschitz algebras which are generated by polynomials, rational functions with poles off X, functions which are analytic in some neighbourhood of X, or functions which are analytic in the interior of X.

MSC (2000): 46J10, 46J15, 46J20, 41A10, 41A20.

Keywords and Phrases: Approximation, Banach function algebras, complex Lipschitz algebras, real Lipschitz algebras, Fréchet Lipschitz algebras,

The approximation problem concerning the equality among the above algebras and the density of these algebras in certain uniform algebras will be investigated. In particular, when X is a compact plane set with planar measure zero, it is shown that  $lip_R(X,\alpha) = lip(X,\alpha)$ , where  $lip_R(X,\alpha)$  is the norm closure of rational functions with poles off X. Another important result is the density of  $D^1(X)$  in  $lip_R(X,\alpha)$ , where  $D^1(X)$  is the Banach function algebra of functions with continuous derivative on the perfect compact plane set X.

We introduce Lipschitz algebras of differentiable complex functions on perfect compact plane sets and then interesting subalgebras of these algebras are defined and the approximation problem among these kinds of Lipschitz algebras will be discussed.

To introduce the real Lipschitz algebras on a compact metric space (X,d) we assume that the map  $\tau: X \longrightarrow X$  is a topological involution on X such that  $d(\tau(x),\tau(y)) \leq Cd(x,y)$  for all  $x,y \in X$ , where C is a positive constant. Let  $\sigma: C(X) \longrightarrow C(X)$  be the algebra involution on C(X) induced by  $\tau$ ; i.e.  $\sigma h = \bar{h} \circ \tau$ . The real Lipschitz algebras are introduced by

$$Lip(X, \tau, \alpha) = \{ h \in Lip(X, \alpha) : \sigma(h) = h \} \quad (0 < \alpha \le 1)$$
  
 
$$\ellip(X, \tau, \alpha) = \{ h \in \ellip(X, \alpha) : \sigma(h) = h \} \quad (0 < \alpha < 1)$$

The extension of the Hedberg's Theorem for the real Lipschitz algebra  $lip(X, \tau, \alpha)$  will be discussed and the density of  $Lip(X, \tau, 1)$  in  $lip(X, \tau, \alpha)$  will be shown.

Another important class of Lipschitz algebras is the Fréchet Lipschitz algebras  $FLip(X,\alpha)$  and  $Flip(X,\alpha)$ , where X is a hemicompact metric space. The density of FLip(X,1) in  $Flip(X,\alpha)$  and the Hedberg's Theorem for the Fréchet Lipschitz algebra  $Flip(X,\alpha)$  will be discussed.

We will try to present a brief historical background on the progress of the subject.

## 1. Historical Background

Although the notion of Lipschitz functions is very old and these interesting functions have been studied for many decades, interest in the Banach space and Banach algebra theory of Lipschitz functions was not developed until 1955. As it is mentioned in the paper of D.R. Sherbert [30], the only work known to us till 1955, which treats the

space of Lipschitz functions as a Banach algebra, was done by S.B. Myers [22]. His paper contains a summary of results only, the proofs never published because of his untimely death in 1955. The proofs of many of the unproved statements of Myers have been supplied by Sherbert in [30], and in some cases, his results have been extended to more general settings. Sherbert also discussed different aspects of Banach algebras of Lipschitz functions in [29].

Since 1964 some other mathematicians have been working on real and complex Lipschitz algebras and obtained interesting results in the following fields: Structure of ideals, maximal ideal spaces, point derivation, automatic continuity, eventual continuity, amenability and weak amenability, differentiability of Lipschitz functions, the Stone-Weierstrass theorem, extreme points, peak points, Shilov boundary and Choquet boundary, isometries between Banach spaces of Lipschitz functions, closed ideals, Lipschitz algebras of differentiable functions, and the approximation problem.

A good collection of the works in Lipschitz algebras can be found in the recent book of Nik Weaver [32], though it does not contain many results in this field. For more advanced account on some aspects of the complex Lipschitz algebras, including the Frechet Lipschitz algebras, one can refer to the recent and interesting monograph of H.G. Dales [7; Sec. 4.4].

In this paper we are going to study an interesting aspect of Lipschitz algebras, called the approximation problem. One of the famous results in approximation is the Hedberg's theorem, which is, in fact, the Stone-Weierstrass theorem for Lipschitz algebras and will be discussed here for real and complex Lipschitz algebras as well as the Frechet Lipschitz algebras.

## 2. Complex Lipschitz Algebras

Let X denote a compact Hausdorff space and C(X) denote the commutative complex unital Banach algebra of all continuous complex-valued functions on X with the pointwise operations and with respect to uniform norm on X.

**Definition 2.1.** A complex Banach function algebra on X, is a complex subalgebra of C(X) which separates the points of X, contains the constant function 1 and it is complete under an algebra norm. If the norm of a complex Banach function algebra is the uniform norm on X, then it is called a *complex uniform* (function) algebra on X.

If A is a complex Banach function algebra on X the space of maximal ideals of A is denoted by  $M_A$ , which is, in fact, the space of all non-zero complex (continuous) homomorphisms on A. Clearly, for each  $x \in X$  the map  $e_x : A \longrightarrow \mathbb{C}$ , defined by  $e_x(f) = f(x)$  is a homomorphism (character) on A which is called the evaluation homomorphism (character) on A at x. It is clear that A is semisimple, since

$$\operatorname{rad}(A) = \{ f \in A : f \in \ker(\phi), \text{ for all } \phi \text{ in } M_A \}$$

$$\subseteq \{ f \in A : f \in \ker(e_x), \text{ for all } x \text{ in } X \}$$

$$= \{ f \in A : f(x) = 0, \text{ for all } x \text{ in } X \}$$

$$= \{ 0 \}.$$

The map  $J: X \longrightarrow M_A$ , defined by  $J(x) = e_x$ , is continuous and injective.

**Definition 2.2** The Banach function algebra A is called *natural* if the map J is surjective, that is, every homomorphism (character) on A is an evaluation homomorphism (character) at some  $x \in X$ .

For example, C(X) is a natural complex uniform algebra on X [4; Sect.17] or [26; Example 11.13]. If  $(A, \|\cdot\|)$  is a complex Banach function

algebra on X, the uniform norm of each element  $f \in A$ , does not exceed from its norm, since

$$||f||_{X} = \sup_{x \in X} |f(x)| \le \sup_{\phi \in M_{A}} |\phi(f)|$$

$$= \sup_{\phi \in M_{A}} |\hat{f}(\phi)| = ||\hat{f}||_{M_{A}} = \rho_{A}(f) = \inf_{n \ge 1} ||f^{n}||^{\frac{1}{n}} \le ||f||.$$

**Definition 2.3** Let (X,d) be a compact metric space and  $\alpha > 0$ . The algebra of all complex-valued functions f on X for which

$$p_{\alpha}(f) = \sup_{\substack{x,y \in X \\ x \neq y}} \frac{|f(x) - f(y)|}{d^{\alpha}(x,y)} < \infty ,$$

is denoted by  $Lip(X,\alpha)$  and the subalgebra of those functions f for which  $\lim_{d(x,y)\to 0} \frac{|f(x)-f(y)|}{d^{\alpha}(x,y)} = 0$ , is denoted by  $\ell ip(X,\alpha)$ . These are called Lipschitz algebras of order  $\alpha$  and were first studied by Sherbert [30].

Clearly  $Lip(X,\alpha)$  is a Banach algebra of continuous complex-valued functions on X under the norm  $||f||_{\alpha} = ||f||_{X} + p_{\alpha}(f)$ , where  $||f||_{X} = \sup_{x \in X} |f(x)|$ . If  $lip(X,\alpha)$  is also equipped with the above norm then it is easily proved that  $lip(X,\alpha)$  is a closed subalgebra of  $Lip(X,\alpha)$  and hence it is also a Banach algebra.

Lipschitz algebras contain the constants for every positive  $\alpha$ . But when X is a connected compact subset of  $\mathbb{C}^n$ ,  $d(x,y) = \|x-y\|$  and  $1 < \alpha \ (\alpha \geq 1)$ , then  $Lip(X,\alpha)$  ( $\ell ip(X,\alpha)$ ) may contain the constant functions only. In general, when  $1 < \alpha$ ,  $Lip(X,\alpha)$  may contain the nonconstant functions. for example, if d is the discrete metric on X then for a fixed  $a \in X$ , the function, f(x) = d(x,a) is an element of  $Lip(X,\alpha)$  which is a nonconstant function.

In most parts of this paper X is usually a compact subset of  $\mathbb{C}^n$ , so we shall concern ourselves with the algebras  $Lip(X,\alpha)$  (respectively,  $\ell ip(X,\alpha)$ ) only when  $0<\alpha\leq 1$  (respectively,  $0<\alpha<1$ ). In this case the algebras contain the family of functions  $f(x)=d(x,\alpha)$ , as a runs

over X, which separates the points of X. Hence  $Lip(X,\alpha)$  for  $\alpha \leq 1$  and  $\ell ip(X,\alpha)$  for  $\alpha < 1$ , are Banach function algebras on the compact metric space X. Since these Lipschitz algebras are self-adjoint and separate the points of X, they are uniformly dense in C(X), by the Stone-Weierstrass Theorem. Hence they are natural Banach function algebras on X, that is, their maximal ideal spaces coincide with X. For the proof one can either follow the same process as the proof of the naturality of C(X), by using the fact that these algebras are inverse-closed, or apply the main theorem in [11].

Note that the function  $f(x) = d^{\alpha}(x,a)$  is an element of  $Lip(X,\alpha)$  but  $\lim_{x \to a} \frac{f(x)}{d^{\alpha}(x,a)} = 1$ . Hence f does not belong to  $\ell ip(X,\alpha)$  when X is infinite and so  $\ell ip(X,\alpha)$  is a proper subalgebra of  $Lip(X,\alpha)$  in this case. Moreover, for every  $\alpha < 1$  we have the inclusion  $Lip(X,1) \subseteq \ell ip(X,\alpha)$ . In fact, Lip(X,1) is dense in  $\ell ip(X,\alpha)$ . This result has been proved in [3], using the measure theory and duality, but it is also followed from a theorem due to Hedberg in [10], which is, in fact, a Stone-Weierstrass theorem for Lipschitz algebras of real-valued functions. However, this theorem does hold even for Lipschitz algebras of complex-valued functions with an extra condition as follows, which we need later for some approximation theorems.

Theorem 2.4 (Hedberg's Theorem) Let (X,d) be a compact metric space, and  $0 < \alpha < 1$ . Let A be a self-adjoint subalgebra of  $\ell ip(X,\alpha)$  which separates the points of X and contains the constant functions. Then A is dense in  $\ell ip(X,\alpha)$  if for every  $a \in X$  there are numbers  $M_a$  and  $\delta_a$  such that for every  $\delta \leq \delta_a$  there is an  $f \in A$  with f(a) = 1, f(x) = 0 on  $S_a(\delta) = \{x \in X : d(x,a) = \delta\}$ , and  $\sup_{y,z \in B_a(\delta)} \frac{|f(y) - f(z)|}{d^{\alpha}(y,z)} < \frac{M_a}{\delta^{\alpha}}$ , where  $B_a(\delta) = \{x \in X : d(x,a) \leq \delta\}$ .

For the proof one can set  $A_1 = \{ f \in A : f \text{ is real-valued} \}$  and see that  $A_1$  is a subalgebra of  $\text{Relip}(X, \alpha)$ , the real Lipschitz algebra, which has

all properties of Hedberg's theorem in the real case. Hence  $A_1$  is dense in  $Relip(X,\alpha)$ . Now by considering the fact that A is self-adjoint and the real and imaginary parts of Lipschitz functions are also Lipschitz functions, it is easy to show that qA is dense in  $lip(X,\alpha)$ .

Corollary 2.5 Since A = Lip(X,1) is a self-adjoint subalgebra of  $lip(X,\alpha)$  and for every  $a \in X$  and arbitrary positive  $\delta$  the function g(x) = 1-  $\frac{d(x,a)}{\delta}$  is an element of Lip(X,1) which satisfies the conditions of Hedberg's theorem, it follows that Lip(X,1) is dense in  $lip(X,\alpha)$  for every  $\alpha < 1$ .

Now we discuss the approximation problem in  $Lip = Lip(X, \alpha)$  for  $\alpha \leq 1$  and in  $\ell ip = \ell ip(X, \alpha)$  for  $\alpha < 1$ , when X is a compact subset of  $\mathbb{C}^n$ . Since the polynomials and the rational functions with poles off X belong to Lip(X,1), we can define the following subalgebras of  $Lip(X,\alpha)$  and  $\ell ip(X,\alpha)$ . From now on we always consider  $Lip(X,\alpha)$  for  $\alpha \leq 1$  and  $\ell ip(X,\alpha)$  for  $\alpha < 1$ .

Before presenting the next definition we prove the following useful result.

**Theorem 2.6** Let A be a natural Banach function algebra on a compact subset X of  $\mathbb{C}^n$ , and suppose A contains the polynomials. Then every function which is analytic in a neighbourhood of X is an element of A.

**proof.** Since  $M_A \cong X$ , i.e. the maximal ideal space of A is homeomorphic with X, the joint spectrum of the coordinate function is precisely the set X. If f is a function analytic in a neighbourhood of X, then by the Functional Calculus Theorem [8; III, 4.5], there exists  $g \in A$  such that  $\hat{g} = f(\hat{z}_1, \hat{z}_2, \ldots, \hat{z}_n)$  on  $M_A \cong X$ , where  $\hat{z}_k$  is the Gelfand transform of the coordinate function  $z_k$   $(1 \le k \le n)$ . Therefore g = f on X and hence  $f \in A$ .

By the above theorem the Lipschitz algebras  $Lip(X, \alpha)$  and  $\ell ip(X, \alpha)$ 

contain any function which is analytic in some neighbourhood of X.

**Definition 2.7** The subalgebra of Lip ( $\ell ip$ ) which is generated by the polynomials in  $z_1, z_2, \ldots, z_n$ , by the rational functions with poles off X, or by the analytic functions in some neighbourhood of X, is denoted by  $Lip_P$  ( $\ell ip_P$ ), by  $Lip_R$  ( $\ell ip_R$ ), or by  $Lip_H$  ( $\ell ip_H$ ), respectively.

**Definition 2.8** The subalgebra of Lip ( $\ell ip$ ) which is generated by those elements of Lip ( $\ell ip$ ) which are analytic in the interior of X, is denoted by  $Lip_A$  ( $\ell ip_A$ ).

These subalgebras are all Banach function algebras on X and they satisfy the following inclusions:

$$Lip_P \subseteq Lip_R \subseteq Lip_H \subseteq Lip_A \subseteq Lip$$
,

$$\ell ip_P \subseteq \ell ip_R \subseteq \ell ip_H \subseteq \ell ip_A \subseteq \ell ip$$
.

Note that continuous functions on X, which are analytic in the interior of X, may not be in  $Lip(X,\alpha)$ . Since for every  $f \in Lip(X,\alpha)$ ,  $||f||_X \leq ||f||_{\alpha}$ , it is easy to see that  $Lip_A(X,\alpha) = A(X) \cap Lip(X,\alpha)$  and  $\ell ip_A(X,\alpha) = A(X) \cap \ell ip(X,\alpha)$ , where A(X) is the algebra of continuous functions on X which are analytic in the interior of X. Hence  $Lip_A(X,\alpha) = Lip(X,\alpha)$  or  $\ell ip_A(X,\alpha) = \ell ip(X,\alpha)$ , if and only if, X has empty interior.

Now we investigate the maximal ideal spaces of these subalgebras, and the equality among some of them. For this purpose, we consider the standard uniform algebras P(X), R(X) and H(X) which are the uniform closures of polynomials, rational functions with poles off X and functions which are analytic in a neighbourhood of X, respectively. For the subalgebra B of C(X), we denote its maximal ideal space by  $M_B$  and its uniform closure by  $\overline{B}$ . Clearly  $\overline{Lip_P} = \overline{\ell ip_P} = P(X)$ ,  $\overline{Lip_R} = \overline{\ell ip_R} = R(X)$ , and  $\overline{Lip_H} = \overline{\ell ip_H} = H(X)$ .

Also for every  $f \in Lip$  we have  $||f^n||_{\alpha} \leq ||f||_X^n + np_{\alpha}(f)||f||_X^{n-1}$ , and so  $||\hat{f}|| \leq ||f||_X$ , where  $\hat{f}$  is the Gelfand transform of f. Thus as a consequence of the Theorem in [11], we obtain the following results:

#### Theorem 2.9

- I.  $M_{Lip_P} = M_{tip_P} \cong \hat{X}$ , where  $\hat{X}$  is the polynomial convex hull of X.
- II.  $M_{Lip_R} = M_{lip_R} \cong R\text{-}hull(X)$ , where R-hull(X) is the rational convex hull of X.
- III.  $M_{Lip_H} = M_{\ell ip_H} = M_{H(X)}$ .
- IV.  $\hat{X} = R$ -hull (X) if and only if  $Lip_P = Lip_R$ .
- V. If R-hull (X) = X then  $Lip_R = Lip_H$ .

Note that whenever A(X) = H(X) we have  $\overline{Lip_A}(X,\alpha) = \overline{lip_A}(X,\alpha)$  = A(X) and so  $M_{Lip_A} = M_{lip_A} = M_{H(X)}$ . But in general these equalities may not be satisfied. However, if X is an arbitrary compact plane set then  $lip_A(X,\alpha)$  and  $Lip_A(X,\alpha)$  are natural Banach function algebras on X [20,16, or1], but when X is an arbitrary compact subset of  $\mathbb{C}^n$  (1 < n) the maximal ideal spaces of  $Lip_A$  and  $lip_A$  are not known yet. For further results on some other subalgebras of  $lip(X,\alpha)$  one can refer to [31].

To establish some results about the approximation problem we need to define the algebra of continuously differentiable functions.

**Definition 2.10** Let X be a perfect compact plane set, a complex-valued function f on X is called differentiable at  $z_0 \in X$ , if

$$f'(z_0) = \lim_{\substack{z \in X \\ z \to z_0}} \frac{f(z) - f(z_0)}{z - z_0}$$

exists, and f is called differentiable on X if it is differentiable at each point of X. The complex algebra of complex-valued functions with derivatives of all orders on X is denoted by  $D^{\infty}(X)$ . Let  $D^{n}(X)$  be the algebra of functions with continuous  $n^{th}$  derivatives on X. For  $f \in D^{n}(X)$ , we define the norm by

$$p_n(f) = || f || = \sum_{k=0}^n \frac{|| f^{(k)} ||_X}{k!},$$

This norm is actually an algebra norm on  $D^n(X)$ . It is interesting to see that  $D^{\infty}(X) = \bigcap_{n=1}^{\infty} D^n(X)$ .

Definition 2.11 Let  $M = \{M_k\}_{k=0}^{\infty}$  be a sequence of positive numbers such that

$$M_0 = 1$$
 and  $\frac{M_k}{M_r.M_{k-r}} \geq \binom{k}{r}$   $r = 0, 1, \ldots, k$ .

The algebra of infinitely differentiable functions f on X such that

$$\sum_{k=0}^{\infty} \frac{\parallel f^{(k)} \parallel_X}{M_k} < \infty$$

is denoted by D(X,M). For convenience, we regard  $D^n(X)$  as being an algebra of the type D(X,M) by setting  $M_r = r!$   $(r = 0,1,\ldots,n)$  and  $\frac{1}{M_r} = 0$   $(r = n + 1, n + 2,\ldots)$ .

From now on whenever we refer to  $M = \{M_k\}$  we mean this sequence satisfies the above conditions.

Now we introduce the type of compact sets which we shall consider next.

**Definition 2.12** Let X be a compact plane set which is connected by rectifiable arcs, and suppose  $\delta(z, w)$  is the geodesic metric on X, the infimum of the lengths of the arcs joining z and w.

- (i) X is called regular if for each  $z_0 \in X$  there exists a constant C such that for all  $z \in X$ ,  $\delta(z, z_0) \leq C|z z_0|$ .
- (ii) X is called uniformly regular if there exists a constant C such that for all  $z, w \in X$ ,  $\delta(z, w) \leq C|z w|$ .

If X is a finite union of regular sets then for each  $z_0 \in X$  there exists a constant C such that for every  $z \in X$  and any  $f \in D^1(X)$ ,

$$| f(z) - f(z_0) | \le C |z - z_0| (||f||_X + ||f'||_X).$$

This inequality implies that  $D^1(X)$  is complete under the norm  $||f||_1 = ||f||_X + ||f'||_X$  [6,14]. It is also interesting to note that the above condition is, in fact, a necessary and sufficient condition for the completeness of  $D^1(X)$ . To see this, let  $D^1(X)$  be complete and define another norm on  $D^1(X)$  by

$$||f|| = ||f||_X + ||f'|| + \sup_{\substack{z \in X \\ z \neq z_0}} \frac{|f(z) - f(z_0)|}{|z - z_0|} \qquad (f \in D^1(X)),$$

where  $z_0$  is a fixed point in X. Then  $D^1(X)$  is also a Banach function algebra on X under this new norm. Thus there exists a constant C such that for all  $f \in D^1(X)$  and for every  $z \in X$ 

$$| f(z) - f(z_0) | \le C |z - z_0| (||f||_X + ||f'||_X).$$

Note that the completeness of  $D^1(X)$  is, in fact, equivalent to the completeness of D(X,M). Moreover,  $D^1(X) \subseteq Lip(X,1)$ , and the norm  $\|.\|_1$  of  $D^1(X)$  and the Lipschitz norm of Lip(X,1) are equivalent on  $D^1(X)$ . Note that  $D^1(X)$  is a proper closed subalgebra of Lip(X,1) for every uniformly regular set X.

From now on we assume that X is a perfect compact plane set such that  $D^1(X)$  is complete, unless otherwise specified.

As mentioned before for each  $\alpha < 1$ , Lip(X, 1) is dense in  $lip(X, \alpha)$ . A question which arises here is that under what conditions  $D^1(X)$  is also dense in  $\ell ip(X,\alpha)$ . If  $D^1(X)$  is dense in  $\ell ip(X,\alpha)$  then  $\ell ip_A(X,\alpha) = \ell ip(X,\alpha)$  and so  $intX = \phi$ . Therefore, a necessary condition for the density of  $D^1(X)$  in  $\ell ip(X,\alpha)$  is that  $intX = \phi$ . But the inverse is not true. However, we shall prove the interesting result that  $D^1(X)$  is actually dense in  $\ell ip_R(X,\alpha)$ .

For this purpose and the next general result we make some preliminaries. Let X be a compact subset of the complex plane and define  $V = \{(z, w) \in X \times X : z \neq w\}$  and  $W = X \cup V$ . We denote the Banach space of bounded continuous functions on W by  $C^b(W)$ , the closed subspace of functions which vanish at infinity by  $C_0(W)$ , and the space of all regular complex Borel measures on W by M(W). It follows from the Riesz representation theorem that the dual space of  $C_0(W)$  is isometrically isomorphic to M(W). As noted in [3, §3] we extend any function  $g \in C(X)$  to a function  $\tilde{g}$  on W by defining

$$\left\{ \begin{array}{ll} \tilde{g}(w) = g(w) & (w \in X), \\ \\ \tilde{g}(\zeta, \eta) = \frac{g(\zeta) - g(\eta)}{|\zeta - \eta|^{\alpha}} & ((\zeta, \eta) \in V). \end{array} \right.$$

The map  $g \mapsto \tilde{g}$ ,  $Lip(X, \alpha) \to C^b(W)$  is a linear isometry and the image of  $\ell ip(X, \alpha)$  is contained in  $C_0(W)$ .

Theorem 2.13 [13] Let X be a compact subset of the complex plane and suppose  $f: X \to \mathbb{C}$  can be extended to a function having continuous partial derivatives in some neighbourhood of X. If either of the following conditions is satisfied, then  $f \in lip_R(X, \alpha)$  for every  $\alpha < 1$ .

(i) m(X) = 0, where m is planar measure.

(ii) 
$$\frac{\partial f}{\partial \bar{z}} = 0$$
 on  $X$ .

**Proof.** Suppose f satisfies the hypotheses of the theorem but  $f \notin lip_R(X, \alpha)$ . By the Hahn-Banach theorem there exists a

continuous linear functional  $\phi$  on  $lip(X,\alpha)$  for which  $\phi(f) \neq 0$  and  $\phi = 0$  on  $lip_R(X,\alpha)$ . Again, by the Hahn-Banach theorem,  $\phi$  has a norm preserving extension to a continuous linear functional  $\psi$  on  $C_0(W)$ , and so there exists a  $\mu \in M(W)$  with  $\phi(g) = \psi(\tilde{g}) = \int_W \tilde{g} d\mu$ , for every  $g \in lip(X,\alpha)$ .

By hypotheses there exists a function F defined and with continuous partial derivatives on  $\mathbb{R}^2$ , having compact support, such that  $F|_X = f$ . By Green's formula

$$F(w) = -\frac{1}{\pi} \iint_{\mathbb{R}^2} \frac{\partial F}{\partial \bar{z}} \cdot \frac{1}{z - w} dx dy \qquad (w \in \mathbb{C}) .$$

So we have

$$\phi(f) = \int_{W} \tilde{f} d\mu = \int_{X} f(w) d\mu(w) + \int_{V} \frac{f(\zeta) - f(\eta)}{|\zeta - \eta|^{\alpha}} d\mu(\zeta, \eta)$$

$$= \int_{X} \left( \frac{-1}{\pi} \iint_{\mathbb{R}^{2}} \frac{\partial F}{\partial \bar{z}} \cdot \frac{1}{z - w} dx dy \right) d\mu(w)$$

$$+ \int_{V} \left( \frac{-1}{\pi} \iint_{\mathbb{R}^{2}} \frac{\partial F}{\partial \bar{z}} \cdot \frac{\frac{1}{z - \zeta} - \frac{1}{z - \eta}}{|\zeta - \eta|^{\alpha}} dx dy \right) d\mu(\zeta, \eta).$$

Since F has a compact support in  $\mathbb{R}^2$ , there exists a closed disk  $\Delta$  in  $\mathbb{C}$  such that  $\frac{\partial F}{\partial z} = 0$  on  $\mathbb{C}\backslash\Delta$ . To show the boundedness of the integral

$$\iint_{\Delta} \left| \frac{\partial F}{\partial \bar{z}} \right| \frac{|\zeta - \eta|^{1 - \alpha}}{|z - \zeta|.|z - \eta|} dx dy$$

over V, we split  $\Delta$  into three parts,  $D_{\zeta}, D_{\eta}$  and  $\Delta \setminus (D_{\zeta} \cup D_{\eta})$ , where  $D_{\zeta}$  and  $D_{\eta}$  are disks with centres  $\zeta$  and  $\eta$ , respectively, having radius  $\frac{|\zeta - \eta|}{2}$ . Clearly the integrals on  $D_{\zeta}$  and  $D_{\eta}$  are bounded. If  $z \in \Delta \setminus (D_{\zeta} \cup D_{\eta})$  then  $|\zeta - \eta| \leq 2|z - \zeta|$  and  $|\zeta - \eta| \leq 2|z - \eta|$  and so  $|\zeta - \eta|^{1-\alpha} \leq 2^{1-\alpha}|z - \zeta|^{\frac{1-\alpha}{2}}.|z - \eta|^{\frac{1-\alpha}{2}}$ . Hence

$$\iint_{\Delta \setminus (D_{\zeta} \cup D_{\eta})} \left| \frac{\partial F}{\partial \bar{z}} \right| \frac{|\zeta - \eta|^{1 - \alpha}}{|z - \zeta||z - \eta|} dx dy \le 2^{1 - \alpha} \iint_{\Delta \setminus (D_{\zeta} \cup D_{\eta})} \left| \frac{\partial F}{\partial \bar{z}} \right| \frac{dx dy}{(|z - \zeta|, |z - \eta|)^{\frac{1 + \alpha}{2}}}.$$

Now, using Hölder's inequality the last integral is bounded over V since  $\alpha < 1$ . Therefore by Fubini's theorem and hypotheses we have

$$\phi(f) = -\frac{1}{\pi} \iint_{\Delta} \frac{\partial F}{\partial \bar{z}} \left( \int_{X} \frac{d\mu(w)}{z - w} + \int_{V} \frac{\frac{1}{z - \zeta} - \frac{1}{z - \eta}}{|\zeta - \eta|^{\alpha}} d\mu(\zeta, \eta) \right) dx dy$$
$$= -\frac{1}{\pi} \iint_{\Delta \setminus X} \frac{\partial F}{\partial \bar{z}} \phi\left(\frac{1}{z - w}\right) dx dy = 0,$$

which contradicts our previous assumption. Therefore  $f \in lip_R(X, \alpha)$  and this completes the proof of the theorem.

As an interesting result on the approximation problem we conclude an extension of the Hartogs-Rosenthal Theorem, which states that R(X) = C(X), when X has planar measure zero.

**Theorem 2.14** [13] If X is a compact subset of the complex plane with planar measure zero, then  $lip_R(X,\alpha) = lip(X,\alpha)$ .

**Proof.** By Theorem 2.13,  $\bar{z} \in lip_R(X,\alpha)$ , where  $\bar{z}$  is the complex conjugate of the coordinate function z. For every  $a \in X$  and  $\delta > 0$  the function  $f(z) = 1 - \frac{|z-a|^2}{\delta^2}$  is an element of  $lip_R(X,\alpha)$ . Hence  $A = lip_R(X,\alpha)$  satisfies the conditions of Hedberg's theorem, and so A is dense in  $lip(X,\alpha)$ , that is,  $lip_R(X,\alpha) = lip(X,\alpha)$ .

Remark 2.15 The above theorem is not true for the algebra  $Lip(X, \alpha)$ , when  $\alpha < 1$  and X is a compact subset of  $\mathbb{C}^n$ . Note that  $Lip_R(X, 1)$  is also a proper closed subalgebra of Lip(X, 1) for every uniformly regular set X.

For similar results, which are related to the second part of Theorem 2.13 and Theorem 2.14, one can refer to [23, 24, 25].

Now we extend a result due to Dales and Davie which states that  $D^1(X) \subseteq R(X)$  if X is uniformly regular [6, Lemma 1.5].

**Theorem 2.16** [13] For every perfect compact plane set X,  $\overline{D^1}(X) = \ell i p_R(X, \alpha)$ .

**Proof.** It is sufficient to prove that  $D^1(X) \subseteq lip_R(X,\alpha)$ . Let  $f \in D^1(X)$ . By the Whitney's extension theorem [21], there exists a function F, defined and with continuous partial derivatives on  $\mathbb{R}^2$ , having a compact support, such that  $F|_X = f$ ,  $\frac{\partial F}{\partial \bar{z}}|_X = 0$  and  $\frac{\partial F}{\partial z}|_X = f'$ . Hence f satisfies the hypotheses of Theorem 2.13, and so  $f \in lip_R(X,\alpha)$ .

Remark 2.17 [13] If  $lip_R(X,\alpha) = lip(X,\alpha)$  then R(X) = C(X). But there exists a Swiss cheese  $X_0$ , with empty interior and positive planar measure, such that  $R(X_0)$  is different from  $C(X_0)$ , [8; II,1]. Hence  $lip_R(X_0,\alpha)$  is different from  $lip(X_0,\alpha)$ . Since every Swiss cheese is a perfect compact plane set we conclude that  $D^1(X_0)$  is not dense in  $lip(X_0,\alpha)$ .

Finally we extend the result of Theorem 2.16 to obtain an extension of a similar result in uniform algebras for Lipschitz algebras.

**Theorem 2.18** [13] Let X be a compact subset of  $\mathbb{C}^n$ , and let  $X_j$   $(1 \le j \le n)$  denote the projection of X onto the jth coordinate plane.

- (i) If  $\ell i p_P(X_j, \alpha) = \ell i p(X_j, \alpha)$  for all j  $(1 \leq j \leq n)$ , then  $\ell i p_P(X, \alpha) = \ell i p(X, \alpha)$ .
- (ii) If  $\ell ip_R(X_j, \alpha) = \ell ip(X_j, \alpha)$  for all j  $(1 \leq j \leq n)$ , then  $\ell ip_R(X, \alpha) = \ell ip(X, \alpha)$ .
- **Proof.** (i) Clearly the complex conjugate of every polynomial in  $z_1, z_2, \ldots, z_n$  is an element of  $\ell i p_P(X, \alpha)$ . Hence the function  $f(z) = 1 \frac{\|z-a\|^2}{\delta^2}$  is an element of  $\ell i p_P(X, \alpha)$  for every  $a \in X$  and  $\delta > 0$ . Therefore by the Hedberg's theorem,  $\ell i p_P(X, \alpha)$  is dense in  $\ell i p(X, \alpha)$ , in other words,  $\ell i p_P(X, \alpha) = \ell i p(X, \alpha)$ .
- (ii) Note that every rational function r is of the form  $r(z) = \frac{p(z)}{q(z)}$ , where p(z) and q(z) are polynomials in  $z_1, z_2, \ldots, z_n$  and  $q(z) \neq 0$  on R-hull (X), which is the maximal ideal space of  $lip_R(X, \alpha)$ . Hence  $\bar{q}(z)$ , the complex conjugate of q(z), as an element of  $lip_R(X, \alpha)$ , is invertible in  $lip_R(X, \alpha)$  and so  $\bar{r}(z)$  is an element of  $lip_R(X, \alpha)$ . The result now

follows as in (i).

As a consequence we get the following.

Corollary 2.19 If for each j  $(1 \le j \le n)$ ,  $X_j$  has planar measure zero, then  $\ell ip_R(X,\alpha) = \ell ip(X,\alpha)$ . If moreover  $\hat{X}_j = X_j$  for all j  $(1 \le j \le n)$  then  $\ell ip_P(X,\alpha) = \ell ip(X,\alpha)$ .

Now we are going to discuss the approximation problem for the Lipschitz algebras of differentiable functions on perfect compact plane sets. For further details one can refer to [19].

**Definition 2.20** The algebra of functions f on X whose derivatives up to order n exist and for each k  $(0 \le k \le n)$ ,  $f^{(k)} \in Lip(X,\alpha)$  is denoted by  $Lip^n(X,\alpha)$ . The algebra  $\ell ip^n(X,\alpha)$  is defined in a similar way.

We now equip these algebras with the norm

$$p_n(f) = || f || = \sum_{k=0}^n \frac{|| f^{(k)} ||_{\alpha}}{k!} = \sum_{k=0}^n \frac{|| f^{(k)} ||_X + p_{\alpha}(f^{(k)})}{k!}.$$

The above algebras have similar properties to  $D^n(X)$ . Clearly for each  $n \geq 1$ ,  $Lip^n(X,1) \subseteq \ell ip^n(X,\alpha) \subseteq Lip^n(X,\alpha) \subseteq D^n(X) \subseteq D^1(X)$ . It is also known that  $D^1(X) \subseteq R(X)$  [6].

**Definition 2.21** The algebra of functions f with derivatives of all orders for which  $f^{(k)} \in Lip(X,\alpha)$  ( $f^{(k)} \in lip(X,\alpha)$ ) for all k is denoted by  $Lip^{\infty}(X,\alpha)$  ( $lip^{\infty}(X,\alpha)$ ).

It is interesting to see that  $D^{\infty}(X) = \bigcap_{n=1}^{\infty} D^{n}(X)$ ,  $Lip^{\infty}(X,\alpha) = \bigcap_{n=1}^{\infty} Lip^{n}(X,\alpha)$  and  $\ell ip^{\infty}(X,\alpha) = \bigcap_{n=1}^{\infty} \ell ip^{n}(X,\alpha)$ . Note that these algebras are not Banach algebras under any norm [7, 12, or 22].

We now introduce certain subalgebras of  $Lip^{\infty}(X,\alpha)$  and  $\ell ip^{\infty}(X,\alpha)$ .

#### Definition 2.22 Let

$$Lip(X,M,\alpha) = \{ f \in Lip^{\infty}(X,\alpha) : \sum_{k=0}^{\infty} \frac{\|f^{(k)}\|_{\alpha}}{M_k} < \infty \},$$

$$\ell ip(X,M,\alpha) = \{ f \in \ell ip^{\infty}(X,\alpha) : \sum_{k=0}^{\infty} \frac{\|f^{(k)}\|_{\alpha}}{M_k} < \infty \},$$

and for f in  $Lip(X, M, \alpha)$  or in  $\ell ip(X, M, \alpha)$  let  $||f|| = \sum_{k=0}^{\infty} \frac{||f^{(k)}||_{\alpha}}{M_k}$ 

Remark 2.23 The above algebras have similar properties to D(X,M). For convenience, we regard  $Lip^n(X,\alpha)$  and  $\ell ip^n(X,\alpha)$  as being algebras of the type  $Lip(X,M,\alpha)$  and  $\ell ip(X,M,\alpha)$ , respectively, by setting  $M_k = k!$   $(k = 0,1,\ldots,n)$  and  $1/M_k = 0$   $(k = n+1,\ldots)$ . Clearly for  $\alpha < 1$ ,  $Lip(X,M,1) \subseteq \ell ip(X,M,\alpha) \subseteq Lip(X,M,\alpha) \subseteq D(X,M)$ ,  $\ell ip(X,M,\alpha) \subseteq \ell ip^{n+1}(X,\alpha) \subseteq \ell ip^n(X,\alpha)$ ,  $Lip(X,M,\alpha) \subseteq Lip^{n+1}(X,\alpha) \subseteq Lip^n(X,\alpha)$ . Since  $Lip(X,\alpha)$  and  $\ell ip(X,\alpha)$  are normed function algebras and  $\|.\|_{\alpha}$  is an algebra norm on them, for every  $f,g \in Lip(X,M,\alpha)$   $(f,g \in \ell ip(X,M,\alpha))$  we have

$$\begin{aligned} &||f.g|| = \sum_{k=0}^{\infty} \frac{||(f.g)^{(k)}||_{\alpha}}{M_{k}} \leq \sum_{k=0}^{\infty} \frac{1}{M_{k}} \sum_{r=0}^{k} {k \choose r} ||f^{(r)}.g^{(k-r)}||_{\alpha} \\ &\leq \sum_{k=0}^{\infty} \sum_{r=0}^{k} \frac{||f^{(r)}||_{\alpha}}{M_{r}} \cdot \frac{||g^{(k-r)}||_{\alpha}}{M_{k-r}} = \left(\sum_{k=0}^{\infty} \frac{||f^{(k)}||_{\alpha}}{M_{k}}\right) \left(\sum_{k=0}^{\infty} \frac{||g^{(k)}||_{\alpha}}{M_{k}}\right) = ||f||.||g||. \end{aligned}$$

Hence  $f,g \in Lip(X,M,\alpha)$  ( $f,g \in lip(X,M,\alpha)$ ) and so any of the above algebras are normed function algebras on X, with respect to the above norm. The completeness of  $D^1(X)$  or, equivalently, D(X,M), implies that  $Lip(X,M,\alpha)$  and  $lip(X,M,\alpha)$  are Banach function algebras on X.

Remark 2.24 When X is a uniformly regular set we have  $D^{n+1}(X) \subseteq Lip^n(X,1) \subseteq lip^n(X,\alpha)$  and for  $\alpha < 1$   $lip(X,M,\alpha) = Lip(X,M,\alpha)$  if  $\frac{1}{M_k}$  is different from zero for infinitely many k.

Remark 2.25 When  $P_k = \sqrt[k]{\frac{M_k}{k!}} \to \infty$  as  $k \to \infty$ , D(X, M) and likewise Lip(X, M, 1) contain all rational functions with poles off X.

From now on we assume that X is a perfect compact plane set such that  $D^1(X)$  is complete, unless otherwise specified.

Now we introduce subalgebras of  $Lip(X, M, \alpha)$  and  $\ell ip(X, M, \alpha)$ .

**Definition 2.26** The closed subalgebra of  $Lip(X, M, \alpha)$  ( $lip(X, M, \alpha)$ ) which is generated by the polynomials, by the rational functions with poles off X that belong to  $Lip(X, M, \alpha)$  ( $lip(X, M, \alpha)$ ), or by those functions of  $Lip(X, M, \alpha)$  ( $lip(X, M, \alpha)$ ) which are analytic in some neighbourhood of X, is denoted by  $Lip_P(X, M, \alpha)$  ( $lip_P(X, M, \alpha)$ ),  $Lip_R(X, M, \alpha)$  ( $lip_R(X, M, \alpha)$ ) or  $Lip_H(X, M, \alpha)$  ( $lip_H(X, M, \alpha)$ ), respectively.

Remark 2.27 Clearly  $Lip_P(X,M,\alpha)$  is uniformly dense in P(X), the uniform closure of polynomials. Moreover, when  $P_k = \sqrt[k]{M_k/k!} \to \infty$  as  $k \to \infty$ , we have  $R_0(X) \subseteq Lip(X,M,1)$ , so Lip(X,M,1) and hence  $lip(X,M,\alpha)$  and  $Lip(X,M,\alpha)$  are uniformly dense in R(X). In particular,  $Lip_R(X,M,\alpha)$  is uniformly dense in R(X). Note that, when  $P_k \to \infty$  as  $k \to \infty$ ,  $Lip_R(X,M,\alpha)$  is generated by the all rational functions with poles off X, and hence it is a natural Banach function algebra on X. So by the Theorem in [11] we have  $M_{Lip_P(X,M,\alpha)} \cong M_{P(X)} \cong \hat{X}$ , where  $M_A$  is the maximal ideal space of the algebra A. Thus when  $P_k \to \infty$  as  $k \to \infty$ ,  $Lip_P(X,M,\alpha) = Lip_R(X,M,\alpha)$  if and only if  $\hat{X} = X$ . Also when  $P_k \to \infty$  as  $k \to \infty$ , by the Functional Calculus Theorem [8; 3.4.5],  $Lip_R(X,M,\alpha)$  contains all analytic functions in a neighbourhood of X, and so  $Lip_R(X,M,\alpha) = Lip_H(X,M,\alpha)$ .

**Theorem 2.28** [14] For each  $n \geq 0$ ,  $\ell ip^n(X, \alpha)$  and  $Lip^n(X, \alpha)$  are natural Banach function algebras on X.

**Proof.** Since the algebras  $lip(X,\alpha)$  and  $Lip(X,\alpha)$  are uniformly dense in C(X) and for  $n \geq 1$ ,  $lip^n(X,\alpha)$  and  $Lip^n(X,\alpha)$  are uniformly dense in R(X), by the naturality of C(X) and R(X), we can show that

for each  $f \in Lip^n(X, \alpha)$  we have  $||\hat{f}|| \le ||f||_X$ , where  $\hat{f}$  is the Gelfand transform of f. Thus by the Theorem in [11] the result follows.

It is clear that for  $\alpha < 1$   $\ell ip_P(X, M, \alpha) = Lip_P(X, M, \alpha)$ ,  $\ell ip_R(X, M, \alpha) = Lip_R(X, M, \alpha)$  and  $\ell ip_H(X, M, \alpha) = Lip_H(X, M, \alpha)$ .

**Theorem 2.29** If X is uniformly regular and  $\ell ip_P(X,\alpha) = \ell ip_A(X,\alpha)$  then  $\ell ip_P^n(X,\alpha) = \ell ip^n(X,\alpha)$  for all  $n \geq 1$  and  $\alpha < 1$ .

**Proof.** As we know if  $f \in D^1(X)$  then  $p_{\alpha}(f) \leq Cd^{1-\alpha}||f'||_X$ , where d = diam(X). Now let  $n \geq 1$  and  $f \in lip^n(X, \alpha)$ . Since  $f^{(n)} \in lip_A(X, \alpha) = lip_P(X, \alpha)$ , for every  $\epsilon > 0$  there exists a polynomial  $P_0$  such that

$$||f^{(n)} - P_0||_{\ell i p(X,\alpha)} = ||f^{(n)} - P_0||_X + p_\alpha(f^{(n)} - P_0) < \epsilon.$$

Let  $z_0$  be a fixed point in X and  $P_1$  be the antiderivative of  $P_0$  with the initial condition  $P_1(z_0) = f^{(n-1)}(z_0)$ .

Since  $f^{(n-1)} - P_1 \in lip^1(X, \alpha) \subseteq D^1(X)$  we have

$$p_{\alpha}(f^{(n-1)} - P_1) \le Cd^{1-\alpha}||f^{(n)} - P_0||_X < Cd^{1-\alpha}\epsilon.$$

Continuing in this manner, we obtain polynomials  $P_2, P_3, \ldots, P_n$  such that  $P'_k = P_{k-1}, P_k(z_0) = f^{(n-k)}(z_0)$ ,  $||f^{(n-k)} - P_k||_X < C^k d^k \epsilon$ , and  $p_{\alpha}(f^{(n-k)} - P_k) \le C^k d^{k-\alpha} \epsilon$ , for  $k = 1, 2, \ldots, n$ . Clearly  $P_n^{(k)} = P_{n-k}$  on X and

$$||f - P_n||_{\ell i p^n(X,\alpha)} \le \sum_{k=0}^{n-1} \frac{C^{n-k} d^{n-k} \epsilon + C^{n-k} d^{n-k-\alpha} \epsilon}{k!} + \frac{\epsilon}{n!} = \lambda \epsilon,$$

for some constant  $\lambda$ . Hence  $f \in \ell ip_P^n(X, \alpha)$ .

Now we investigate rational approximation on circles and annuli. We note that when X is uniformly regular and  $\alpha < 1$ , then  $lip(X, M, \alpha) = Lip(X, M, \alpha)$  if  $1/M_k \neq 0$  for infinitely many k.

**Theorem 2.30** [14] If  $T = \{ z \in C : |z-z_0| = R \}$  then  $\ell i p_R(T, M, \alpha) = \ell i p(T, M, \alpha)$ .

**Proof.** We assume that  $z_0 = 0$  and R = 1. Let  $f \in lip(T, M, \alpha)$  and  $\sum_{-\infty}^{\infty} a_j z^j$  be the Fourier series generated by f, where  $a_j = \frac{1}{2\pi} \int_{-\pi}^{\pi} f(e^{i\theta}) e^{-ij\theta} d\theta$ . The Cesaro means of this series are

$$\sigma_n(e^{i\theta}) = \frac{1}{2\pi} \int_{-\pi}^{\pi} f(e^{it}) K_n(\theta - t) dt = \frac{1}{2\pi} \int_{-\pi}^{\pi} f(e^{i(\theta - t)}) K_n(t) dt,$$

where  $K_n(t)$  is the Fejer kernel. It is known that  $\sigma_n$  is a rational function on T with the only pole z=0, and  $\|\sigma_n-f\|_T\to 0$  as  $n\to\infty$ . Since for each  $k\geq 0$ ,  $f^{(k)}$  is continuous on T we have

$$\sigma_n^{(k)}(z) = \frac{1}{2\pi} \int_{-\pi}^{\pi} e^{-ikt} f^{(k)}(ze^{-it}) K_n(t) dt \qquad (z \in T),$$

and so  $\|\sigma_n^{(k)}\|_T \leq \|f^{(k)}\|_T$ .

On the other hand,

$$\frac{|\sigma_n^{(k)}(z) - \sigma_n^{(k)}(w)|}{|z - w|^{\alpha}} \le \frac{1}{2\pi} \int_{-\pi}^{\pi} \frac{|f^{(k)}(ze^{-it}) - f^{(k)}(we^{-it})|}{|ze^{-it} - we^{-it}|^{\alpha}} K_n(t) dt \le p_{\alpha}(f^{(k)}).$$

Hence  $p_{\alpha}(\sigma_n^{(k)}) \leq p_{\alpha}(f^{(k)})$  and so  $\sigma_n \in lip_R(T, M, \alpha)$ .

Now we prove that  $\|\sigma_n - f\|_{\ell i p(T, M, \alpha)} \to 0$  as  $n \to \infty$ . Since

$$\|\sigma_n - f\|_{\ell i p(T, M, \alpha)} \le 2 \|f\|_{\ell i p(T, M, \alpha)},$$

by the dominated convergence theorem, it is enough to show that for each  $k \geq 0$ ,  $\|\sigma_n^{(k)} - f^{(k)}\|_T + p_{\alpha}(\sigma_n^{(k)} - f^{(k)}) \to 0$  as  $n \to \infty$ .

By the uniform continuity of each  $f^{(k)}$  on T we have  $\|\sigma_n^{(k)} - f^{(k)}\|_T \to 0$  as  $n \to \infty$ .

Since  $f^{(k)} \in \ell ip(T, \alpha)$ , for  $\epsilon > 0$  there exists  $\delta > 0$  such that for all  $z, w \in T$ , if  $0 < |z - w| < \delta$  then  $|f^{(k)}(z) - f^{(k)}(w)|/|z - w|^{\alpha} < \epsilon/2$ . Let  $k \geq 0$  and  $z, w \in T$ ,  $(z \neq w)$ . If  $|z - w| < \delta$ , then

$$\frac{|\sigma_n^{(k)}(z) - f^{(k)}(z) - \sigma_n^{(k)}(w) + f^{(k)}(w)|}{|z - w|^{\alpha}} < \epsilon.$$

If  $|z - w| \ge \delta$  and n is large enough, then

$$\frac{|\sigma_n^{(k)}(z) - f^{(k)}(z) - \sigma_n^{(k)}(w) + f^{(k)}(w)|}{|z - w|^{\alpha}} \le \frac{2||\sigma_n^{(k)} - f^{(k)}||_T}{\delta^{\alpha}} < \epsilon.$$

Hence  $p_{\alpha}(\sigma_n^{(k)} - f^{(k)}) \to 0$  as  $n \to \infty$ . This completes the proof of the theorem.

Note that the following results are not satisfied when  $lip(X, M, \alpha)$  reduces to  $lip(X, \alpha)$ .

**Theorem 2.31** [14] If  $X = \{z \in C : r \le |z - z_0| \le R\}$ , where 0 < r < R, then  $\ell i p_R(X, M, \alpha) = \ell i p(X, M, \alpha)$ . But in these cases we have  $\ell i p_R(X, \alpha) = \ell i p_A(X, \alpha)$ .

**Proof.** Without loss of generality we can assume that  $z_0=0$ . Let  $f\in \ell ip(X,M,\alpha)$ . Since f is analytic in r<|z|< R it has a Laurent series of the form  $f(z)=\sum_{-\infty}^{\infty}a_jz^j$  on r<|z|< R, where  $a_j=(2\pi\rho^j)^{-1}\int_{-\pi}^{\pi}e^{-ijt}f(\rho e^{it})dt$ , for  $r<\rho< R$ . The Cesaro means of the Laurent series of f is

$$\sigma_n(z) = \frac{1}{2\pi} \int_{-\pi}^{\pi} f(ze^{-it}) K_n(t) dt \qquad (r < |z| < R),$$

where  $K_n(t)$  is the Fejer kernel. Clearly for each  $k \geq 0$  we have

$$\sigma_n^{(k)}(z) = \frac{1}{2\pi} \int_{-\pi}^{\pi} e^{-ikt} f^{(k)}(ze^{-it}) K_n(t) dt \qquad (r < |z| < R),$$

and so  $|\sigma_n^{(k)}(z)| \leq ||f^{(k)}||_X$  for r < |z| < R. Since  $\sigma_n$  is a rational function with the only pole z = 0,  $\sigma_n^{(k)}$  is analytic in  $r \leq |z| \leq R$ . Therefore the above inequality holds for all z in  $r \leq |z| \leq R$ . Hence  $||\sigma_n^{(k)}||_X \leq ||f^{(k)}||_X$  and  $p_\alpha(\sigma_n^{(k)}) \leq p_\alpha(f^{(k)})$  for all  $k \geq 0$  and for every positive integer n, and so  $\sigma_n \in lip_R(X, M, \alpha)$ . Now we can proceed exactly the same as in the proof of theorem 2.30 to show that  $||\sigma_n - f||_{lip(X,M,\alpha)} \to 0$  as  $n \to \infty$ . Therefore  $f \in lip_R(X,M,\alpha)$  and this completes the proof of the theorem.

If  $r \to 0$ , the above theorem implies the following result.

Corollary 2.32 [14] If  $X = \{z : |z| \leq R\}$  then  $\ell i p_P(X, M, \alpha) = \ell i p(X, M, \alpha)$ .

**Theorem 2.33** [14] Let X be a regular set for which there exists  $z_0 \in X$  such that for  $0 \le \beta < 1$ ,  $\beta(z-z_0)+z_0 \in int X$  for all  $z \in X$ . Or, equivalently, the segment  $[z_0,z)$  is contained in the interior of X for all  $z \in X$ . If  $P_k \to \infty$  as  $k \to \infty$ , then  $\ell ip_P(X,M,\alpha) = \ell ip(X,M,\alpha)$ . **Proof.** Clearly X is star-shaped and so it is polynomially convex. Thus

$$\ell i p_P(X, M, \alpha) = \ell i p_R(X, M, \alpha) = \ell i p_H(X, M, \alpha).$$

Without loss of generality we can assume that  $z_0=0$ . By the hypothesis for each positive integer n and every  $z\in X$ ,  $r_nz\in intX$ , where  $r_n=n/(n+1)$ . Let  $f\in lip(X,M,\alpha)$  and define the sequence  $\{f_n\}$  on X by  $f_n(z)=f(r_nz)$ . Each  $f_n$  is analytic in a neighbourhood of X and so  $f_n\in lip_H(X,M,\alpha)$ . Moreover for each  $k\geq 0$ ,  $f_n^{(k)}(z)=r_n^kf^{(k)}(r_nz)$  and so  $\|f_n^{(k)}\|_X\leq \|f^{(k)}\|_X$ ,  $p_\alpha(f_n^{(k)})\leq p_\alpha(f^{(k)})$  for all  $k\geq 0$  and every n. By the uniform continuity of each  $f^{(k)}$  on  $f^{(k)}$  on  $f^{(k)}$  on  $f^{(k)}$  on  $f^{(k)}$  or each  $f^{(k)}$  or each  $f^{(k)}$  or each  $f^{(k)}$  or each  $f^{(k)}$  or  $f^{(k)}$  or each  $f^{(k)}$  or  $f^{(k)$ 

Corollary 2.34 If X is a compact convex set with non-empty interior and  $P_k \to \infty$  as  $k \to \infty$ , then  $\ell ip_P(X, M, \alpha) = \ell ip(X, M, \alpha)$ .

# 3. Real Lipschitz Algebras

We first present a brief general description of real uniform algebras and real Banach function algebras.

### 3.1 Real Uniform (Function) Algebras

**Definition 3.1.1** Let X be a topological space. A map  $\tau: X \longrightarrow X$  is called a *topological involution* on X if  $\tau$  is continuous and  $\tau(\tau(x)) = x$  for all  $x \in X$ .

Note that,  $\tau$  is a homeomorphism of X onto X.

**Examples 3.1.2** (i) The identity map on a topological space X is a topological involution on X.

- (ii) Let X be the closed interval [a, b]. The map  $\tau: X \longrightarrow X$ , defined by  $\tau(x) = b a x$  is a topological involution on X.
- (iii) Let X be a compact plane set such that X is symmetric about the real axis. The map  $\tau: X \longrightarrow X$ , defined by  $\tau(z) = \bar{z}$ , is a topological involution on X.

Let X be a compact Hausdorff space and  $\tau$  be a topological involution on X. Let

$$C(X,\tau) := \{ f \in C(X) : f \circ \tau = \bar{f} \}.$$

Then  $C(X,\tau)$  is a uniformly closed real subalgebra of C(X) which contains the constant function 1.

Remark 3.1.3 If  $\tau$  is the identity map on a compact Hausdorff space X then  $C(X,\tau)=C_{\mathbb{R}}(X)$ , the real uniform algebra of continuous realy valued functions on X. On the other hand, for a given compact Hausdorff space X, we suppose  $Y=X\times\{0,1\}$  and the map  $\tau:Y\longrightarrow Y$  defines by  $\tau(x,j)=(x,(j+1) \text{ modulo 2})$ . Identifying X with  $X\times\{0\}$ , X can be treated as a subspace of Y. Then every f in C(X) can be extended uniquely to Y by requiring that the extension belongs to  $C(Y,\tau)$ . In fact, the map  $\psi:C(X)\longrightarrow C(Y,\tau)$ , defined by  $\psi(f)(x,0)=f(x)$ ,  $\psi(f)(x,1)=\overline{f(x)}$ , is an isometrical isomorphism of C(X), as a real Banach algebra, onto  $C(Y,\tau)$ . Thus  $C(X,\tau)$  is a more general object than  $C_{\mathbb{R}}(X)$  and C(X).

Theorem 3.1.4 [18; Theorem 1.3.5]. Let X be a compact Hausdorff space and  $\tau$  be a topological involution on X. Define the map  $\sigma: C(X) \longrightarrow C(X)$  by  $\sigma(f) = \bar{f} \circ \tau$ . Then:

(i)  $\sigma$  is an algebra involution on C(X) and

$$C(X,\tau) = \{ f \in C(X) : \sigma(f) = f \}.$$

- (ii)  $C(X) = C(X, \tau) \oplus iC(X, \tau)$ , that is, every h in C(X) can be expressed uniquely as f + ig with f, g in  $C(X, \tau)$ .
- (iii)  $\sigma$  is an isometry.

Definition 3.1.5 Let X be a compact Hausdorff space and  $\tau$  be a topological involution on X. The algebra involution  $\sigma$  on C(X) defined by  $\sigma(f) = \bar{f} \circ \tau$ , is called the algebra involution on C(X) induced by  $\tau$ .

Theorem 3.1.6 [18; Theorem 1.3.5(v)] Let X be a compact Hausdorff space. If  $\sigma$  is an algebra involution on C(X) then:

- (i) There exists a topological involution on X such that  $\sigma(f) = \bar{f} \circ \tau$ , for all f in C(X).
- (ii)  $\sigma$  is an isometry.

Theorem 3.1.7 [18; Lemma 1.3.7] Let X be a compact Hausdorff space and  $\tau$  be a topological involution on X. Let  $x, y \in X$  and  $x \neq y$ .

- (i) If  $y = \tau(x)$ , then there exists a function f in  $C(X, \tau)$  such that f(x) = i and f(y) = -i.
- (ii) If  $y \neq \tau(x)$ , then there exists a function f in  $C(X, \tau)$  such that f(x) = 1 and f(y) = 0.

In particular,  $C(X, \tau)$  separates the points of X.

**Definition 3.1.8** Let X be a compact Hausdorff space and  $\tau$  be a topological involution on X. A real uniform (function) algebra on  $(X, \tau)$ , is a uniformly closed subalgebra of  $C(X, \tau)$  which contains 1 and separates the points of X.

It is clear that  $C(X,\tau)$  is a uniform (function) algebra on  $(X,\tau)$ .

**Theorem 3.1.9** [18; Theorem 1.3.20]. Let X be a compact Hausdorff space,  $\tau$  be a topological involution on X and  $\sigma$  be the algebra involution induced by  $\tau$  on C(X). Let A be a real subspace of  $C(X,\tau)$  and define

$$B:=\{f+ig:f,g\in A\}.$$

Then

- (i)  $\sigma(B) = B$  and  $A = \{h \in B : \sigma(h) = h\} = B \cap C(X, \tau)$ .
- (ii)  $B = A \oplus iA$ , that is, every  $h \in B$  can be written uniquely as f + ig with f, g in A.
- (iii) For  $f, g \in A$ ,

$$\max\{\|f\|_X, \|g\|_X\} \le \|f + ig\|_X \le \|f\|_X + \|g\|_X.$$

- (iv) B is uniformly closed (self-adjoint, separates the points of X, contains 1) if and only if A has the same property, respectively. Moreover, B is a complex algebra if and only if A is a real algebra.
- (v) If A is a real uniform algebra on  $(X, \tau)$  then the map  $\alpha: M_A \longrightarrow M_B$ , defined by

$$\alpha(\phi)(f+ig) = \phi(f) + i\phi(g),$$

is a homeomorphism of  $M_A$  onto  $M_B$ 

By the above theorem, if  $\tau$  is a topological involution on a compact Hausdorff space X then

$$M_{C(X,\tau)} = \{e_x : x \in X\},\,$$

where  $e_x$  is the evaluation character on X at  $x \in X$  [18; Corollary 1.3.21].

In some problems, we need to know when a given complex uniform algebra on  $(X,\tau)$  can be viewed as a complexification of a real uniform algebra. The following theorem gives a criterion.

**Theorem 3.1.10** [18; Theorem 1.3.22] Let X be a compact Hausdorff space,  $\tau$  be a topological involution on X and  $\sigma$  be the algebra involution induced by  $\tau$  on C(X). If B is a complex uniform algebra on X with  $\sigma(B) = B$  and suppose  $A = \{h \in B : \sigma(h) = h\}$  then A is a real uniform algebra on  $(X, \tau)$  and B can be regarded as the complexification of A.

We now state the Stone-Weierstrass theorem for real subalgebras of  $C(X,\tau)$ .

Theorem 3.1.11 [18; Proposition 1.2, Corollary 2.1.14] Let X be a compact Hausdorff space and  $\tau$  be a topological involution on X. If A is a self-adjoint real subalgebra of  $C(X,\tau)$  containing 1 and separating the points of X then  $\bar{A} = C(X,\tau)$ , where  $\bar{A}$  is the uniform closure of A.

## 3.2 Real Banach Function Algebras

Real uniform (function) algebras were first defined and studied by S.H. Kulkarni and B.V. Limaye in 1981. In general, a complex uniform (function) algebra on a compact Hausdorff space X may not be a real uniform (function) algebra on X, with the topological involution  $\tau$  as the identity function. But, it is interesting to note that every complex uniform (function) algebra can be regarded as a real uniform (function)

algebra on a compact Hausdorff space with a suitable topological involution  $\tau$ . Hence, the class of real uniform algebras is larger than that of complex uniform algebras.

Now we extend the notion of real uniform (function) algebras, by introducing a larger class, which is called the *real Banach function* algebras.

We will show that every complex Banach function algebra can be viewed as a real Banach function algebra with some topological involution  $\tau$ . Hence the class of real Banach function algebras is larger than the class of complex Banach function algebras. We are going to extend some general properties of the complex Banach function algebras for the real Banach function algebras. Then the real Lipschitz algebras of complex functions are introduced and some properties of this interesting class of real Banach function algebras are discussed.

Let X be a compact Hausdorff space and  $\tau$  be a topological involution on X. In this section, we first define a real Banach function algebra on  $(X,\tau)$ . We prove that for each complex Banach function algebra  $(B,\|\cdot\|)$  on X, there exist a compact Hausdorff space Y, a topological involution  $\tau$  on Y and a real Banach function algebra  $(A,\|\|\cdot\|\|)$  on  $(Y,\tau)$  such that  $(B,\|\cdot\|)$ , as a real Banach algebra, is isometrically isomorphic to  $(A,\|\|\cdot\|\|)$ . Then we extend the Theorems 3.1.9 and 3.1.10 to real Banach function algebras. Finally we give some results concerning the carrier space (maximal ideal space) of real Banach function algebras.

**Definition 3.2.1** Let X be a compact Hausdorff space and  $\tau$  be a topological involution on X. A real Banach function algebra on  $(X,\tau)$  is a real subalgebra A of  $C(X,\tau)$  which contains the constant function 1 and separates the points of X and there exists an algebra norm  $\|\cdot\|$  on A such that  $(A,\|\cdot\|)$  is a real Banach algebra.

Note that if the norm of a real Banach function algebra A on  $(X, \tau)$ 

is the same as the uniform norm on X, then A is called a real uniform (function) algebra on  $(X, \tau)$ .

Let X be a compact Hausdorff space and  $\tau$  be a topological involution on X. We recall that the map  $\sigma: C(X) \longrightarrow C(X)$  which is defined by  $\sigma(f) = \bar{f} \circ \tau$ , is an isometric algebra involution on C(X) and is called the algebra involution on C(X) induced by  $\tau$ .

**Theorem 3.2.2** [2; Theorem 1.1] Let X be a compact Hausdorff space,  $\tau$  be a topological involution on X and  $\sigma$  be the algebra involution on C(X) induced by  $\tau$ . If  $(B, \|\cdot\|)$  is a complex Banach function algebra on X such that  $\sigma(B) = B$  and  $A := \{h \in B : \sigma(h) = h\}$ , then

- (i) A is a real subalgebra of B and  $A = B \cap C(X, \tau)$ .
- (ii) Every h in B can be expressed uniquely as f + ig with  $f, g \in A$ .
- (iii) There exists a constant  $C \ge 1$  such that  $\|\sigma(h)\| \le C\|h\|$  for every  $h \in B$  and  $\max\{\|f\|, \|g\|\} \le C\|f + ig\|$  for all  $f, g \in A$ .
  - (iv)  $(A, \|\cdot\|)$  is a real Banach function algebra on  $(X, \tau)$ .
- (v) For  $\phi$  in  $M_A$ , define  $\alpha(\phi)(f+ig) := \phi(f) + i\phi(g)$  for  $f,g \in A$ . Then  $\alpha(\phi) \in M_B$  and the map  $\alpha$  is a homeomorphism from  $M_A$  onto  $M_B$ . In particular, A is natural if and only if B is natural.

Remark 3.2.3 In the above theorem, if B is a complex uniform algebra on X then A is a real uniform algebra on  $(X,\tau)$  and we can choose C=1. Therefore, the above theorem is a generalization of a similar result for the real uniform algebras.

**Theorem 3.2.4** [2; Theorem 1.2] Let X be a compact Hausdorff space,  $\tau$  be a topological involution on X and  $\sigma$  be the algebra involution on C(X) induced by  $\tau$ . Let  $(A, \|\cdot\|)$  be a real Banach function algebra on  $(X, \tau)$  and define  $B := \{f + ig : f, g \in A\}$ . Then:

- (i)  $\sigma(B) = B$  and  $A = \{h \in B : \sigma(h) = h\} = B \cap C(X, \tau)$ .
- (ii) B is complex subalgebra of C(X) and  $B = A \oplus iA$ .

(iii) There is an algebra norm  $|||\cdot|||$  on B such that ||f||=|||f||| for all  $f\in A$  and

$$\max\{\|f\|,\|g\|\} \le \||f+ig\|| \le 2\max\{\|f\|,\|g\|\} \quad (f,g \in A).$$

- (iv)  $(B, ||| \cdot |||)$  is a complex Banach function algebra on X.
- (v) For  $\phi \in M_A$  define  $(\alpha(\phi))(f+ig) = \phi(f) + i\phi(g)$  for  $f,g \in A$ . Then  $\alpha(\phi) \in M_B$  and  $\alpha$  is a homeomorphism from  $M_A$  onto  $M_B$ . In particular, A is natural if and only if B is natural.

**Remark 3.2.5** In the above theorem, if A is a real uniform algebra on  $(X, \tau)$  then B is a complex uniform algebra on X. Therefore, this result is a generalization of a similar result for the real uniform algebras.

## 3.3 Real Lipschitz Algebras of Complex Functions

In this section we first define C-quasicontraction and in particular d-isometric topological involutions on a compact metric space (X,d). Then we show that if  $\sigma$  is the algebra involution on C(X) induced by a C-quasicontraction topological involution  $\tau$  on a compact metric space (X,d) and  $\alpha \in (0,1]$  then  $\sigma(Lip(X,\alpha)) = Lip(X,\alpha)$  and  $\sigma(lip(X,\alpha)) = lip(X,\alpha)$ . By using this result we define the real Lipschitz algebras of complex functions  $Lip(X,\tau,\alpha)$  and  $lip(X,\tau,\alpha)$ , which are real Banach function algebras on  $(X,\tau)$ . Next, we show that for a compact metric space (X,d), there exists a compact metric space  $(Y,\rho)$  and a  $\rho$ -isometric topological involution  $\tau$  on Y such that  $Lip(X,\alpha)$  (respectively,  $lip(X,\alpha)$ ), as a real Banach algebra, is isometrically isomorphic to  $Lip(Y,\tau,\alpha)$  (respectively,  $lip(Y,\tau,\alpha)$ ) for  $\alpha \in (0,1]$  (respectively,  $\alpha \in (0,1)$ ). Finally, we study the approximation problem of the real Lipschitz algebras  $Lip(X,\tau,\alpha)$  and  $lip(X,\tau,\alpha)$  and certain subalgebras of these algebras.

. M **Definition 3.3.1** Let (X,d) be a compact metric space and  $\tau$  be a topological involution on X.

- (i)  $\tau$  is called a C- quasicontraction on X if there exists a constant C > 0 such that  $d(\tau(x), \tau(y)) \leq Cd(x, y)$ , for all x, y in X.
- (ii)  $\tau$  is called a *d-isometric topological involution on* Xif  $d(\tau(x), \tau(y)) = d(x, y)$  for all x, y in X.

Remark 3.3.2 If  $\tau$  is a C-quasicontraction topological involution on a compact metric space (X,d) then it is easy to see that  $C \geq 1$ . Moreover if C = 1 then  $\tau$  is d-isometric.

**Lemma 3.3.3** [2; Lemma 2.4] Let (X, d) be a compact metric space and  $\tau$  be a C-quasicontraction topological involution on X and  $\sigma$  be the algebra involution induced by  $\tau$  on C(X). Then

- (i) For every  $\alpha \in (0,1]$  and  $f \in Lip(X,\alpha)$ ,  $p_{\alpha}(\sigma(f)) \leq C^{\alpha}p_{\alpha}(f)$ .
- (ii) For every  $\alpha \in (0,1]$ ,  $\sigma(Lip(X,\alpha)) = Lip(X,\alpha)$ .
- (iii) For every  $\alpha \in (0,1]$ ,  $\sigma(lip(X,\alpha)) = lip(X,\alpha)$ .
- (iv) For every  $\alpha \in (0,1]$  and every  $f \in Lip(X,\alpha)$ ,  $\|\sigma(f)\|_{\alpha} \leq C^{\alpha} \|f\|_{\alpha}$ .
- (v) If  $\tau$  is a d-isometric topological involution then for every  $\alpha \in (0,1]$  and  $f \in Lip(X,\alpha)$ ,  $p_{\alpha}(\sigma(f)) = p_{\alpha}(f)$  and  $\|\sigma(f)\|_{\alpha} = \|f\|_{\alpha}$ . Hence  $\sigma$  is an isometric algebra involution on  $Lip(X,\alpha)$ .

**Proof.** (i) Let  $\alpha \in (0,1]$  and  $f \in Lip(X,\alpha)$ . Then

$$p_{\alpha}(\sigma(f)) = \sup \{ \frac{|\sigma(f)(x) - \sigma(f)(y)|}{d^{\alpha}(x,y)} : x, y \in X, x \neq y \}$$

$$= \sup \{ \frac{|f(\tau(x)) - f(\tau(y))|}{d^{\alpha}(x,y)} : x, y \in X, x \neq y \}$$

$$\leq \sup \{ \frac{|f(\tau(x)) - f(\tau(y))|}{C^{-\alpha}d^{\alpha}(\tau(x),\tau(y))} : x, y \in X, x \neq y \}$$

$$= C^{\alpha}p_{\alpha}(f).$$

(ii) By (i),  $\sigma(Lip(X,\alpha)) \subseteq Lip(X,\alpha)$ . Since  $\sigma(\sigma(f)) = f$  for every  $f \in Lip(X,\alpha)$ ,  $\sigma(Lip(X,\alpha)) = Lip(X,\alpha)$ .

(iii) Let  $\alpha \in (0,1)$  and  $f \in lip(X,\alpha)$ . Let  $\varepsilon > 0$  be given. Since  $f \in lip(X,\alpha)$ , there is a  $\delta_1 > 0$  such that  $\frac{|f(x)-f(y)|}{d^{\alpha}(x,y)} < C^{-\alpha}\varepsilon$ , whenever  $0 < d(x,y) < \delta_1$ . Set  $\delta = C^{-1}\delta_1$  and suppose  $x,y \in X$  such that  $0 < d(x,y) < \delta$ . Since  $\tau$  is a C-quasicontraction topological involution on (X,d), therefore

$$0 < d(\tau(x), \tau(y)) \le Cd(x, y) < C\delta = \delta_1.$$

Hence  $\frac{|f(\tau(x))-f(\tau(y))|}{d^{\alpha}(\tau(x),\tau(y))} < C^{-\alpha}\varepsilon$ . This implies that  $\frac{|\sigma(f)(x)-\sigma(f)(y)|}{d^{\alpha}(x,y)} < \varepsilon$  and so  $\sigma(f) \in lip(X,\alpha)$ . Therefore,  $\sigma(lip(X,\alpha)) \subseteq lip(X,\alpha)$  and since  $\sigma$  is an algebra involution on C(X),  $\sigma(lip(X,\alpha)) = lip(X,\alpha)$ .

(iv) Let  $\alpha \in (0,1]$  and  $f \in Lip(X,\alpha)$ . By (i) and  $C \geq 1$ , we have

$$\|\sigma(f)\|_{\alpha} = \|\sigma(f)\|_{X} + p_{\alpha}(\sigma(f))$$

$$= \|f\|_{X} + p_{\alpha}(\sigma(f))$$

$$\leq C^{\alpha}\|f\|_{X} + C^{\alpha}p_{\alpha}(f) = C^{\alpha}\|f\|_{\alpha}.$$

(v) If  $\tau$  is a d-isometric topological involution on X then C=1 and so  $p_{\alpha}(\sigma(f))=p_{\alpha}(f)$ . Hence  $\|\sigma(f)\|_{\alpha}=\|f\|_{\alpha}$ .  $\square$ 

Remark 3.3.4 Let (X,d) be a compact metric space and  $\alpha \in (0,1]$ . If  $B = Lip(X,\alpha)$  or  $lip(X,\alpha)$ , then the map  $f \mapsto \bar{f}$  is an isometric algebra involution on B.

Theorem 3.3.5 [2; Theorem 2.7] Let (X, d) be a compact metric space,  $\tau$  be a C-quasicontraction topological involution on X and  $\sigma$  be the algebra involution on C(X) induced by  $\tau$ . We define

$$Lip(X,\tau,\alpha):=\{h\in Lip(X,\alpha):\sigma(h)=h\}\quad (\alpha\in(0,1]),$$

and

$$lip(X,\tau,\alpha):=\{h\in lip(X,\alpha):\sigma(h)=h\}\quad (\alpha\in(0,1)).$$

If  $A = Lip(X, \tau, \alpha)$  and  $B = Lip(X, \alpha)$   $(A = lip(X, \tau, \alpha))$  and  $B = lip(X, \alpha)$ , respectively), then

- (i)  $B = A \oplus iA$ .
- (ii) For every  $f, g \in A$ ,  $\max\{\|f\|_{\alpha}, \|g\|_{\alpha}\} \leq C^{\alpha}\|f + ig\|_{\alpha}$ .
- (iii)  $(A, \|\cdot\|_{\alpha})$  is a real Banach function algebra on  $(X, \tau)$ .
- (iv) A is self-adjoint and  $\bar{A} = C(X, \tau)$ , where  $\bar{A}$  is the uniform closure of A.
  - $(v) M_A = \{e_x : x \in X\}.$

**Proof.** Since  $(B, \|\cdot\|_{\alpha})$  is a complex Banach function algebra on X and by Lemma 3.3.3, we have  $\sigma(B) = B$ , therefore (i) and (iii) hold by Theorem 3.2.2.

Since  $\sigma$  is an isometric involution on C(X) and by Lemma 3.3.3,  $\|\sigma(h)\|_{\alpha} \leq C^{\alpha} \|h\|_{\alpha}$  for each  $h \in B$ , we can easily show that (ii) holds.

Since A is self-adjoint,  $\bar{A}=C(X,\tau)$ . Thus (iv) holds. Since B is a natural Banach function algebra on X, A is also natural by Theorem 3.2.2(v) and hence (v) holds.  $\Box$ 

Theorem 3.3.6 [2; Theorem 2.8] Let (X,d) be a compact metric space and  $\alpha \in (0,1]$ . Then there exist a compact metric space  $(Y,\rho)$ , a  $\rho$ -isometric topological involution  $\tau$  on Y such that the complex Lipschitz algebra  $Lip(X,\alpha)$  (respectively,  $lip(X,\alpha)$ ), regarded as a real Banach algebra, is isometrically isomorphic to the real Lipschitz algebra  $Lip(Y,\tau,\alpha)$  (respectively,  $lip(Y,\tau,\alpha)$ ).

**Proof.** Let  $Y := X \times \{0,1\}$ . We define the map  $\rho: Y \times Y \longrightarrow \mathbb{R}$  by

$$\rho((x,j),(y,k)) = \max\{d(x,y),|j-k|\} \quad (x,y \in X, j,k \in \{0,1\}).$$

Then  $\rho$  is a metric on Y and the topology induced by  $\rho$  on Y coincides to the product topology on Y. Now, we define the map  $\tau: Y \longrightarrow Y$  by

$$\tau(x,0) = (x,1), \quad \tau(x,1) = (x,0) \quad (x \in X).$$

It is easy to see that  $\tau$  is a  $\rho$ -isomeric topological involution on Y. Now, we define the map  $\psi: Lip(X,\alpha) \longrightarrow Lip(Y,\tau,\alpha)$  by

$$\psi(f)(x,0) = f(x), \quad \psi(f)(x,1) = \bar{f}(x) \quad (x \in X).$$

We can easily show that  $\psi$  is an isometrically homomorphism from  $(Lip(X,\alpha), \|\cdot\|_{\alpha})$ , as a real Banach algebra, into  $(Lip(Y,\tau,\alpha), \|\cdot\|_{\alpha})$ . Now, let  $g \in Lip(Y,\tau,\alpha)$  and define the map  $f: X \longrightarrow \mathbb{C}$  by

$$f(x) = g(x,0).$$

It is easy to see that  $f \in Lip(X,\alpha)$  and  $\psi(f) = g$ . Therefore,  $\psi$  is an isometrically isomorphism from  $(Lip(X,\alpha),\|\cdot\|_{\alpha})$ , as a real Banach algebra, onto  $Lip(Y,\tau,\alpha)$ . Moreover, if  $f \in lip(X,\alpha)$  then we conclude that  $\psi(f) \in lip(Y,\tau,\alpha)$  and if  $g \in lip(Y,\tau,\alpha)$ ,  $\psi^{-1}(g) \in lip(X,\alpha)$ . Therefore  $\psi|_{lip(X,\alpha)}$  is an isometrically isomorphism from  $(lip(X,\alpha),\|\cdot\|_{\alpha})$ , as a real Banach algebra, onto  $lip(Y,\tau,\alpha),\|\cdot\|_{\alpha}$ .  $\square$ 

Remark 3.3.7 If (X,d) is a compact metric space and  $\tau$  is the identity map on X then  $\tau$  is a d-isometric topological involution on X and

$$Lip(X, \tau, \alpha) = Lip_{\mathbb{R}}(X, \alpha) = \{ f \in Lip(X, \alpha) : f \text{ is real-valued} \},$$
  
$$lip(X, \tau, \alpha) = lip_{\mathbb{R}}(X, \alpha) = \{ f \in lip(X, \alpha) : f \text{ is real-valued} \}.$$

Therefore the class of real Lipschitz algebras of complex functions  $Lip(X,\tau,\alpha)$  ( $lip(X,\tau,\alpha)$ , respectively) is larger than the class of the real Lipschitz algebras  $Lip_{\mathbb{R}}(X,\alpha)$  ( $lip_{\mathbb{R}}(X,\alpha)$ , respectively). Also, by Theorem 3.3.6 we conclude that the class of real Lipschitz algebras of complex functions  $Lip(X,\tau,\alpha)$  ( $lip(X,\tau,\alpha)$ , respectively) is larger than the class of complex Lipschitz algebras  $Lip(X,\alpha)$  ( $lip(X,\alpha)$ , respectively), as real Banach algebras.

**Theorem 3.3.8** [2; Theorem 2.9] Let (X,d) be a compact metric space and  $\tau$  be a C-quasicontraction topological involution on X. Then

- (i) If  $0 < \alpha < \beta \le 1$ , then  $Lip(X, \tau, \beta)$  is a subalgebra of  $lip(X, \tau, \alpha)$ .
- (ii) If  $0 < \alpha < 1$  and X is an infinite set then  $lip(X, \tau, \alpha)$  is a proper subalgebra of  $Lip(X, \tau, \alpha)$ .
- **Proof.** (i) Since  $Lip(X,\tau,\beta) = Lip(X,\beta) \cap C(X,\tau)$ ,  $lip(X,\tau,\alpha) = lip(X,\alpha) \cap C(X,\tau)$ ,  $Lip(X,\beta)$  is a complex subalgebra of  $lip(X,\alpha)$  and  $C(X,\tau)$  is a real subalgebra of C(X), we conclude that  $Lip(X,\tau,\beta)$  is a real subalgebra of  $lip(X,\tau,\alpha)$ .
- (ii) Since  $lip(X, \alpha)$  is a proper subalgebra of  $Lip(X, \alpha)$ , there exists  $h \in Lip(X, \alpha) \setminus lip(X, \alpha)$ . Since  $h \in Lip(X, \alpha)$  and by Theorem 3.3.5 we have

$$Lip(X, \alpha) = Lip(X, \tau, \alpha) \oplus iLip(X, \tau, \alpha),$$

thus h can be expressed uniquely as h = f + ig with f, g in  $Lip(X, \tau, \alpha)$ . Since  $h \notin lip(X, \alpha)$  and by Theorem 3.3.5 we have

$$lip(X, \alpha) = lip(X, \tau, \alpha) \oplus ilip(X, \tau, \alpha),$$

thus  $f \notin lip(X, \tau, \alpha)$  or  $g \notin lip(X, \tau, \alpha)$  and the proof (ii) is now complete.  $\square$ 

Let (X,d) be a compact metric space and take  $\alpha \in (0,1)$ . A type of Stone-Weierstrass theorem in real Lipschitz algebra  $lip_{\mathbb{R}}(X,\alpha)$  was first given by L. I. Hedberg in 1969 [10; Theorem 1]. Let  $\tau$  be a C-quasicontraction topological involution on X. We now extend the Hedberg's theorem in real Lipschitz algebra of complex functions  $lip(X,\tau,\alpha)$ , without using the complexification technique.

Theorem 3.3.9 (Hedberg's theorem in real Lipschitz algebras of complex functions) [2; Theorem 2.10] Let (X, d) be a compact metric space,  $\tau$  be a C-quasicontraction topological involution on X, and take  $\alpha \in (0,1)$ . Let A be a self-adjoint real subalgebra of  $lip(X, \tau, \alpha)$  which

separates the points of X and contains the real-valued constant functions on X. Then A is dense in  $lip(X, \tau, \alpha)$  if for every  $a \in X$ , there are positive numbers  $M_a$  and  $\delta_a$  such that for every  $\delta \leq \delta_a$ , there is a  $f \in A$ , with f(a) = 1, f(x) = 0 on  $S_{\delta}(a) = \{x \in X : d(x, a) = \delta\}$ , and

$$\sup\left\{\frac{|f(y)-f(z)|}{d^{\alpha}(y,z)}:y,z\in B_{\delta}(a),y\neq z\right\}<\frac{M_a}{\delta^{\alpha}},$$

where  $B_{\delta}(a) = \{x \in X : d(x, a) \leq \delta\}.$ 

Notice that if (X,d) is a compact metric space and  $\tau$  is the identity map on X then  $Lip(X,\tau,\alpha)=Lip_{\mathbb{R}}(X,\alpha)$  and  $lip(X,\tau,\alpha)=lip_{\mathbb{R}}(X,\alpha)$ . Hence, the theorem 3.3.9 is a generalization of the Hedberg's theorem in real Lipschitz algebra  $lip_{\mathbb{R}}(X,\alpha)$ . Note that T. G. Honary and H. Mahyar stated the Hedberg's theorem in complex Lipschitz algebras  $lip(X,\alpha)$  in [13].

As a consequence of Theorem 3.3.9, the Hedberg's theorem in complex Lipschitz algebra  $lip(X, \alpha)$  is obtained, which is stated as follows:

Corollary 3.3.10 [2; Corollary 2.11] Let (X,d) be a compact metric space and take  $\alpha \in (0,1)$ . Let B be a self-adjoint complex subalgebra of  $lip(X,\alpha)$  which separates the points of X and contains the complex-valued constant functions on X. Then B is dense in  $lip(X,\alpha)$  if for  $a \in X$ , there are positive numbers  $M_a$  and  $\delta_a$  such that for every  $\delta \leq \delta_a$ , there is a  $f \in B$  such that f(a) = 1, f(x) = 0 on  $S_{\delta}(a) = \{x \in X : d(x,a) = \delta\}$ , and

$$\sup\left\{\frac{|f(x)-f(y)|}{d^{\alpha}(x,y)}:y,z\in B_a(\delta),y\neq z\right\}<\frac{M_a}{\delta^{\alpha}},$$

where  $B_{\delta}(a) = \{x \in X : d(x, a) \leq \delta\}.$ 

As an application of Theorem 3.3.9, we prove that the real Lipschitz algebra  $Lip(X,\tau,1)$  is dense in  $lip(X,\tau,\alpha)$  for  $\alpha \in (0,1)$  without using the complexification technique.

Corollary 3.3.11 [2; Corollary 2.12] Let (X, d) be a compact metric space and  $\tau$  be a C-quasicontraction topological involution on X. If  $\alpha \in (0,1)$  then  $Lip(X,\tau,1)$  is dense in  $(lip(X,\tau,\alpha), \|.\|_{\alpha})$ .

## 4. Fréchet Lipschitz Algebras

First we introduce some elementary definitions and known results concerning Fréchet algebras. For further details refer to [12, 27, 15 or 7].

**Definition 4.1** A Fréchet algebra is an algebra which is a complete metrizable topological vector space and has a neighbourhood basis  $(V_n)_{n\in\mathbb{N}}$  of zero consisting of convex sets  $V_n$  such that  $V_n.V_n\subseteq V_n$  for all  $n\in\mathbb{N}$ . We always assume that the Fréchet algebra contains the unit 1.

The topology of a Fréchet algebra A can be generated by a sequence  $(p_n)_{n\in\mathbb{N}}$  of separating submultiplicative seminorms  $(p_n(f,g) \leq p_n(f).p_n(g))$  for each  $n\in\mathbb{N}$  and every  $f,g\in A$  such that  $p_n(f)\leq p_{n+1}(f)$  for all  $n\in\mathbb{N}$  and  $f\in A$ . If A has a unit,  $p_n$  can be chosen such that  $p_n(1)=1$  [9]. We may denote the Fréchet algebra A with the above generating sequence of seminorms by  $(A,(p_n))$ . Clearly a sequence  $(f_k)$  in A converges to  $f\in A$  if and only if  $p_n(f_k-f) \xrightarrow[k\to\infty]{} 0$  for each  $n\in\mathbb{N}$ .

**Definition 4.2** The spectrum of a Fréchet algebra A is the set of all non-zero continuous complex-valued homomorphisms on A and it is denoted by  $M_A$ . We endow  $M_A$  with the Gelfand topology.

It can be shown that a complex homomorphism  $\varphi$  on the commutative Fréchet algebra  $(A,(p_n))$  is continuous if and only if there exists  $n \in \mathbb{N}$  such that  $|\varphi(f)| \leq p_n(f)$  for all  $f \in A$  [9; Remark 3.2.2(ii)].

**Definition 4.3** The radical of a Fréchet algebra A, denoted by rad(A), is the intersection of all maximal left (right) ideals in A. The Fréchet algebra A is called semisimple if  $rad(A) = \{0\}$ .

If A is a commutative Fréchet algebra then rad(A) is the intersection

of all closed maximal ideals; i.e.  $rad(A) = \bigcap_{\varphi \in M_A} \ker \varphi$  [9; Proposition 8.1.2].

The following interesting result for Fréchet algebras, which is due to Carpenter [5], is similar to the uniqueness theorem of Johnson [17] for semisimple Banach algebras.

**Theorem 4.5** If A is a commutative semisimple Fréchet algebra then A has a unique topology as a Fréchet algebra.

**Definition 4.6** A Hausdorff space X is called *hemicompact* if there exists a sequence  $(X_n)_{n\in\mathbb{N}}$  of compact subsets of X such that  $X_n\subseteq X_{n+1}$  for each  $n\in\mathbb{N}$  and every compact subset K of X is contained in some  $X_n$ . Such sequence  $(X_n)_{n\in\mathbb{N}}$  is called an admissible exhaustion of X.

If  $(A,(p_n))$  is a Fréchet algebra and  $A_n$  is the completion of  $A/\ker p_n$  with respect to the norm  $p'_n(f + \ker p_n) = p_n(f)$ ,  $f \in A$ , then  $A_n$  is a Banach algebra.

**Definition 4.7** Let X be a hemicompact space, and let A be a subalgebra of C(X) which contains the constants and separates the points of X. Then A is called a *Fréchet function algebra* on X if it is a Fréchet algebra with respect to some topology such that the evaluation homomorphisms are continuous on A, i.e.  $\varphi_x \in M_A$  for all  $x \in X$ .

Clearly with the above definition every uniform Fréchet algebra is a Fréchet function algebra which is equipped with the compact-open topology.

It is easy to see that every Fréchet function algebra is semisimple. Moreover, every Banach function algebra on a compact Hausdorff space X is a Fréchet function algebra on X. For further details see [27].

Let X be a perfect compact plane set and for  $n \in \mathbb{N}$ ,  $B_n(X)$  be any of  $D^n(X)$ ,  $Lip^n(X,\alpha)$  or  $\ell ip^n(X,\alpha)$ . Suppose A is any of the algebras  $D^{\infty}(X)$ ,  $Lip^{\infty}(X,\alpha)$  or  $\ell ip^{\infty}(X,\alpha)$  and  $(p_n)_n$  is the corresponding

sequence of algebraic norms defined on  $D^n(X)$ ,  $Lip^n(X,\alpha)$  or  $lip^n(X,\alpha)$ , respectively, which was defined in the Definitions 2.10 and 2.20. We endow A with the metric topology defined by the sequence  $(p_n)_n$  and denote it by  $(A,(p_n))$ . Then we get the following result, which can be found in [12] or [27]. Moreover, by the Carpenter's Theorem there exists a unique topology for each of these algebras as a Fréchet algebra.

**Theorem 4.8** If X is a perfect compact plane set such that  $D^n(X)$  is complete, then  $(A,(p_n))$  is a Fréchet function algebra on X which is not a Banach algebra.

**Theorem 4.9** If X is a uniformly regular space then

$$D^{\infty}(X) = Lip^{\infty}(X, 1) = \ell ip^{\infty}(X, \alpha) = Lip^{\infty}(X, \alpha).$$

**Proof.** If  $f \in D^{n+1}(X)$  then  $f^{(k)} \in D^1(X)$  for  $0 \le k \le n$ . Since X is uniformly regular,  $D^1(X)$  is complete and so there exists a constant M such that for all  $z, w \in X$  and for each  $k \ (0 \le k \le n)$ ,

$$|f^{(k)}(z) - f^{(k)}(w)| \le M|z - w| ||f^{(k+1)}||_X$$
.

Hence for all  $k(0 \le k \le n)$ ,

$$p_1(f^{(k)}) = \sup_{\substack{z,w \in X \\ z \neq w}} \left| \frac{f^{(k)}(z) - f^{(k)}(w)|}{z - w} \right| \le M \parallel f^{(k+1)} \parallel_X < \infty,$$

and so for every  $k(0 \le k \le n)$ ,  $f^{(k)} \in Lip(X,1)$ . Thus  $f \in Lip^n(X,1)$  and it follows that  $D^{n+1}(X) \subseteq Lip^n(X,1)$ . It is easy to check that

$$Lip^n(X,1) \subseteq lip^n(X,\alpha) \subseteq Lip^n(X,\alpha) \subseteq D^n(X),$$

and so

$$D^{\infty}(X) \subseteq Lip^{\infty}(X,1) \subseteq \ell ip^{\infty}(X,\alpha) \subseteq Lip^{\infty}(X,\alpha) \subseteq D^{\infty}(X).$$

Hence the result follows.

Corollary 4.10 If X is uniformly regular then the metric topologies of  $D^{\infty}(X)$ ,  $Lip^{\infty}(X,1)$  and  $Lip^{\infty}(X,\alpha)$  are equivalent.

**Proof.** It is immediate by the Carpenter's Theorem [5].

Now we introduce the Fréchet function algebras  $FLip(X,\alpha)$  and  $Flip(X,\alpha)$  and some important subalgebras of them, where X is a hemicompact metric space. We also extend some known results and theorems about the Banach function algebras  $Lip(X,\alpha)$  and  $lip(X,\alpha)$  to these Fréchet function algebras.

Let (X,d) be an arbitrary metric space (not necessarily compact) and  $0 < \alpha \le 1$ . Then, as before, we take  $\text{Lip}(X,\alpha)$  as the space of all bounded Lipschitz functions of order  $\alpha$  on X, which is a Banach algebra under the same norm as defined in the Definition 2.3.

Now we define a new topology  $\tau$  on  $Lip(X,\alpha)$  under which it is an LMC-algebra so that when X is a hemicompact metric space,  $(Lip(X,\alpha),\tau)$  is metrizable. For a compact subset K of X we define

$$p_{\kappa}(f) = ||f||_{K} + p_{\alpha}(f|_{K}) \quad (f \in Lip(X, \alpha)).$$

Clearly  $p_{\kappa}$  is a submultiplicative seminorm on  $Lip(X,\alpha)$  so that the family  $(p_{\kappa})_K$  is a separating family of seminorms. Let  $\tau$  be the topology on  $Lip(X,\alpha)$  which is defined by the family  $(p_{\kappa})_{\kappa}$  of seminorms. Then we have the following interesting result, which can be found in [27;2.4.1]

**Theorem 4.11** Let (X,d) be a hemicompact metric space. Then  $(Lip(X,\alpha),\tau)$  is a Fréchet algebra if and only if X is compact (in this case  $(Lip(X,\alpha),\tau)$  is indeed a Banach function algebra).

**Definition 4.12** For the hemicompact metric space (X, d) the completion of  $Lip(X, \alpha)$  with respect to  $\tau$  is denoted by  $FLip(X, \alpha)$  and it is called the Fréchet Lipschitz algebra of order  $\alpha$  on X.

We can characterize the elements of  $FLip(X,\alpha)$  as the continuous functions f (not necessarily bounded) such that  $f|_K \in Lip(K,\alpha)$  for each compact subset K of X. To prove this result one can refer to [27; Theorem 2.4.3], which is stated partly in the following.

**Theorem 4.13** Let (X,d) be an arbitrary metric space and let A be the algebra of all continuous functions f on X for which

$$p_{K}(f) = ||f||_{K} + p_{\alpha}(f|_{K})$$

is finite for each compact subset K of X. We endow A with the topology defined by the family  $(p_{\kappa})_K$  of seminorms. Then

- (i)  $Lip(X, \alpha)$  is dense in A with respect to this topology.
- (ii)  $(A,(p_{\kappa}))$  is a Fréchet algebra if and only if X is hemicompact.
- (iii)  $(A,(p_{\kappa}))$  is a Banach algebra if and only if X is compact.

Remark 4.14 Note that Parts (i) and (iii) of the above theorem also imply Theorem 4.11. From now on we assume that X is a hemicompact metric space and  $(K_n)$  is an admissible exhaustion of X. The theorem shows that

$$FLip(X,\alpha) = \{ f \in C(X) : f|_{K_n} \in Lip(K_n,\alpha), n \in \mathbb{N} \}.$$

Here it is not required to assume that  $f \in C(X)$ . Because X is a k-space [9] and so if f satisfies the second condition, which implies the continuity of f on each  $K_n$ , then it is continuous on X.

Another interesting property is that the algebra  $FLip(X, \alpha)$  is dense in C(X) with respect to the compact-open topology. To see this we argue as follows:

It is clear that for each n,  $FLip(X,\alpha)|_{K_n} = Lip(K_n,\alpha)$  and  $Lip(K_n,\alpha)$  is dense in  $C(K_n)$ . So if  $f \in C(X)$  and  $U = \{g \in C(X) : ||g-f||_{K_n} < \varepsilon\}$  is a

neighbourhood of f in C(X) for  $n \in \mathbb{N}$ , then there exists  $g \in Lip(K_n, \alpha)$  with  $||f|_{K_n} - g||_{K_n} < \varepsilon$ . Now we can extend g to a  $\tilde{g} \in Lip(X, \alpha) \subseteq FLip(X, \alpha)$  and conclude that  $\tilde{g} \in U$ .

Now as in the compact case we introduce some interesting subalgebras of  $FLip(X, \alpha)$ . As before X is assumed to be a hemicompact metric space with the admissible exhaustion  $(K_n)$  and  $0 < \alpha \le 1$ .

**Definition 4.15** For  $0 < \alpha < 1$  we denote the set of all  $f \in C(X)$  for which  $f|_{K_n} \in lip(K_n, \alpha)$  by  $Flip(X, \alpha)$ .

Clearly  $Flip(X, \alpha)$  is a subalgebra of  $FLip(X, \alpha)$  and it is easy to see that  $Flip(X, \alpha)$  is indeed a closed subalgebra of  $FLip(X, \alpha)$ . Obviously  $Flip(X, \alpha)$  is also a Frechet function algebra (Ff-algebra) on X.

The inclusion  $Lip(K_n,1) \subseteq lip(K_n,\alpha)$ , for each n, implies easily that  $FLip(X,1) \subseteq Flip(X,\alpha)$ . Likewise the compact case we can show that FLip(X,1) is dense in  $Flip(X,\alpha)$ . More generally, we have the following result, which is found in [27].

**Theorem 4.16** The algebra Lip(X,1) and hence FLip(X,1) is dense in  $Flip(X,\alpha)$ , for each  $\alpha < 1$ .

**Proof.**Clearly  $Lip(X,1) \subseteq FLip(X,1) \subseteq Flip(X,\alpha)$ . Let  $f \in Flip(X,\alpha)$  and  $U = \{g \in Flip(X,\alpha) : p_{\kappa_n}(g-f) < \varepsilon\}$  for  $n \in \mathbb{N}$  be a neighbourhood of f in  $Flip(X,\alpha)$ . Then  $f|_{K_n} \in lip(K_n,\alpha) = \overline{Lip(K_n,1)}$ . Hence there exists  $g \in Lip(K_n,1)$  with  $||f-f|_{K_n}||_{K_n} + p_{\alpha}(g-f|_{K_n}) < \varepsilon$ . Let  $\tilde{g} \in Lip(X,1)$  be an extension of g to X so that  $\tilde{g}|_{K_n} = g$  and therefore  $\tilde{g} \in U \cap Lip(X,1)$ .

Now we present a Fréchet algebra version of the Hedberg's theorem for the Ff-algebra  $Flip(X,\alpha)$ . It is interesting to see that we can also prove the density of FLip(X,1) in  $Flip(X,\alpha)$  by using this theorem.

We recall that for  $a \in X$  and  $\delta > 0$ 

$$S_a(\delta) = \{x \in X : d(x, a) = \delta\}$$
  
$$B_a(\delta) = \{x \in X : d(x, a) < \delta\}$$

Theorem 4.17 [27] Let (X,d) be a hemicompact metric space,  $0 < \alpha < 1$  and A be a self-adjoint subalgebra of  $Flip(X,\alpha)$ , which separates the points of X and contains the constants. Then A is dense in  $Flip(X,\alpha)$  if for every  $a \in X$  there exist numbers  $M_a$  and  $\delta_a$  such that for every  $\delta \leq \delta_a$  there is an  $f \in A$  with f(a) = 1, f(x) = 0 on  $S_a(\delta)$  and

$$\sup_{\substack{y,z \in B_{\alpha}(\delta) \\ y \neq z}} \frac{|f(y) - f(z)|}{d^{\alpha}(y.z)} < \frac{M_{\alpha}}{\delta^{\alpha}}.$$

**Proof.** It is enough to show that for each compact subset K of X, the subalgebra  $A|_K = \{f|_K : f \in A\}$  of  $lip(K,\alpha)$  is dense in  $lip(K,\alpha)$ . Since  $\overline{A|_{K_n}} = lip(K_n,\alpha)$ , for an arbitrary element  $f \in Flip(X,\alpha)$  and a neighbourhood  $U = \{g \in Flip(X,\alpha) : p_{K_n}(g-f) < \varepsilon\}$  of f, we can choose a  $g \in A$  with  $p_{K_n}(g-f) < \varepsilon$  so that  $g \in U \cap A$ .

Now let  $K \subseteq X$  be compact. We try to verify the hypothesis of the Hedberg's theorem for the subalgebra  $A|_K$  of  $lip(K,\alpha)$ . Clearly  $A|_K$  is a self-adjoint subalgebra of  $lip(K,\alpha)$ , which separates the points of K and contains the constants. For an arbitrary element  $a \in K$ , choose  $M_a$  and  $\delta_a$  as in the hypothesis of the theorem. So for every  $\delta \leq \delta_a$  we can choose  $f \in A$  with f(a) = 1, f(x) = 0 on  $S_a(\delta)$  and

$$\sup_{\substack{y,z\in B_a(\delta)\\y\neq z}}\frac{|f(y)-f(z)|}{d^{\alpha}(y,z)}<\frac{M_a}{\delta^{\alpha}}.$$

Then  $f|_K \in A|_K$  has the desired properties, i.e.,  $f|_K(a) = 1$ ,  $f|_K(x) = 0$  on  $S_a(\delta) \cap K = \{x \in K : d(x,a) = \delta\}$  and  $\sup \frac{|f(y)-f(z)|}{d^a(y,z)} < \frac{M_a}{\delta^a}$ , where the supremum is taken over all distinct elements

 $y, z \in B_a(\delta) \cap K = \{x \in K : d(x, a) \leq \delta\}$ . Therefore,  $\overline{A|_K} = \ell i p(K, \alpha)$  by the Hedberg's theorem in the complex case, and hence the claim is now established.

As an application of this theorem one can prove that  $\overline{FLip(X,1)} = Flip(X,\alpha)$  by taking the function  $f(x) = 1 - \frac{d(x,a)}{\delta}$  for arbitrary  $\delta > 0$  and  $M_a = 2^{1-\alpha}$  in the above theorem.

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Proceedings of the Monthly Colloquium of the Iranian Mathematical Society Vol. 2, pp 105-145 (2004)

# An Empirical Bayes View of Inference in Inverse Gaussian Distribution

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Abstract: The Inverse Gaussian distribution has been around for more than a century. It has been treated by classical methods for estimation, hypothesis testing, developing models for analysis of variance and regression by many authors. Its possible usefulness, when treated by Bayesian and empirical Bayes methods, first came to the attention of the present author about a decade ago. It was inspired by a short communication dealing with it via Bayesian approach Achcar and Rosales (1993). Since then we have been involved in model building for this distribution exploiting the empirical Bayes approach. First, the estimation problems were studied. Later, we dealt with two-way analysis of variance, and finally with regression problems. The analysis of covariance is in progress. In this paper, we review the main points of this continued research. some applications are presented as illustrations.

MSC (2000): 62JXX, 62KXX, 62C12

Keywords: Inverse Gaussian distribution, estimation, regression, Bayes, empirical Bayes, prior, posterior.

# 1. Historical Review Of The Inverse Gaussian Distribution

Robert Brown (1773-1858), one of the greatest botanists of England was the one who began the story of the Inverse Gaussian distribution. Interested in pollens, he found a swimming, dancing motion of pollen particles when these particles were immersed in water. He repeated this experiment with a vide variety of plant pollens, finding a similar motion in every case. Later on, he tried particles from dead plants, fossils, and even mineral specimens of all sorts, in fact virtually everything he could imagine from the soot of London to a fragment from the Sphinx. Brown apparently believed he had discovered a new type of particle, common to all matter, organic and inorganic. Several researchers before him had observed and noted the motion of microscopic organic particles in fluids, but his work led to the realization that it was a physical phenomenon, not a biological one. Whether he himself initially recognized this fact is open to debate. During the rest of the century many researchers conducted experiments and theorized about the nature of the so-called Brownian motion. The law which governs the position of a single particle performing one-dimensional Brownian motion was derived by Bachelier in 1900. It turned out as a normal distribution. Wiener in 1923 amended Bachelier's work by providing a measure on the path space. By Einestion's (1905) work who also derived the normal distribution as the model for Brownian motion, the theory of Brownian motion was firmly launched. Schrodinger (1915) considered the Brownian motion with a positive drift and derived the distribution of the first passage time. This distribution was also obtained by Smouluchowski (1915).

Tweedie (1941) and Wald (1944) encountered the first passage distribution. Tweedie noticed that there is an inverse relationship between

the cumulant - generating function of the time to cover unit distance and the cumulant - generating function of the distance covered in unit time. Tweedie (1945) also observed this type of relationship between the binomial and the negative binomial, and between the poisson and the exponential distribution. He proposed to call them inverse statistical variates. In 1956, he coined the name Inverse Gaussian for the first passage time distribution of the Brownian motion. In 1957, he published a detailed study of this distribution, establishing many of its important statistical properties. A special case of the distribution was given by Wald in 1947, which is an approximation of the sample size distribution in a sequential probability ratio test. Folks and Chhikara (1978) and Chhikara and Folks (1989) are the main sources for the statistical properties of the Inverse Gaussian model. The Inverse Gaussian model has been applied in reliability, marketing, and analysis of experiments. The main references concerning the classical approach are Fries and Bhatacharrya (1983), Chhikara and Gmttman (1982), Banerjee and Bhattacharyya (1979). On the Bayesian front the noteable works are Banerjee and Bhattacharrya (1979), Achcar et al. (1991), and Achcar and Rosales (1992, 1993).

For regression analysis under the Inverse Gaussian model, we refer to Whitmore (1983, 1986) Seto and Iwase (1985), and Chhikara and Folks (1989). Iwase (1989), Woldie and Folks (1994, 1995) and lately Seshadri (1999) are the relevant material.

The present author began his efforts with an enpirical Bayes approach. In 1994, one of his graduate students was engaged in the parameter estimation, Homayun - Aria, (1996). At the same time he himself developed methodology for analysis of variance, Meshkani (1996). Following this line of research, he later developed a methodology for the regression analysis, Meshkani (1999). Currenty he is working on

the analysis of covariance. In this paper, we intend to summarize the findings from these works and show some of its applications in different settings. Some directions for further research will also be mentioned.

# 2. Empirical Bayes estimation of the parameters in an Inverse Gaussian distribution

# 2.1 Introduction

A random variable Y with the probability distribution function

$$f(y|\theta,\lambda) = \left[\frac{\lambda}{2\pi y^3}\right]^{\frac{1}{2}} \exp\left\{\frac{-\lambda(y-\theta)^2}{2y\theta^2}\right\} \qquad y,\lambda,\theta > 0 \qquad (2-1)$$

is called an Inverse Gaussain variate, where  $\theta$  is the mean and  $\lambda$  is the scale parameter. This function will be denoted by  $IG(\theta, \lambda)$ . The  $IG(\theta, \lambda)$  for various values of  $\theta$  and  $\lambda$  provides a rich family of positively skewed densities which are suitable for modelling many types of positive random variables, Figures 2-1 and 2-2. It has the following cumulant - generating function

$$k_x(t) = \frac{\lambda}{\theta} \left[ 1 - \left( 1 - \frac{2\theta^2 t}{\lambda} \right)^{\frac{1}{2}} \right], \qquad t < \frac{\lambda}{2\theta^2}$$
 (2-2)

which provides the moments  $E(x) = \theta$ ,  $var(x) = \frac{\theta^3}{\lambda}$ , and generally

$$K_r = 1 \cdot 3 \cdot 5 \cdots (2r - 3)\theta^{2r - 1}\lambda^{r - 1}, \quad r \ge 2$$

The measures of skewness and kurtosis are  $\sqrt{\beta_1} = 3\sqrt{\frac{\theta}{\lambda}}$  and  $\beta_2 = 15(\frac{\theta}{\lambda}) + 3$ , which both are positive. For a normal variate,  $V \sim N(v, \sigma^2)$ , one has the cumulant - generating function

$$k_v(t) = vt - \frac{1}{2}\sigma^2 t^2, \quad t \in \mathbb{R}$$
 (2-3)

since (2-2) and (2-3) are inverse functions of each orther hence the Inverse Gaussian name for (2-1). Moreover, there are striking analogies between the sampling distributions for  $IG(\theta, \lambda)$  and those for  $N(v, \sigma^2)$ , see Chhikara and Folks (1989). The Inverse Gaussian distribution belongs to the exponential family.

## 2.2 Estimation

We suppose a random sample of  $Y = (Y_1, \ldots, Y_n)$  is available from the  $IG(\theta, \lambda)$  distribution. The aim is to obtain an estimator for  $(\theta, \lambda)$  based on this sample. Most of estimation procedures exploit the likelihood function which is

$$L(\theta, \lambda | \boldsymbol{y}) = \left(\frac{\lambda}{2\pi}\right)^{\frac{n}{2}} \left(\prod_{i=1}^{n} y_i^{-\frac{3}{2}}\right) \exp\left\{-\frac{\lambda}{2} \sum_{i=1}^{n} \frac{(y_i - \theta)^2}{y_i \theta^2}\right\}, \quad \theta, \lambda > 0$$

$$(2 - 4)$$

The maximum likelihood estimators are based on maximization of (2-4) with respect to  $\theta$  and  $\lambda$ , simultaneously. This approach yields:

$$\hat{\theta}_{ML} = \bar{Y} = \frac{1}{n} \sum_{i=1}^{n} Y_i, \quad \hat{\lambda}_{ML} = \frac{n}{\sum_{i=1}^{n}} \left( \frac{1}{Y_i} - \frac{1}{\bar{Y}} \right)$$
 (2-5)

If we define the sample harmonic mean as

$$\bar{Y}_H = \frac{1}{\left[\frac{1}{n}\sum_{i=1}^n \frac{1}{Y_i}\right]}$$

Them  $\hat{\lambda}_{ML}$  in (2-5) is non-negative, since  $\bar{Y}_H \leq \bar{Y}$ . Among many variants of the Bayes estimators relative to various priors, we report a case where it is assumed that  $\theta$  and  $\lambda$  are independent and exponentially distributed, viz.

$$\theta \sim \exp(a)$$
 and  $\lambda \sim \exp(b)$ 

Then the joint posterior of  $(\theta, \lambda)$  will be

$$P(\theta, \lambda | y) \propto \lambda^{\frac{n}{2}} \exp\left\{-\frac{\lambda v}{2} - \frac{n\lambda v\bar{Y}}{2} \left(\theta - \frac{1}{\lambda}\right)^2 - a\theta - b\lambda\right\}$$
 (2-6)

with  $v = \sum_{i=1}^{n} \left(\frac{1}{y_i} - \frac{1}{\bar{y}}\right)$ . Upon integration, we can obtain the marginal posteriors of  $P(\theta|\mathbf{y})$  and  $P(\lambda|\mathbf{y})$ . From (2-6), the Bayes estimators of  $\theta$  and  $\lambda$  relative to the squared error loss are the posterior means of  $P(\theta|\mathbf{y})$  and  $P(\lambda|\mathbf{y})$ . That is,

$$\hat{\theta}_B = \hat{\theta}(a, b) = C\Gamma\left(\frac{n}{2} + 1\right) \int_0^\infty \frac{\theta e^{-a\theta}}{\frac{v}{2} + \frac{n\overline{Y}}{2}(\theta - \frac{1}{\overline{Y}})^2 + b} d\theta \qquad (2 - 7)$$

where  $C = \left(\frac{1}{2\pi}\right)^{\frac{n}{2}} \prod_{i=1}^{n} Y_i^{-\frac{3}{2}}$ . Likewise,

$$\hat{\lambda}_B = \hat{\lambda}_B(a,b) =$$

$$\sqrt{2\pi}Ce^{-\frac{a}{\bar{Y}}} \int_{0}^{\frac{a}{n}} \left\{ \frac{\lambda^{\frac{n}{2}+1}}{\sqrt{n\lambda\bar{Y}}} \exp\{-\lambda(\frac{v}{2}+b) - \frac{a^{2}}{2n\lambda\bar{Y}}\} \Phi\left(\frac{n\lambda - a}{\sqrt{n\lambda\bar{Y}}}\right) \right\} d\lambda + \int_{\frac{a}{n}}^{\infty} \left\{ \frac{\lambda^{\frac{n}{2}+1}}{\sqrt{n\lambda\bar{Y}}} \exp\{-\lambda(\frac{v}{2}+b) - \frac{a^{2}}{2n\lambda\bar{Y}}\} \Phi\left(\frac{a-n\lambda}{\sqrt{n\lambda\bar{Y}}}\right) \right\} d\lambda \qquad (2-8)$$

The Bayes estimators (2-7) and (2-8) have to be evaluated by numerical methods. Details of these derivations can be found in Homayun-Aria (1996). By a similar operation, we find the posterior variances  $var(\theta|y)$  and  $var(\lambda|y)$ . The Empirical Bayes estimators are derived from the Bayes estimators by replacing the prior distribution by its estimated version. In case of a parametric prior, this strategy amounts to estimation of the parameters of the prior from the observed data and then using them in the Bayes estimators. This estamation can be done employing either the method of maximum likelihood (ML) or the method moments (MM). For the method of ML, we first obtain the marginal distribution of the sample values under the above exponential priors for  $(\theta, \lambda)$ :

$$m(y|a,b) = \int_0^\infty \int_0^\infty L(\theta,\lambda|y)P(\theta)P(\lambda) d\theta d\lambda$$

$$\frac{ab}{(2\pi)^{\frac{n}{2}}\dot{y}} \int_0^\infty \int_0^\infty \exp\left\{\frac{-\lambda}{2} \sum_{i=1}^n \frac{(y_i - \theta)^2}{y_i \theta^2} - a\theta - b\lambda\right\} d\theta d\lambda$$

Where  $\dot{y} \equiv \prod_{i=1}^n y_i^{\frac{3}{2}}$ ,

$$m(\boldsymbol{y}|a,b) = \frac{2ab\bar{y}_H}{(2\pi)^{\frac{n}{2}}\dot{y}(2b\bar{y}_H + n)} \int_0^\infty \frac{\theta^2 e^{-a\theta}}{\theta^2 - r\theta + s} d\theta$$

with 
$$r = \frac{2n\vec{y}_H}{2b\vec{y}_H + n} > 0, S = \frac{r\vec{y}}{2} > 0.$$

After evaluation of this integral, as I(a,b) say, we obtain

$$m(y|a,b) = \frac{2ab\bar{y}_H}{(2\pi)^{\frac{n}{2}}\dot{y}(2b\bar{y}_H + n)}I(a,b)$$

To obtain the ML estimates of a and b, we have to maximize m(y|a,b) with respect to a, and b. Suppose  $\hat{a}_{ML}$  and  $\hat{b}_{ML}$  are the solutions. Then, we obtain the empirical Bayes estimators as

$$\hat{\theta}_{EBML} = \hat{\theta}_B(\hat{a}_{ML}, \hat{b}_{ML}) \tag{2-9}$$

$$\hat{\lambda}_{EBML} = \hat{\lambda}_B(\hat{a}_{ML}, \hat{b}_{ML}) \tag{2-10}$$

whose variances are estimated by  $\operatorname{var}(\hat{\theta}_{EBML}|\boldsymbol{y})$  and  $\operatorname{var}(\hat{\lambda}_{EBML}|\boldsymbol{y})$ , respectively. The above results do not have closed from and are obtainable only by numerical integrations. To obtain an explicit solution, we may use the method of moments to estimate a and b. This procedure leads to

$$\hat{a}_{MM} = \bar{y}$$
 and  $\hat{b}_{MM} = \frac{\left[\frac{1}{n} \sum_{i=1}^{n} y_i^2 - \bar{y}^2\right]}{2\bar{y}^3}$ .

These estimates provide an alternative version of the empirical Bayes estimators of  $\theta$  and  $\lambda, i, e$ .,

$$\hat{\theta}_{EBMM} = \hat{\theta}_B(\hat{a}_{MM}, \hat{b}_{MM}) \tag{2-11}$$

$$\hat{\lambda}_{EBMM} = \hat{\lambda}_B(\hat{a}_{MM}, \hat{b}_{MM}) \tag{2-12}$$

with corresponding variances  $var(\hat{\theta}_{EBMM}|y)$  and  $var(\hat{\lambda}_{EBML}|y)$ .

# 3. Linear Models under the Inverse Gaussian Model

In scientific activities one encounters problems which are broadly classified as comparing the means of a certain variable among several subpopulations. Examples are the average yields from different process formulas, average number of defectives from various machines, etc. The Technical name for these problems in statistics is the Analysis of Variance (ANOVA). Occasionalyy, one is interested in finding the functional relation betwen a dependent (response) variable and some independent (explanatory) variables. These types of problems are called regression analysis. Yet a third type of problem is encountered when one has the regression and ANOVA at the same time which is referred to Analysis of Covariance (ANCOVA). These three type of problems are subsumed under the general title of linear models, whenever the functional relations imvolve parameters of the first degree. A number of authors have tried to develope statistical linear models for the Inverse Gaussian model akin to those under the normal model. However, the results are limited in the realm of classical approach. But Bayesian and Empirical Bayes approaches seem promising. we report some results obtained for ANOVA and regression.

# 4. Empirical Bayes Analysis of Variance

### 4.1 Introduction

There are many types of experimental setups in science and engineering where the normal theory is inappropriate for the analysis of factorial

experiments. One important class is related to the highly skewed nature of the data which cannot be removed by the usual transformations. Alternatively, the Inverse Gaussian family of distributions are flexible enough to provide a suitable model for these types of data. Tweedie (1957 a,b) pioneered work in providing an analogue to analysis of variance for nested classifications concerning observations from an Inverse Gaussian model. Despite the quite striking resemblance between normal analysis of variance and what he called, the Inverse Gaussian analysis of reciprocals, in one-way or nested classifications, the possibilities of developing analogous results for other classifications appeared to be limited, Folks and Chhikara (1978), and Chkikara and Folks (1989). However, Shuster and Muira (1972) succeeded in providing tests for balanced two - way classifications. Their approach has the disadvantage that it requires many observations in each cell. Such a requirement is hard to fulfil in most experiments. Fries and Bhattacharyya (1983) treated the analysis of two-factor experiment with no interaction and obtained explicit solutions to the likelihood equations. They also proved asymptotic consistency and normality of their estimators. A few authors have contributed to the Bayesian analysis of the Inverse Gaussian distribution. Achcar and Rosales (1992,1993) are the only records in print for Bayesian analyis of two-factor experiments under an Inverse Gaussian model. They actually follow the approach presented in Fries and Bhattachatyya (1983), assuming a non-informative prior density. Consequently, as one expects, because of heavy reliance on the likelihood, their results are not much different from those obtained by the maximum likelihood method of estimation.

In this paper, we present an empirical Bayes analysis of two-factor experiments under an Inverse Gaussian model. A real-life example previously analyzed by Shuster and Muira (1972), and later by Achcar and Rosales

(1993), is reworked.

### 4.2 THE MODEL

Consider an experiment with two factors, factor A with I levels indexed by i, and factor B with J levels, numbered by j, with each treatment combination being repeated n times. Observations from this experiment, denoted by  $y_{ijk}$ , are assumed to follow an Inverse Gaussian model,  $IG(\theta_{ij}, \lambda)$ ,

$$Y_{ijk} \sim IG(\theta_{ij}, \lambda), \quad i = 1, \ldots, I \quad j = 1, \ldots, J \quad k = 1, \ldots, n.$$

For each i,j, the random variables  $Y_{ijk}$  are iid with mean  $\theta_{ij}$  and shape parameter  $\lambda$ . The two-parameter Inverse Gaussian density is

$$f(y_{ijk}; \theta_{ij}, \lambda) = \left\{ \lambda/(2\pi y_{ijk}^3) \right\}^{\frac{1}{2}} \exp\left\{ -\lambda (y_{ijk} - \theta_{ij})^2 / 2y_{ijk} \theta_{ij}^2 \right\},$$

$$y_{ijk} > 0, \quad \theta_{ij} > 0, \quad \lambda > 0, \quad i = 1, ..., I \quad j = 1, ..., J, \quad k = 1, ..., n.$$

$$(4-1)$$

In a two-factor experiment with interaction, each cell mean is assumed to be inversely proportional to the drift, while the drift is considered as the sum of factor main effects  $(\alpha, \beta)$  and their interaction $(\gamma)$ . Thus, it is assumed that

$$\theta_{ij}^{-1} = \mu + \alpha_i + \beta_j + \gamma_{ij}, \quad i = 1, ..., I, \qquad j = 1, ..., J,$$

$$(4-2)$$

$$\sum_{i=1}^{I} \alpha_i = \sum_{j=1}^{J} \beta_j = 0, \quad \sum_{i=1}^{I} \gamma_{ij} \sum_{j=1}^{J} \gamma_{ij} = 0.$$
 (4-3)

Here,  $\mu$  denotes the reciprocal of each cell mean when there is no drift. To incorporate the constraints (4-3) in the model, we can define the  $IJ \times 1$  parameter vector  $\Phi$  as

$$\Phi = [\mu \mid \alpha_{1}, \dots, \alpha_{I-1} \mid \beta_{1}, \dots, \beta_{J-1} \mid \gamma_{11}, \dots, \gamma_{1,J-1} \mid \dots | \gamma_{I-1,J}, \dots, \gamma_{I-1,J-1} \mid] = [\mu, \alpha, \beta \gamma_{1}, \dots \gamma_{i}, \dots \gamma_{I-1}].$$
(4-4)

Then, the likelihood for the whole experiment can be written as

$$L(\Phi, \lambda \mid y) \propto \lambda^{nIJ/2} \exp\{-\lambda [R_{+++} - 2n\Phi'd + n\Phi'M\Phi]/2\}. \quad (4-5)$$

In (4-5), the convention used by Fries and Bhattacharyya (1983) has been utilized, where we have set

$$y = [y_{111}, \dots, y_{11k}, \dots, y_{11n}, \dots, y_{ij1}, \dots, y_{ijk}, \dots, y_{ijn}, \dots, y_{IJn}],$$

$$y_{ijk}, \dots, y_{ijn}, \dots, y_{IJ1}, \dots, y_{IJk}, \dots, y_{IJn}],$$

$$\theta_{ij}^{-1} = \mu + \alpha_i + \beta_j + \gamma_{ij} = X'_{ij}\Phi,$$

$$X' = (X_{11}, X_{12}, \dots, X_{IJ}) = design \quad matrix,$$

$$D = \text{diag} \{y_{11}, y_{12}, \dots, y_{IJ}\}, \qquad y_{ij} = \sum_{k=1}^{n} y_{ijk}/n,$$

$$M = X'DX, \qquad d = \sum_{i=1}^{I} \sum_{j=1}^{J} X_{ij},$$

$$R_{ijk} = y_{ijk}^{-1}. \qquad (4-6)$$

Summing over and index is shown by a plus sign while averaging is denoted by a dot. Thus, we shall use  $R_{i++}$ ,  $R_{+j+}$ ,  $R_{ij+}$ ,  $R_{+++}$ ,  $R_{i..}$ ,  $R_{.j.}$ ,  $R_{ij}$ , and  $R_{...}$  as sums and averages, respectively.

We intend to use a conjugate prior for  $\lambda$  and  $\Phi$ . The following priors have been proposed, see Chhikara and Folks(1989), and Banerjee and Bhattacharyya (1979). The prior for  $\lambda$  is chosen from the gamma family and given  $\lambda$ , a normal prior is assumed or  $\Phi$ . Thus,

$$\pi(\lambda) \propto \lambda^{a-1} \exp\{-b\lambda/2\}, \qquad \lambda, a, b > 0,$$
 (4-7)

and, given  $\lambda$ , elements of  $\Phi$  are considered independent with either of the following two priors.

Case 1. Unrestricted parameter space.

In the unrestricted case, the prior distribution for  $\Phi$  is

$$\Phi\Big|_{\lambda} \sim N(\eta, \lambda^{-1}\Delta)$$
 (4-8)

with

$$\eta = [\eta_1, \dots, \eta_{IJ}], \qquad \Delta = \operatorname{diag}\{\delta_1^2, \dots, \delta_{IJ}^2\}.$$

Then, the posterior is

$$q_1(\Phi, \lambda | y) \propto |\Psi|^{\frac{-1}{2}} \lambda^{\nu-1} \exp\{-\lambda [Q_1(\eta) + Q_2(\Phi)]/2\}$$
 (4-9)

where

$$\Psi = (n\Delta M + \Delta^{-1})^{-1},$$
 $\nu = a + (n+1)IJ/2,$ 
 $Q_1(\eta) = R_{+++} + b + \eta'\Delta^{-1}\eta - \eta^{*'}\Psi^{-1}\eta^{*},$ 
 $\eta^* = (n\Delta M + I)^{-1}(n\Delta d + \eta),$ 
 $Q_2(\Phi) = (\Phi - \eta^*)'\Psi^{-1}(\Phi - \eta^*).$ 

It is evident from (4-9) that

$$\pi(\lambda|\mathbf{y}) \propto \lambda^{k-1} \exp\{-\lambda Q_1(\mathbf{\eta})/2\}, \quad \lambda > 0,$$
 (4-10)  
 $\kappa = (nIJ + 2a)/2, \qquad Q_1(\mathbf{\eta}) > 0,$ 

and

$$q_1(\Phi|\lambda, \boldsymbol{y}) \propto |\lambda^{-1}\Psi|^{-1/2} \exp\{-\lambda Q_2(\Phi)/2\},$$

that is, conjugacy holds. Therefore, we can write

$$q_{1}(\Phi \mid y) \propto \frac{1}{[Q_{1}(\eta) + Q_{2}(\Phi)]^{\nu}} \qquad \Phi \in \mathcal{R}^{IJ},$$

$$\propto \left[1 + \frac{(\Phi - \eta^{*})' \sum^{-1} (\Phi - \eta^{*})}{2a + nIJ}\right]^{-\frac{(2a + nIJ) + IJ}{2}} \qquad (4 - 11)$$

with

$$\sum = \frac{Q_1(\eta)\Psi}{2a+nIJ}$$

which is a multivariate T-type distribution, with 2a + nIJ > 2 degrees of reedom.

## Case 2. Restricted parameter space.

Strictly speaking, one should have  $\phi_1 = \mu > 0$ . This restriction on  $\phi_1$  is observed in the prior assigned to  $\phi_1$ . Thus, a normal distribution truncated at zero is considered for  $\phi_1$ , which has density

$$q_2(\phi_1 \mid \lambda) \propto \lambda^{1/2} [\delta_1 \mathcal{N}(\lambda^{1/2} \eta_1 / \delta_1)]^{-1} \exp\{-\lambda (\phi_1 - \eta_1)^2 / 2\delta_1^2\}, \qquad \phi_1 > 0$$

$$(4 - 12)$$

and the remaining is as in case 1. In (4-12),  $\mathcal{N}(.)$  is the standard normal distribution function. The restriction imposed on  $\phi_1$  results in a posterior proportional to (4-9).

# 4.3 Bayes Estimates

In case 1, from (4-10) we have

$$E(\lambda^{m}|\mathbf{y}) = [2/Q_{1}(\eta)]^{m} [\Gamma(\kappa+m)/\Gamma(\kappa)]$$
 (4-13)

which provides the Bayes estimate of  $\lambda$  relative to the squared error loss:

$$\lambda_{B1} = E(\lambda \mid y) = nIJ + 2a)/Q_1(\eta),$$
 (4-14)

$$V_{\lambda} = Var(\lambda|y) = 2(2a + nIJ)/[Q_1(\eta)]^2.$$
 (4-15)

Upon using (3.11), we arrive at

$$\Phi_{B1} = E(\Phi|y) = \eta^* = (n\Delta M + I)^{-1}(n\Delta d + \eta)$$
 (4-16)

$$V_{B1} = Var(\Phi|y) = \frac{Q_1(\eta)\Psi}{2(a-1) + IJ}$$
 (4-17)

For case 2, the posterior moments of  $\lambda$  remain unchanged from those for case 1. However, for  $\Phi$ , the restriction on  $\phi_1$  renders results on  $\Phi$  different from those in (4-16) and (4-17) for case 1. To this end, let  $\Phi$ ,  $\eta$  and  $\Psi$  be partitioned as

$$\eta = \begin{bmatrix} \eta_1 \\ \eta^{(2)} \end{bmatrix}, \eta^* = \begin{bmatrix} \eta_1^* \\ \eta^{*(2)} \end{bmatrix}, \Phi = \begin{bmatrix} \phi_1 \\ \Phi^{(2)} \end{bmatrix}, \Psi = \begin{bmatrix} \psi_{11} & \Psi_{12} \\ \Psi_{21} & \Psi_{22} \end{bmatrix}.$$

Then by virtue of the facts  $\Psi_{22.1} = \Psi_{22} - \Psi_{21} \psi_{11}^{-1} \Psi_{12}$  and  $|\Psi| = \psi_{11} |\Psi_{22.1}|$ ,  $Q_2(\Phi)$  in (4-9) can be written as

$$Q_2(\Phi) = \psi_{11}^{-1} (\phi_1 - \eta_1^*)^2 + (\Phi^{(2)} - \eta_{2.1}^*)' \Psi_{22.1}^{-1} (\Phi^{(2)} - \eta_{2.1}^*)$$
(4.18)

where

$$\eta_{2.1}^* = E(\Phi^{(2)}|\phi_1, \lambda, y) = \eta^{*(2)} + \psi_{11}^{-1}(\phi_1 - \eta_1)\Psi_{21}$$
(4 - 19)

and

$$Var(\Phi^{(2)}|\phi_1,\lambda,y) = \lambda^{-1}\Psi_{22.1}.$$
 (4.20)

In (4-19) , we take successive expectations with respect to  $\phi_1$  and  $\lambda$ , noting that

$$q_2(\phi_1|\lambda, y) \propto |\lambda^{-1}\psi_{11}|^{1/2} \exp\{-\lambda \psi_{11}^{-1}(\phi_1 - \eta_1^*)^2/2\}, \qquad \phi_1 > 0.$$

which is a normal density truncated at zero. For this distribution,

$$E(\phi_1|\lambda, y) = \eta_1^* + \lambda^{-1}\psi_{11})^{1/2}\omega$$

with

$$\omega = \varphi \big[ \eta_1^* (\lambda^{-1} \psi_{11})^{1/2} \big] \mathcal{N} \big[ \eta_1^* / (\lambda^{-1} \psi_{11})^{1/2} \big]$$

where  $\varphi(.)$  is the standard normal density function, and

$$Var(\phi_1|\lambda,y) = \lambda^{-1}\psi_{11}[1-\omega^2] + \eta_1^*\omega(\lambda^{-1}\psi_{11})^{1/2}.$$

These moments, however, are too complicated to be useful for estimation purposes. To simplify them, we observe that  $\omega(x) = \varphi(x)/\mathcal{N}(x)$  is a smooth decreasing function of x. In the literature there are a host of approximations to  $\mathcal{N}(x)$  [Patel and Read (1996, Chapter 3)], which can be used to approximate  $\omega(x)$  with desired precision. Here, we choose to use the simpler one due to Shah(1985),

$$\mathcal{N}(x) = \begin{cases} 0.5 + x(4.4 - x)/10, & 0 \le x \le 2.2, \\ 0.99, & 2.2 \le x < 2.6, \\ 1.00, & x \ge 2.6. \end{cases}$$

Consequently,

$$\omega(x) = \begin{cases} \frac{2(2.2-x)}{5+x(4.4-x)}, & 0 \le x \le 2.2\\ 0, & x > 2.2. \end{cases}$$

In our problem,  $0 \le x \le \infty$  and  $0 \le \omega(x) \le 0.8$ . Thus, we can approximate  $\omega(x)$  by its average value, which is about 0.3. Of course, if one has a better guess of  $x = \eta_1^*/(\lambda^{-1}\psi_{11})^{1/2}$ , a closer approximation could be obtained. Using this approximation, we have

$$E(\phi_1|\lambda, y) = \eta_1^* + 0.3(\lambda^{-1}\psi_{11})^{1/2}$$

$$Var(\phi_1|\lambda, y) = 0.91\lambda^{-1}\psi_{11} + 0.3\eta_1^*(\lambda^{-1}\psi_{11})^{1/2}.$$

Now, we take expectations with respect to the posterior distribution of  $\lambda$ , employing (4-13) and obtain

$$\phi_{1B2} = E(\phi_1|\mathbf{y}) = \eta_1^* + 0.21[\psi_{11}Q_1(\mathbf{\eta})]^{1/2}W$$
 (4-21)

with  $W = \Gamma(\kappa - 0.5)/\Gamma(\kappa) \simeq [2.72(\kappa - 0.5)/\kappa]^{\frac{1}{2}}$ , by Stirling's formula and

$$V_{1B2} = Var(\phi_1|y) = 0.45\psi_{11}Q_1(\eta)/(\kappa-1) + 0.045\psi_{11}Q_1(\eta)[11/(\kappa-1) - W^2]$$

$$+0.21\eta_1^*[\psi_{11}Q_1(\eta)]^{1/2}W.$$
 (4 - 22)

Now, we substitute the posterior moments of  $\phi_1$  into (4-9) to obtain a simpler form as:

$$\Phi_{2B2} = \eta^{*(2)} + \{0.21[\psi_{11}^{-1}Q_1(\eta)]^{1/2}\}\Psi_{21}W \qquad (4-23)$$

$$V_{2B2} = [Q_1(\eta)/2(\kappa - 1)][\Psi_{22} - 0.09(\kappa - 1)W^2\Psi_{21}\psi_{11}^{-1}\Psi_{12}]$$
$$+0.21W\eta_1^*[\psi_{11}^{-1}Q_1(\eta)]^{1/2}\Psi_{21}\psi_{11}^{-1}\Psi_{12}]. \qquad (4-24)$$

The expressions (4-21) - (4-24) provide the Bayes estimates relative to the restricted prior given in (4-12). As long as the prior distribution can be assessed, the above Bayes estimators, could be put into application for two-way classifications. Unfortunately, the situations where these priors can reasonably be assessed are rare. In such cases, we can utilize the empirical Bayes procedure to estimate the prior distributions from the data. By this, we borrow strength from Bayesian logic and objectivity from classical method.

# 4.4 Empirical Bayes estimates

To estimate the prior parameters from the marginal distribution of the observations, one can use any method of estimation. Two more common methods are the method of moments and maximum likelihood. To provide explicit expressions for estimates of  $a,b,\eta$  and  $\Delta$  from the data, y, we shall first use the method of moments. To this end, we have from Chhikara and Folks (1989),

$$V_{ij} = \sum_{k=1}^{n} (Y_{ijk}^{-1} - Y_{ij.}^{-1}) \sim \lambda^{-1} \chi_{n-1}^{2}.$$

Thus,

$$E(V_{ij}) = E[E(V_{ij})|\lambda)] = (n-1)b/2(a-1),$$

$$Var(V_{ij}) = E[Var(V_{ij})|\lambda] + Var[E(V_{ij}|\lambda)]$$
  
=  $(n-1)(n-3+2a)b^2/4(a-1)^2(a-2)$ .

Let

$$V = \sum_{i=1}^{I} \sum_{j=1}^{J} V_{ij} / IJ, \qquad S_V^2 = \sum_{i=1}^{I} \sum_{j=1}^{J} [V_{ij} - V]^2 / (IJ - 1);$$

and let  $C = S_V/V$  be the sample coefficient of variation for  $V_{ij}$ . Then,

$$a_0 = \frac{[2(n-1)C^2 + n - 3]}{[(n-1)C^2 - 2]}, \qquad b_0 = \frac{2(a_0 - 1)V}{(n-1)}$$

which are valid positive estimates of a and b for n > 1, if one has  $C^2 > 2/(n-1)$ , otherwise, take a and b equal to zero, i.e., use a noninformative prior.

For estimation of  $\eta$  and  $\Delta$ , we need 2IJ equations. These are provided by the following considerations. Let  $Z_i, i = 1, \dots, n$ , be iid andom variables distributed as  $IG(\theta, \lambda)$ . The distributional relations between  $Z_i$ ,  $\bar{Z}$  and  $Z_i^{-1}$  have been found in Chhikara and Folks (1989). That is,  $\bar{Z} \sim IG(\theta, n\lambda)$  and  $Z_i^{-1}$  has mean and variance as stated below:

$$Z_i^{-1} \sim [\theta^{-1} + \lambda^{-1} = E(Z_i^{-1}), \quad (\lambda \theta)^{-1} + 2\lambda^{-2} = Var(Z_i^{-1})].$$

Using these results in our model (4-1) - (4-6), for the reciprocals and their means, we have:

$$S_{ij}^{2} = \sum_{qk=1}^{n} (R_{ijk} - R_{ij.})^{2} / n(n-1),$$

$$S_{i}^{2} = \sum_{j=1}^{J} (R_{ij.} - R_{i..})^{2} / J(J-1),$$

$$S_{j}^{2} = \sum_{i=1}^{J} (R_{ij.} - R_{.j.})^{2} / I(I-1),$$

$$S^{2} = \sum_{i=1}^{J} \sum_{j=1}^{J} \sum_{k=1}^{n} (R_{ijk} - R_{ij.})^{2} / nIJ(nIJ-1).$$

We obtain:

$$\eta_1^0 = R_{...} - b_0/2(a_0 - 1)$$

$$\begin{split} \delta_{1,0}^2 &= 2(a_0-1)S^2/b_0 - \eta_1^0/nIj - b_0[2(a_0-1)+nIJ](a_0-1)(a_0-2) \\ \text{For} \qquad i = 1, \cdots, I-1, \\ \eta_{i+1}^0 &= R_{i..} - \eta_1^0 - b_0/2(a_0-1) = R_{i..} - R_{...}, \\ \delta_{i+1,0}^2 &= 2(a_0-1)S_i^2/b_0 - (\eta_1^0 + \eta_{i+1}^0)/nJ - \delta_1^0 - b_0[2(a_0-1)+nJ]/2nJ(a_0-1)(a_0-2). \\ \text{For} \qquad j = 1, \cdots, J-1, \\ \eta_{I+j}^0 &= R_{.j.} - \eta_1^0 - b_0/2(a_0-1) = R_{.j.} - R_{...}, \\ \delta_{I+j,0}^2 &= 2(a_0-1)S_j^2/b_0 - (\eta_1^0 + \eta_{I+j}^0)/nI - \delta_1^0 \end{split}$$

Finally, for  $i = 1, \dots, I-1$ , and  $j = 1, \dots, J-1$ ,

 $-b_0[2(a_0-1)+nI]2nI(a_0-1)(a_0-2).$ 

$$\begin{split} \eta_{I+i(J-1)+j}^2 &= R_{ij.} - \eta_1^0 - \eta_{i+1}^0 - \eta_{I+j}^0 - b_0/2(a_0 - 1) \\ &= R_{ij.} - R_{i..} - R_{.j.} + R_{...}, \\ \delta_{I+i(J-1)+j,0}^2 &= 2(a_0 - 1)S_{ij}^2/b_0 - [\eta_1^0 + \eta_{i+1}^0 + \eta_{I+j}^0 + \eta_{I+i(J-1)+j}^0]/n \\ &- (\delta_{1,0}^2 + \delta_{i+1,0}^2 + \delta_{I+j,0}^2) - b_0[2(a_0 - 1) + n]/2n(a_0 - 1)(a_0 - 2) \end{split}$$

which provide

$$m{\eta}^0 = [\eta_1^0, \eta_2^0, \dots, \eta_{IJ}^0],$$
  $\Delta^0 = ext{diag}\{\delta_{1,0}^2, \delta_{2,0}^2, \dots, \delta_{IJ,0}^2\}.$ 

Now, we subtitute these estimates into (4-14) and (4-16) to obtain the empirical Bayes estimate relative to unrestricted prior distribution. This gives us

$$\lambda_{EB1} = (nIJ + 2a_0)/Q_1(\eta^0), \qquad (4-25)$$

$$\Phi_{EB1} = (n\Delta^0 M + I)^{-1} (n\Delta^0 d + \eta^0) = \eta^{*0}.$$
 (4.26)

Posterior variances are estimated by

$$Var(\lambda|y) = 2(2a_0 + nIJ)/[Q_1\eta^0]^2$$
 (4.27)

$$Var(\Phi|\mathbf{y}) = [Q_1(\eta^0)/2(\kappa-1)]\Psi^0, \qquad \Psi^0 = [n\Delta^0 M + (\Delta^0)^{-1}]^{-1}.$$
 (4.28)

In case 2, the only difference in prior is that  $\phi_1$  has a truncated normal prior. Accordingly, we should alter the posterior and the marginal moments for differences in moments of  $\phi_1$ . In this case,

$$E(\phi|\lambda) = \eta_1 + 0.3\lambda^{-1/2}\delta_1,$$

$$Var(\phi_1|\lambda) = 0.91\lambda_1^{-1}\delta_1^2 + 0.3\eta_1\lambda^{-1/2}\delta_1.$$

These differ from the respective moments of  $\phi_1$  in case 1. To account for this difference, the previous moment equations should be modifed accordingly. Omitting the details which can be found in Meshkani (1996), we shall give the final results.

Let 
$$k(a_0) = 0.21w(a_0)b_0^{1/2}$$
  $w(a_0) = \Gamma(a_0 - 0.5)/\Gamma(a_0)$ 

$$m_0 = [b_0^2/4nIJ(a_0-1)^2(a_0-2)][2(a_0-1)+nIJ]$$

Then,

$$\eta_1 = R_{...} - b_0/2(a_0 - 1) - k(a_0)\delta_1$$

$$S^2 = m_0 + [b_0/2(a_0 - 1)]\delta_1^2 + [b_0/2nIJ(a_0 - 1)][\eta_1 + k(a_0 - 1)\delta_1] - k^2(a_0)\delta_1^2$$

Absorbing  $\eta_1$  into  $S^2$  leads to the quadratic equation

$$A\delta_1^2 + B\delta + D = 0$$

with

$$A = 1 - 4(a_0 - 1)k^2(a_0)/b_0,$$

$$B = k(a_0)\{(1+2nIJ)[\omega(a_0-1)-\omega(a_0)]/IJ\omega(a_0) + 2(a_0-1)R_{...}/b_0\} - 1,$$

and

$$D = (R_{...}/nIJ) + b_0(a_0 + nIJ)/2nIJ(a_0 - 1)(a_0 - 2) - 2(a_0 - 1)S^2/b_0.$$

Thus, we obtain an estimate for  $\delta_1^2$  as

$$\tilde{\delta}_1^2 = \left\{ egin{array}{ll} (-B/2A)^2 & {
m if} & B^2 - 4AD \geq 0, \\ D/A & {
m if} & B^2 - 4AD < 0. \end{array} 
ight.$$

This gives  $\tilde{\eta}_1 = R_{\cdots} - b_0/2(a_0 - 1) - k_0\tilde{\delta}_1$ , while other elements of  $\tilde{\eta}$  being equal to those given for case 1. However, for  $i = 1, \dots, I-1$ ,

$$\begin{split} \tilde{\delta}_{i+1} &= 2(a_0-1)S_i^2/b_0 - [\tilde{\eta}_1 + \tilde{\eta}_{i+1} + \tilde{\delta}_1 k(a_0-1)]/nJ \\ &- b_0[2(a_0-1) + nJ]/2nJ(a_0-1)(a_0-2) - \tilde{\delta}_1^2 + \\ &2k(a_0)[\tilde{\delta}_1 k(a_0) - \tilde{\eta}_1]\tilde{\delta}_1(a_0-1)/b_0 - 2\tilde{\delta}_1[k(a_0-1) - k(a_0)]; \end{split}$$

for j = 1, ..., J - 1,

$$\begin{split} \tilde{\delta}_{I+j} &= 2(a_0-1)S_j^2/b_0 - [\tilde{\eta}_1 + \tilde{\eta}_{I+j} + \tilde{\delta}_1 k(a_0-1)]/nI \\ &- b_0[2(a_0-1) + nI]/2nI(a_0-1)(a_0-2) - \tilde{\delta}_1^2 + \\ &2k(a_0)[\tilde{\delta}_1 k(a_0) - \tilde{\eta}_1]\tilde{\delta}_1(a_0-1)/b_0 - 2\tilde{\delta}_1[k(a_0-1) - k(a_0)]; \end{split}$$

and finally, for  $i = 1, \dots, I-1$ , and  $j = 1, \dots, J-1$ ,

$$\begin{split} \tilde{\delta}_{I+i(J-1)+j} &= 2(a_0-1)S_{ij}^2/b_0 - [\tilde{\eta}_1 + \tilde{\eta}_{i+1} + \tilde{\eta}_{I+j} + \tilde{\eta}_{I+i(J-1)+j} + \tilde{\delta}_1 k(a_0-1)]/nI \\ &- b_0[2(a_0-1)+n]/2n(a_0-1)(a_0-2) - [\tilde{\delta}_1^2 + \tilde{\delta}_{i+1}^2 + \tilde{\delta}_{I+j}^2] \\ &+ 2k(a_0)[\tilde{\delta}_1 k(a_0) - \tilde{\eta}_1]\tilde{\delta}_1(a_0-1)/b_0 - 2\tilde{\delta}_1[k(a_0-1)-k(a_0)]. \end{split}$$

Thus, we have  $\tilde{\eta} = [\tilde{\eta}_1, \dots, \tilde{\eta}_J]$ ,  $\tilde{\Delta} = \text{diag}\{\tilde{\delta}_1^2, \dots, \tilde{\delta}_{IJ}^2\}$ , which provide the respective empirical Bayes estimates:

$$\lambda_{EB2} = (nIJ + 2a_0)/Q_1(\tilde{\eta})$$
 (4 - 29)

$$Var(\lambda|y) = 2(nIJ + 2a_0)/[Q_1(\tilde{\eta})]^2$$
 (4 - 30)

$$\tilde{\phi}_{1,EB2} = \tilde{\eta}_1^* + 0.21\omega(k)[\tilde{\psi}_{11}Q_1(\tilde{\eta})]^{1/2}$$
 (4 - 31)

$$Var(\phi|\mathbf{y}) = 0.045\tilde{\psi}_{11}Q_1(\tilde{\boldsymbol{\eta}})[11/(k-1) - \omega^2(k)] + 0.21\tilde{\boldsymbol{\eta}}_1^*\omega(k)[\tilde{\psi}_{11}Q_1(\tilde{\boldsymbol{\eta}})]^{1/2}$$

$$(4-32)$$

with

$$\begin{split} \tilde{\boldsymbol{\eta}}^{\star} &= [\tilde{\eta}_{1}^{\star}, \tilde{\boldsymbol{\eta}}^{\star}(2)] = [n\tilde{\Delta}M + I]^{-1}[n\tilde{\Delta}d + \tilde{\boldsymbol{\eta}}] \\ \tilde{\Psi} &= [n\tilde{\Delta}M + \tilde{\Delta}^{-1}] = \begin{bmatrix} \tilde{\psi}_{11} & \tilde{\Psi}_{12} \\ \tilde{\Psi}_{21} & \tilde{\Psi}_{22} \end{bmatrix} \end{split}$$

Moreover,

$$\Phi_{EB2}^{(2)} = \tilde{\eta}^{*(2)} + \{0.21\omega(k)[\psi_{11}^{-1}Q_1(\tilde{\eta})]^{1/2}\}\tilde{\Psi}_{21} \qquad (4-33)$$

$$Var(\Phi|\mathbf{y}) = [Q_1(\boldsymbol{\eta}/2(k-1))][\tilde{\Psi}_{22} - 0.09(k-1)\omega^2(k)\tilde{\Psi}_{21}\tilde{\psi}_{11}^{-1}\tilde{\Psi}_{12}] + 0.21\omega(k)\tilde{\eta}_1^*[\tilde{\psi}_{11}^{-1}Q_1(\tilde{\eta})]^{1/2}\tilde{\Psi}_{21}\tilde{\psi}_{11}^{-1}\tilde{\Psi}_{12}. \qquad (4-34)$$

Although we have used the method of moments to reach explicit solutions, we could have alternatively used the maximum likelihood procedure to obtain estimates of  $a,b,\eta$  and  $\Delta$ . This method needs numerical maximization which can be done by usual routines. Here, we only outline the procedure and leave the detail for practical data analysis. From (4-5) - (4-8),

$$\begin{split} \mathcal{L} &= l(\boldsymbol{y}|a,b,\eta,\Delta) = K \bigg\{ \int_0^\infty \bigg[ \lambda^{\nu-1} b^a |\Delta|^{-1/2} / \Gamma(a) \bigg] \exp \bigg\{ -\frac{\lambda}{2} Q_1(\eta) \bigg\} \\ &\times \int \exp{-\frac{\lambda}{2} Q_2(\phi) d\phi} \bigg\} \\ &= \frac{\Gamma(a+nIJ/2)}{\Gamma(a)} \cdot \frac{|\Psi|^{1/2}}{|\Delta|^{1/2}} \frac{(2b)^a}{[Q_1(\eta)]^{nIJ/2+a}} \end{split}$$

Maximizing  $\mathcal{L}$  with respect to a,b, $\eta$  and  $\Delta$  would provide the maximum likelihood estimates, denoted by  $\hat{a}, \hat{b}, \hat{\eta}, \text{and} \hat{\Delta}$ . Applying them in (4-16) and (4-17) would result in empirical Bayes estimates based on the

maximum likelihood procedure. Again, if we observe the estriction on  $\phi_1$ , we shall have the corresponding results. Let the result be xpressed as

$$\phi_{EBL} = (n\hat{\Delta}M + I)^{-1}(n\hat{\Delta} + \hat{\eta}), \qquad (4 - 35)$$

$$Var(\Phi|\mathbf{y}) = [Q_1(\hat{\boldsymbol{\eta}})/2(\kappa - 1)]\hat{\Psi}, \qquad (4 - 36)$$

with

$$\hat{\Psi} = (n\hat{\Delta}M + \hat{\Delta}^{-1})^{-1}.$$

The above derivation remains valid for the 2-factor ANOVA model without interaction, as well as for one-way ANOVA. In these cases, one only needs to reduce the order of the vector of parameters and the design matrix, according to the model used and follow the above procedure.

### 4.5 AN EXAMPLE

To illustrate our proposed estimators and compare them with other estimators, we analyze an experiment originally reported by Ostle(1963) and analyzed by Shuster and Muira (1972), and later by Achcar and Rosales (1993). Data in Table 5.1 ave resulted from a randomized  $2 \times 5$  layout with 10 replicates in each cell. The responses consist of the impact resistance of 5 kinds of insulators to shocks when they are cut lengthwise or widthwise. There are 10 replicates for each combination.

#### Maximum likelihood estimates:

It can be shown that the maximum likelihood estimates are

$$\hat{\Phi}(M\mathcal{L}) = M^{-1}d, \qquad d = [1, 0, \cdots, 0]$$

$$\hat{\lambda}(M\mathcal{L}) = IJ/[nR - d'M^{-1}d]$$

whose large - sample variances are

pectively.

$$Var[\hat{\Phi}(M\mathcal{L})] = (n\hat{\lambda})^{-1}M^{-1}$$
 
$$Var[\hat{\lambda}(M\mathcal{L})] = 2(nIJ)^{-1}[\hat{\lambda}(M\mathcal{L})]^{2}$$
 
$$Cov[\hat{\Phi}(M\mathcal{L}), \hat{\lambda}(M\mathcal{L})] = 0$$

In this example, the diagonal elements of  $\mathbf D$  are given in the last column of Table 4.1, and X' is

$$X' = \begin{bmatrix} 1'_4 & 1 & 1'_4 & 1 \\ 1'_4 & 1 & -1'_4 & -1 \\ I_4 & -1_4 & I_4 & -1_4 \\ I_4 & -1_4 & -I_4 & 1_4 \end{bmatrix}$$

which yield the estimates and their standard errors (S.E.) given in Table 4.2. The asymptotic 95 percent confidence intervals (CI) are also provided in Table 4.2. It is clear that except for  $\lambda$  no parameter can be taken different from zero at 5 percent level. However, only the constant  $\mu$  is different from zero at 0 percent level. But, due to wide confidence intervals one should feel uncertain about these inferences. Better inferences are possible by the empirical Bayes procedure presented below. Empirical Bayes estimates of a and b are  $a_0 = 6.95$  and  $b_0 = 0.26$ , es-

The estimates of the model parameters for each of the two cases (unrestricted nd restricted) along with the standard errors, are shown in Table 4.3. In Table 4.4, the 95 percent credible intervals (CI) based on the marginal posteriors are given. Again, we observe that only  $\lambda$  and  $\mu$  are infered to be different from zero at 5 percent level. Although we have reached the same conclusion as the one relative to MLE, but here we have much smaller standard errors which make the inference more precise. In fact, comparing Table 4.2 and 4.3, we note a striking

consequence of exploiting the empirical Bayes procedure. The credible intervals in both cases (unrestricted and restricted) are very much shorter than the corresponding confidence intervals given in Table 4.2. The only exception is the intervals for  $\lambda$  which have become somewhat longer for the empirical Bayes procedure.

Table 4.1 Observations from the experiment

$K \rightarrow$	1	2	3	4	5	6	7	8	9	10	Mean
(i,j) ↓									L-Verrous Marie Ma		$(y_{ij.})$
1,1	1.15	0.84	0.88	0.91	0.86	0.88	0.92	0.87	0.93	0.95	0.919
1,2	1.16	0.85	1.00	1.08	0.80	1.01	1.14	0.87	0.97	1.09	0.999
1,3	0.79	0.68	0.64	0.72	0.63	0.59	0.81	0.65	0.64	0.75	0.690
1,4	0.96	0.82	0.98	0.93	0.81	0.79	0.79	0.86	0.84	0.92	0.870
1,5	0.49	0.61	0.59	0.51	0.53	0.72	0.67	0.47	0.44	0.48	0.551
2,1	0.89	0.69	0.46	0.85	0.73	0.67	0.78	0.77	0.80	0.79	0.743
2,2	0.86	1.17	1.18	1.32	1.03	0.84	0.89	0.84	1.03	1.06	1.022
2,3	0.52	0.52	0.80	0.64	0.63	0.58	0.65	0.60	0.71	0.59	0.623
2,4	0.86	1.06	0.81	0.97	0.90	0.93	0.87	0.88	0.89	0.82	0.899
2,5	0.52	0.53	0.47	0.47	0.57	0.54	0.56	0.55	0.45	0.60	0.526

Table 4.2 ML Estimates of  $\lambda$  and  $\Phi$  and their symptotic 95 percent confidence intervals

Parameter	MLE	S.E.	95 percent	CI	$CI\ length$
λ	0.98	0.09	0.8	1.16	0.36
$\mu$	1.34	0.82	-0.27	2.95	3.22
$\alpha_1$	0.04	0.82	-1.57	1.65	3.22
$oldsymbol{eta_1}$	-0.13	1.58	-3.23	2.96	6.19
$eta_2$	-0.35	1.47	-2.27	2.53	4.80
$eta_3$	0.18	1.72	-3.19	3.55	6.74
$eta_4$	-0.21	1.54	-3.23	2.81	6.04
711	-0.09	1.58	-3.19	3.01	6.20
γ <sub>12</sub>	-0.06	1.47	-2.94	2.82	5.76
γ <sub>13</sub>	-0.03	1.71	-3.38	3.32	6.70
714	0.06	1.54	-2.96	3.08	6.04

Table 4.3 Empirical Bayes estimates

Prior	Unrestri	cted	Restri	cted
parameter	Estimated	S.E.	Estimate	S.E.
λ	0.7973	0.1057	1.2198	0.1616
μ	1.1502	0.2324	0.9858	0.1462
$\alpha_1$	-0.0004	0.1302	-0.0006	0.1239
$eta_1$	0.0216	0.2440	0.0177	0.2406
$eta_2$	-0.0555	0.3223	0.0001	0.3131
$eta_3$	0.0212	0.3733	0.0923	0.2107
$eta_4$	-0.0467	0.3193	-0.0001	0.1025
$\gamma_{11}$	0.0005	0.1781	-0.0014	0.3202
γ <sub>12</sub>	0.0012	0.2324	-0.0003	0.1422
γ <sub>13</sub>	-0.0001	0.2392	0.0005	0.1552
γ14	0.0016	0.2362	-0.0022	0.1471

Table 4.4 The 95 percent credible intervals for empirical Bayes stimates

Prior	Unrestri	icted		Restricted		
parameter	95 percent CI		CI length	95 percent	CI	CI length
λ	0.5901	1.0044	0.4139	0.9031	1.5365	0.6334
μ	0.6947	1.6057	0.9110	0.6992	1.2724	0.5732
$lpha_1$	-0.2556	0.2548	0.5104	-0.2434	0.2422	0.4856
$eta_1$	-0.4566	0.4998	0.9564	-0.4539	0.4893	0.9432
$eta_2$	-0.6872	0.5763	0.5950	-0.6136	0.6138	1.2274
$eta_3$	-0.7105	0.7529	1.4664	-0.3207	0.5053	0.8260
$eta_4$	-0.6725	0.5721	1.2516	-0.2010	0.2008	0.4018
711	-0.3486	0.3496	0.6982	-0.6290	0.6262	1.2552
<b>γ</b> 12	-0.4543	0.4567	0.9110	-0.2790	0.2784	0.5574
<i>γ</i> 13	-0.4689	0.4687	0.9376	-0.3037	0.3047	0.6084
714	-0.4614	0.4646	0.926	-0.2905	0.2861	0.5766

# 5. Empirical Bayes regression

#### 5.1 Introduction

In this section we look at the regression analysis under the Inverse Gaussian model. This problem has also been treated mainly from the classical point of view. Whitmore (1983) considers the censored data and estimates the regression coefficient, using the EM algorithm. Seto and Iwase (1985) derive the minimum variance unbiased estimator. However, Hsieh and Korwar (1990) prove that it is inadmissible. Chhikara and Folks (1989) and Seshadri (1999) each devote a chapter to the regression problem. Woldie and Folks (1999, 1994) are the most relevant

works, though they adopt the classical approach, too. We intend to extend their results by exploiting the Bayesian and empirical Bayes approaches.

#### 5.2 The model

Suppose we have observed the pairs of observations  $(y_i, x_i)$ , i = 1, ..., n where  $y_i$  are realizations of random variable  $Y_i$ which has the Inverse Gaussian distribution with mean  $(x'_i\beta)^q$  and the scale parameter  $\lambda_i$  where  $(x'_i$  and  $\beta$  are p-vectors with  $(x'_i\beta > 0)$ ,  $\alpha$  is an integer, and  $\alpha$  as simple inear regression on  $\alpha$  and  $\alpha$  provide different models such as simple inear regression through origin  $\{p = q = 1, \lambda_i = \lambda, x_i = (x_i), \beta = (\beta)\}$ , linear regression with intercept  $\{p = 2, q = 1, \lambda_i = \lambda, x_i = (1, x_{i2}), \beta = (\beta_1, \beta_2)\}$ , non-linear regression  $\{q = -1, \lambda_i = \lambda\}$ ; and general non-linear regression  $\{q = -1, \lambda_i = \lambda\}$ . The problem is to estimate  $\beta$  and  $\alpha$  from the observations  $\alpha$  of variance which we discussed in the previous section. The difference is in that here  $\alpha$  are real vectors while there in ANOVA they were indicators showing the presence (1) or absence (0) of a factor. Thus, the approach adopted there can be helpful in dealing with this problem.

# 5.3 Bayesian analysis

According to Bayesian way of thinking, we assume

$$Y_i|\beta, \lambda \sim IG\{(x_i'\beta)^q, \lambda\}, \qquad i = 1, \dots, n$$

where the parameters are treated as random variables, having some sort of probability distribution, which are called the prior distribution. To perform a Bayesian analysis, we need to specify the prior distributions  $\pi_1(\beta|\lambda)$  and  $\pi_2(\lambda)$ . While various authors have employed different prior distributions for  $\beta$  nd  $\lambda_i$ , we choose to use the conjugate priors for these

parameters. Those are explicitly stated as

$$\pi_1(\lambda) = \left(\frac{b}{2}\right)^{\frac{a}{2}} \lambda^{\frac{a}{2}-1} \exp \frac{\left\{-\frac{b\lambda}{2}\right\}}{\Gamma(\frac{a}{2})} \quad \lambda, a, b > 0, \tag{5-1}$$

and

$$\pi_2(\boldsymbol{\beta}|\lambda) = \text{const.} \lambda^{\frac{p}{2}} |\Delta|^{-\frac{1}{2}} \exp\left\{\frac{-(\boldsymbol{\beta} - \boldsymbol{\eta}')\Delta^{-1}(\boldsymbol{\beta} - \boldsymbol{\eta})}{2}\right\}$$
 (5 - 2)

where

$$egin{aligned} & oldsymbol{\eta} = [\eta_1, \ldots, \eta_p], oldsymbol{\eta} \in R^p \ & \Delta = \mathrm{diag}\{\delta_1, \ldots, \delta_p\}, \delta_j \in R^+. \end{aligned}$$

The likelihood function is reduced to

$$L(\beta, \lambda|) = \text{constant} + \lambda^{\frac{\pi}{2}} \exp\left\{\frac{-\lambda Q_1(\beta)}{2}\right\}$$
 (5-3)

with

$$Q_1(\beta) = \sum_{i=1}^{n} (y_i - x_i'\beta)^2 / y_i (x_i'\beta)^2$$
$$= [YX\beta - 1_n]'Y^{-1}[YX\beta - 1_n]$$

with

$$X = [x_1, \ldots, x_n], Y = \text{diag}\{y_1, \cdots, y_n\}, 1_n = [1, \ldots, 1].$$

Thus, the posterior distribution of  $(\lambda, \beta)$  is obtained as

$$q(\lambda, \beta|y) \propto \frac{b^{\frac{a}{2}} \lambda^{v-1}}{\Gamma(\frac{a}{2})|\Delta|^{\frac{1}{2}}} \exp\left\{\frac{-\lambda}{2} [Q_1(\beta) + Q_2(\eta) + b]\right\}$$
 (5-4)

$$v=\frac{(a+n+p)}{2}, \qquad \lambda>0, \qquad \beta\in R^p.$$

It makes easier to use the identity

$$Q_1(\beta) + Q_2(\eta) + b \equiv g(\eta, \Delta) + h(\eta, \Delta, b)$$

with

$$g(\eta, \Delta) = (\beta - m)' \sum_{i=1}^{-1} (\beta - m)$$

$$m = [I + \Delta X'YX]^{-1} [\eta + n\Delta \bar{x}]$$

$$\sum_{i=1}^{-1} [\Delta^{-1} + X'YX]^{-1}$$

$$h(\eta, \Delta) = \eta' \Delta^{-1} \eta + 1'_n Y^{-1} 1_n + b$$

$$-[\eta + n\Delta \bar{x}'] \Delta^{-1} [I + \Delta X'YX]^{-1} [\eta + n\Delta \bar{x}]$$

From (5-4), the marginal posterior distribution of  $\lambda$  is obtained by integrating on  $\beta$ . It follows that

$$q(\lambda|\boldsymbol{y}) \propto \lambda^{\frac{a+n}{2}-1} \exp\left\{\frac{-\lambda}{2}[h(\eta+\Delta+b)]\right\}$$

That is,

$$\lambda | \boldsymbol{y} \sim \Gamma \left[ \frac{a+n}{2}, \frac{h(\boldsymbol{\eta}, \boldsymbol{\Delta}, b)}{2} \right].$$
 (5-5)

Hence, the Bayes estimator of  $\lambda$  with respect to the squared error loss function is

$$\begin{cases} \hat{\lambda}_{B} = E(\lambda|\mathbf{y}) = \frac{a+n}{h(\boldsymbol{\eta}, \boldsymbol{\Delta}, b)} \\ var(\lambda|\mathbf{y}) = \frac{2(a+n)}{h(\boldsymbol{\eta}, \boldsymbol{\Delta}, b)^{2}} q \end{cases}$$
 (5-6)

For  $\beta$ , we integrate on  $\lambda$  in (5-4) and obtain

$$q(\boldsymbol{\beta}|\boldsymbol{y}) \propto \frac{1}{\left[(a+n) + (\boldsymbol{\beta} - \boldsymbol{m})'\Psi^{-1}(\boldsymbol{\beta} - \boldsymbol{m})\right]^{\frac{(a+n+p)}{2}}}$$
 (5-7)  
$$\Psi_{,} = \frac{h(\boldsymbol{\eta}, \boldsymbol{\Delta}, b)}{a+n} [\boldsymbol{\Delta}^{-1} + \boldsymbol{X}' \boldsymbol{Y} \boldsymbol{X}]^{-1}$$

That is,

$$\beta|y \sim T_p[(a+n)m, V]$$

Thus, for a + n > 2 we have

$$\begin{cases} \hat{\beta}_{B} = E(\beta|\mathbf{y}) = \mathbf{m} = [I + \Delta X'YX^{-1}[\mathbf{\eta} + n\Delta\bar{\mathbf{x}}] \\ \\ var(\beta|\mathbf{y}) = V = \frac{a+n}{a+n-2}\Psi \end{cases}$$
(5-8)

By the distributional properties of the multivariate T-distribution, each subvector of  $\beta$  also has a T-distribution with corresponding mean and variance. Even each linear transformation of  $\beta$  has a T-distribution, too. These facts prove useful in model selection, below. Therefore, the simple regression is estimated as

$$\hat{\mu}_{i,B}^{-1} = x'_{i}\hat{\beta}_{B}, \qquad i = 1, \dots, n$$
 (5-9)

which marginally has a T-distribution. This fact helps us to find a Highest Posterior Density (HPD) region akin to a confidence interval in the classical approach.

#### 5.4 Model selection

Selecting a particular model is tantamount to excluding some independent variables  $(x_j)$  from the full regression equation. This task can be accomplished either by testing the significance of various coefficients or by utilizing the Bayes factor. The former is the usual method in the classical approach. The latter method compares the two competing models  $M_{j-1}(y)$  and  $M_j(y)$  while  $M_{j-1}(y)$  is a reduced from of  $M_j(y)$ , lacking some variables. Then

$$P_{\cdot}O_{\cdot} = posterior \ odds = \frac{\alpha_{j}}{1 - \alpha_{j}} \frac{q(\boldsymbol{\beta}|\boldsymbol{y}, M_{j})}{q(\boldsymbol{\beta}|\boldsymbol{y}, M_{j-1})}, \ j = 1, 2, \dots, p$$

where  $0 < \alpha_j < 1$  is the prior probability that the  $M_j(y)$  is the correct model. If  $P_j(0) > 1$ , the model  $M_j(y)$  is more reasonable than  $M_{j-1}(y)$ ,

otherwise vice-versa.

## 5.5 Empirical Bayes analysis

In order to use the result (5-9) in practice, one has to be able to specify the actual values of the parameters in (5-1) and (5-2). This is not an easy task because there may not be a consensus among the users about the specified parameters. One way out of this difficulty is to let the data speak for themselves and determine those parameters. The operational aspect of this procedure is to estimate the parameters of the prior distributions from the observed data and then follow the Bayesian rules. The marginal distribution of the random sample which is observed is

$$f(\boldsymbol{y}|\boldsymbol{a},\boldsymbol{b},\boldsymbol{\eta},\Delta) = \int \cdots \int \bigg\{ \int_0^\infty f(\boldsymbol{y}|\boldsymbol{\beta},\lambda) \pi_2(\boldsymbol{\beta}|\lambda) \pi_1(\lambda) d\lambda \bigg\} d\boldsymbol{\beta}$$

which by virtue of (5-1)-(5-3) and the properties of the gamma and multivariate T distributions reduces to

$$f(\boldsymbol{y}|a,b,\boldsymbol{\eta},\Delta) = constant \cdot \frac{b^a [h(\boldsymbol{\eta},\Delta,b)]^{\frac{p}{2}} \Gamma(\frac{a+n}{2})}{[a+n]^{\frac{a+n+p}{2}} \Gamma(a)|I+\Delta X'YX|^{\frac{1}{2}}}$$
$$\boldsymbol{y} \in [\mathbb{R}^+]^n \tag{5-10}$$

That is, (5-10) is in fact the likelihood function of  $(a,b,\eta,\Delta)$  which contains 2(p+1) parameters. Now, similar to the previous section, we can maximize (5-10) in terms of  $(a,b,\eta,\Delta)$  to obtain the MLE of  $(a,b,\eta,\Delta)$  which are denoted by  $(\hat{a},\hat{b},\hat{\eta},\hat{\Delta})$ . Next we substitute these estimates in (5-6)-(5-8) to obtain the  $\hat{\lambda}_{EBML}$  and  $\hat{\beta}_{EBML}$  which in turn yield

$$\hat{\mu}_{i,EBML}^{-1} = x_i' \hat{\beta}_{EBML} \tag{5-11}$$

which being a linear function of  $\hat{oldsymbol{eta}}_{EBML}$  has the variance as

$$var(\hat{\mu}_{i,EBML}^{-1}) = x_i'var(\hat{\beta}_{EBML})x$$

Since the asymptotic variance of  $\hat{\beta}_{EBML}$  is obtained by the inverse of the Fisher information matrix  $I^{-1}(\hat{\beta}_{EBML})$ , it follows that

$$var(\hat{\mu}_{i,EBML}) = x_i'I^{-1}(\hat{\beta}_{EBML})x.$$

This martrix is used in finding the HPD region, or confidence bands for the regression equation (5-11).

The above procedure needs numerical computations to find  $(\hat{a}, \hat{b}, \hat{\eta}, \hat{\Delta})$  from (5-10). Thus he solution has no closed from. In order to provide an explicit solution we may use the method of moment to find the estimates of  $(a, b, \eta, \Delta)$ . If  $Y \sim IG(\mu, \lambda)$  with the density  $f(y|\mu, \lambda)$  then  $R_i = Y_i^{-1}$  has the density

$$k(r|\mu,\lambda) = \mu r f(r|\mu^{-1},\lambda\mu^{-2}), , \mu,\lambda > 0$$

with moments of all order, Chhikara and Folks (1989, P. 43). Using this fact we find the first two marginal moments of  $R_i$ , i.e.

$$E(R_i) = \gamma + x_i' \eta \tag{5-12}$$

and

$$var(R_i) = \gamma^2 \xi + \gamma x_i' \eta + x_i' \Delta x_i', \quad i = 1, \dots, n$$
 (5 - 13)

with  $\gamma = \frac{b}{2(a-2)}$ ,  $\xi = \frac{2(a-1)}{(a-4)}$ . The equation (5-12) implies that

$$R_i = \gamma + x'\eta + \epsilon_i, \quad i = 1, \dots, n$$

which is similar to the multiple linear regression and can be expressed as

$$R = Z\eta + \epsilon \tag{5-14}$$

with  $R = [R_1, \ldots, R_n]$ ,  $Z = [1_n, X]$ ,  $X = (x_{ij})$   $\eta = [\gamma, \eta]$ , and  $\epsilon = [\epsilon_1, \ldots, \epsilon_n]$ . Then the least square solution of (5-13) is

$$\tilde{\boldsymbol{\eta}}_1 = (Z'Z)Z'R \tag{5-15}$$

A similar treatment of (5-13) leads to the system of equations

$$S_i^* = S_{(i)}^2 - \bar{\gamma} x_i' \bar{\eta} = \bar{\gamma}^2 \xi + \sum_{j=1}^p x_{ij}^2 \cdot \delta_j, \quad i = 1, \dots, n$$
 (5 - 16)

where, for  $\bar{R}_{(i)} = \sum_{j \neq i}^{n} \frac{R_{j}}{(n-1)}$ , we have defined

$$S_{(i)}^2 = \sum_{j \neq i}^n \frac{[R_j - \bar{R}_{(i)}]^2}{(n-2)}, \quad i = 1, \dots, n.$$

Now, we define

$$S^* = [S_1^*, \cdots, S_n^*],$$
 
$$T = [\tilde{\gamma}^2 1_n, X^*], \quad X^* = (x_{ij}^2)$$
  $\epsilon = [\epsilon_1, \dots, \epsilon_n], \quad \delta_c = \Delta 1_p = [\delta_1, \dots, \delta_p]$ 

and  $\xi_1 = [\xi, \delta_c]$ . Thus, (5-16) is in fact as

$$S^* = Txi_1 + \epsilon \tag{5-17}$$

From (5-17), through the least squares method, one obtains

$$\tilde{\xi}_1 = (T'T)^{-1}T'S^* \tag{5-18}$$

If it happens that in (5-15) and (5-18) the inverse matrices do not exist, their Moor-Penrose inverse would be used. We have, therefore, the estimates of  $\gamma$ ,  $\xi$ ,  $\eta$  and  $\Delta$ . The pair  $\gamma$  and  $\xi$  gives us  $a = \frac{2(2\xi-1)}{(\xi-2)}$  and  $b = \frac{2\gamma(\xi+1)}{(\xi-2)}$ . These estimates are denoted by  $(\hat{a}, \hat{b}, \hat{\eta}, \hat{\Delta})$  which will be used n the Bayes estimates already obtained as (5-9). Thus, as an alternative to (5-11) we have

$$\mu_{i,EBMM}^{-1} = \mathbf{x}_i' \hat{\boldsymbol{\beta}}_{EBMM} \tag{5-19}$$

with

$$var(\hat{\mu}_{i,EBMM}) = x'_i var(\hat{\beta}_{EBMM})x.$$
 (5 - 20)

The estimales obtained above either by using the ML or MM method are in fact Bayes estimates relative to the estimated proir distributions. Thus any inference concerning  $\lambda$  and  $\beta$  should be based on thier posterior distributions. For example in the case of the MM method we have

$$(\lambda|\mathbf{y}) \sim \Gamma\left[\frac{(\tilde{a}+n)}{2}, \frac{h(\tilde{\eta}, \tilde{\Delta}, \tilde{b})}{2}\right],$$
 (5 – 21)

$$(\boldsymbol{\beta}|\boldsymbol{y}) \sim T_P[(\tilde{a}+n), \tilde{\boldsymbol{m}}, \tilde{V}]$$
 (5 – 22)

Hence, the variance in (5-20) is found from (5-22). Likewise the distribution, and the HPD region for  $\mu_{i,EBMM}^{-1}$  is found from (5-20). Table 5.1 shows the data set relsuted from an experiment on the turnip plant. It is believed that the explanatory variables.  $X_1$ =sunlight,  $X_2$ =soil humidity, and  $X_3$ =air temperature affect y =the content of vitamin  $B_2$  in the leaves of the turnip plant. We would like to establish a relation between  $\mu_i$  and these 3 variables in the from of  $\mu_i = [x_i'\beta]^{-1}$ . To this end, we define  $Z = [1_{27}, X]$ , where X is the last three coumns of Table 5.1. Using (5-15), we abtain the estimates

$$[\tilde{\gamma}, \tilde{\eta_1}, \tilde{\eta_2}, \tilde{\eta_3}] = 10^{-6} [12581, 2, 134, -28].$$

To obtain  $\hat{\beta}$ , we first compute the statistics

$$\bar{R}_{(i)} = \frac{(0.339 - R_i)}{26}, S_{(i)}^2 = 0.04[0.005 - R_i^2 - 26\bar{R}_{(i)}^2], \quad i = 1, \dots, 27.$$

From these, we find  $S^*$  and T.Finally (5-18) gives us  $\xi$  which in turn provides

$$(\boldsymbol{\beta}|\boldsymbol{y}) \sim T_3[28, \tilde{\boldsymbol{m}}, \tilde{V}]$$

with

$$\hat{\boldsymbol{\beta}}_{EBMM} = \widetilde{\boldsymbol{m}} = 10^{-7} [-87, 4762, 318]$$

$$\tilde{V} = \text{var}(\boldsymbol{\beta}|\boldsymbol{y}) = 10^{-6} \begin{bmatrix} 43505 & -205205 & 36960 \\ & 21945 & 2695 \\ & & 23870 \end{bmatrix}$$

and finally, from (5-19), we arrive at

$$\tilde{\mu}_{i,EBMM}^{-1} = 10^{-7} [-87x_{i1} + 4762x_{i2} + 318x_{i3}]$$

The use and interpretation of this regression equation is like the usual regression model. Hence we do not elaborate on it.

# 6. Concluding remarks

We have journyed a relatively long path starting from the genesis of the Inverse Gaussian distribution and heading towards inference about its parameters. Along the way we used the Bayesian and empirical Bayes vehicle to reach estimates, to do analysis of variance and to construct a regression equation. It is not the end of journey. we could have visited the subject of analysis of covariance, or could have studied the reliability problems under this model. the study of asymptotic properties of various estimators is yet another common ground to work on. The posibilities are vast and the problems endless. One could conjecture that there are as many problem under the Inverse Gaussian model as there are under the normal (Gaussian) model. Most of them are not yet explored and await for eager and perseverant workers.

Table 5.1: the data regarding the vitamin  $B_2$  content of the turnip leaves and sunlight, soil humidity and air emperature

plant	Vitamin B <sub>2</sub>	Sunlight	Soil Hmidity	Air temperature
number	Y	$X_1$	$X_2$	$X_3$
1	110.4	176	7.0	78
2	102.8	155	7.0	89
3	101.0	273	7.0	89
4	108.4	273	7.0	72
5	100.7	256	7.0	84
6	100.3	280	7.0	87
7	102.2	280	7.0	74
8	93.7	184	7.0	87
9	98.9	216	7.0	88
10	96.6	198	2.0	76
11	99.4	59	2.0	65
12	96.2	80	2.0	67
13	99.0	80	2.0	62
14	88.4	105	2.0	70
15	75.3	180	2.0	73
16	92.0	180	2.0	65
17	82.4	177	2.0	76
18	77.1	230	2.0	82
19	74.0	203	47.4	76
20	65.7	191	47.4	83
21	56.8	191	47.4	82
22	62.1	191	47.4	69
23	61.0	76	47.4	74
24	53.2	213	47.4	76
25	59.4	213	47.4	69
26	58.7	151	47.4	75
27	58.0	205	47.4	76

Source: Draper and Smith (1981, P.406)

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Proceedings of the Monthly Colloquium of the Iranian Mathematical Society Vol. 2, pp 147-154 (2004)

# A Note on the Concept of Limit

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Abstract: The paper investigates the problems that students usually have when beginning the study of the notion of limit in its understanding and usage.

# I. Convergence in terms of sequences

A very common problem that most beginners will have with the definition of limit is to adjust the static nature of the definition with the dynamic nature of the words "tending", "converging", that they have in their minds.

Another problem is the difficulty of establishing the inequality  $|x_n - x| < \varepsilon$  for large values of n in order to prove  $\{x_n\}$  converges to x as  $n \to \infty$ .

MSC (2000): 26A03, 97C90

Keywords: Limit, Continuity,  $\epsilon - \delta$  definition, Sequential approach, Heuristic definition of limit

A common method to solve the latter inequality for n is to prove first that  $|x_n - x| \le a_n^{-1}$  for some unbounded increasing sequence  $\{a_n\}$  and then replace the inequality  $|x_n - x| < \varepsilon$  by the sufficient condition  $a_n > \varepsilon^{-1}$ . The new inequality has a solution of the form  $n > N_{\varepsilon}$  for some natural number  $N_{\varepsilon}$ . In most practical examples the sequence  $\{a_n\}$  and the number  $N_{\varepsilon}$  can be obtained through algebraic or other simple methods. For example, for  $\lim_{n \to \infty} \frac{n + \sin n}{2n - \cos n + 5} = \frac{1}{2}$ , we have

$$\left| \frac{n + \sin n}{2n - \cos n + 5} - \frac{1}{2} \right| \le \left( \frac{n}{2} + 1 \right)^{-1} < \varepsilon,$$

and hence

$$N_{\varepsilon} = \max\{1, [[2(\varepsilon^{-1} - 1)]]\},$$

where [[t]] denotes the greatest integer in t.

In general, the sequence  $\{a_n\}$  may not have a simple expression but, as the following proposition shows, it always exists.

**I.1. Proposition.** Let  $\{x_n\}$  be a sequence in a metric space X converging to  $x \in X$ ; i.e., for every  $\varepsilon > 0$  there exists  $N_{\varepsilon} \in \mathbb{N}$  such that  $d(x_n, x) < \varepsilon$  for all  $n > N_{\varepsilon}$ , where d denotes the metric on X. Then there exists an unbounded increasing sequence  $\{a_n\}$  of positive numbers such that  $d(x_n, x) \le a_n^{-1}$  (n = 1, 2, ...).

**Proof.** Let  $N_1$  be a positive integer such that  $d(x_n, x) < 1$  for all  $n > N_1$  and define

$$a_1 = a_2 = \cdots = a_{N_1} = (\max\{d(x_1, x), d(x_2, x), \ldots, d(x_{N_1}, x), 1\})^{-1}.$$

Assume by induction that  $N_1, N_2, \ldots, N_k$  and  $a_1, a_2, \ldots, a_{N_k}$  are defined for some  $k \geq 1$ . Let  $N_{k+1} > N_k$  be a positive integer such that  $d(x_n, x) < 1/(k+1)$  for all  $n > N_{k+1}$ . Define

$$a_{N_k+1}=a_{N_k+2}=\cdots=a_{N_{k+1}}=k.$$

In view of the Archimedes principle or the axiom of completeness, the increasing sequence  $\{a_n\}$  thus obtained is unbounded. Moreover,  $d(x_n, x) \le a_n^{-1}$  (n = 1, 2, ...)

It is clear that if a sequence  $\{x_n\}$  in a metric space (X,d) satisfies  $d(x_n,x) \leq a_n^{-1}$   $(n=1,2,\ldots)$  for some  $x \in X$  and some unbounded increasing sequence  $\{a_n\}$ , then  $\lim_{n\to\infty} x_n = x$ . Thus, one can see that the axiom of completeness and the definition of  $\lim_{n\to\infty} x_n$  are equivalent to the following Axiom I.2 and Definition I.3.

- **I.2.** Axiom. If  $y_1 < y_2 < ...$  is bounded in R, then there exists  $y \in R$  such that  $(y y_n)^{-1}$  is unbounded.
- **I.3. Definition.** A sequence  $\{x_n\}$  in a metric space X is said to be convergent to some  $x \in X$  if  $d(x_n, x) \leq 1/a_n$  (n = 1, 2, ...) for some unbounded increasing sequence  $\{a_n\}$  of positive numbers.

Axiom I.2 guarantees the unboundedness of the increasing sequences  $\{n\}$ ; if  $\{n\}$  is bounded then there exists  $y \in R$  such that  $\{(y-n)^{-1}\}$  is unbounded. Hence  $1 = (y-n) - (y-n-1) \le y-n$  for all  $n \in N$ ; a contradiction. This shows that, in the light of Axiom I.2, Definition I.3 is not vacuous. The inequality  $0 \le d(x_n, x) \le a_n^{-1}$  reflects the compression of a sequence  $\{x_n\}$  to x by the boundary of a shrinking sphere of radius  $a_n^{-1}$  centered at x. The terms "compress" and "shrink" have dynamical natures which one expects in the notion of limit.

If the standard  $\varepsilon - N$  definition of  $\lim_{n \to \infty} x_n = x$  is replaced by the new definition, then one can successfully prove and discuss all necessary results in a calculus or advanced calculus course related to limits of sequences. (The definition of  $\lim_{n \to +\infty} x_n = \pm \infty$  can be easily given as  $x_n > a_n$  or  $x_n < -a_n$  for some unbounded increasing sequence  $\{a_n\}$ .)

Moreover, limit, continuity, and differentiability of functions can be given in terms of sequences. In fact  $\lim_{x\to c} f(x) = L$  can be defined as the requirement  $\lim_{n\to\infty} f(x_n) = L$  for all sequences  $\{x_n\}$  in the domain of f

such that  $x_n \neq c$  and  $\lim_{n\to\infty} x_n = c$ . (For continuity of f at c we require  $\lim_{n\to\infty} f(x_n) = f(c)$  for any sequence  $\{x_n\}$  in the domain of f converging to c.)

Thus, to prove the composition  $f \circ g$  of continuous functions f and g is continuous, we simply observe  $\lim_{n\to\infty} f(g(x_n)) = f(\lim_{n\to\infty} g(x_n)) = f(g(c))$  whenever  $\{x_n\}$  is in the domain of g converging to g. The intermediate value theorem, the compactness of a closed interval, and the convergence of a Cauchy sequence, etc. can be proved in similar sequential fashions. For example, if a continuous function  $f:[a,b]\to R$  satisfies  $f(a)\leq k\leq f(b)$ , one can construct a nest of intervals  $[a,b]\supset [a_1,b_1]\supset [a_2,b_2]\supset \ldots$  such that  $b_n-a_n=(b-a)/2^n$  and  $f(a_n)\leq k\leq f(b_n)$ . Hence, if  $c=\lim a_n=\lim b_n$ , then  $0\leq (k-f(c))^2=\lim_{n\to\infty} (k-f(a_n))(k-f(b_n))\leq 0$  and thus k=f(c).

## II. Shortcomings of the sequential definitions

As we mentioned in the previous section, the sequential definition of convergence provides a dynamic spirit for the concept of limit. Even in the case of functions, students expect a dynamical approach. In fact, most textbooks and instructors of calculus respond to this need by beginning the subject of limit with tables showing the values of a function as simple as  $x^2$  at various points  $x_1, x_2, \ldots$  which cluster around a given point x = c and demonstrate that as  $x_n$  approaches c,  $x_n^2$  approaches  $c^2$ . This means that while the mathematicians of the nineteenth century regarded continuity of a curve as the free movement of a point on a plane, the twentieth century mathematicians regard the continuity of a function f as the convergence of the sequence  $\{f(x_n)\}$  to f(c) for any sequence  $\{x_n\}$  converging to c in the domain of f. It is well known that this definition of continuity is equivalent with the standard  $\varepsilon - \delta$  definition: For all  $\varepsilon > 0$ , there exists  $\delta = \delta(\varepsilon, c) > 0$  such that  $|f(x) - f(c)| < \varepsilon$ 

whenever  $|x-c| < \delta$ . (Absolute values can be replaced by distances in metric spaces.) If  $\delta$  is independent of the point c, the continuity of f is said to be uniform. It is not easy to give a definition of uniform continuity purely in terms of sequences; even if it is done, it is not handy to be used in the proofs based on uniform continuity. For example, one may define uniform continuity as having  $\lim_{n\to\infty} |f(x_n)-f(y_n)|=0$  whenever  $|x_n-y_n|\to 0$ . But such a definition is not convenient for proving the integrability of continuous functions. Similarly, uniform convergence of sequences of functions are easily defined and handled by the  $\varepsilon-N$  definition; we found it quite artificial to use sequential definitions to prove, for instance, that uniform limit of a sequence of continuous functions is a continuous function.

In general, we believe the  $\varepsilon-N$  or  $\varepsilon-\delta$  definitions of limit (for sequences or functions) are indispensable parts of mathematics. They are not only suitable for problems involving uniformity, but are the gates to important subjects such as topology. The  $\varepsilon-\delta$  definition introduces the concept of neighborhoods and intimates the beginner with the manipulation of abstract topological concepts.

With all these, we still believe the (dynamic) sequential definitions of limit and continuity are worth to be included in calculus textbooks. This is what a student expects when for the first time he/she encounters the notion of convergence. Proposition I.1 and the paragraph preceding Axiom I.2 showed that the  $\varepsilon-N$  definition is equivalent to Definition I.3. Thus, a beginner mathematician may begin the subject of limit by Definition I.2 and then observe the standard  $\varepsilon-N$  definition as a theorem. The difference between something to be regarded as a "definition" or as a "theorem" is that in the first case one has to accept a metafore of a concept which may differ from one's expectation of that concept, while in the second case one defends a claim that one has successfully

proved.

However, once the equivalence of the two methods of approach are established, it is easy to switch from one method to another. The following section shows that how combination of the two methods may help to shorten the proofs of the theorems.

# III. Uniformity and integrals

The Riemann integral in calculus is a concept which benefits lot from uniformity. As mentioned before, a function f is uniformly continuous if and only if  $d(f(x_n), f(y_n)) \to 0$  whenever  $d(x_n, y_n) \to 0$  and  $\{x_n\}$  and  $\{y_n\}$  are sequences in the domain of f. A sequence  $\{f_n\}$  is said to be uniformly convergent to a function f on a set A, if for every  $\varepsilon > 0$  there exists  $N \in \mathbb{N}$  such that n > N implies that  $|f_n(x) - f(x)| < \varepsilon$  for all  $x \in A$ . This is equivalent with saying that  $|f_n(x) - f(x)| \le a_n^{-1}$  for all  $x \in A$ , where  $\{a_n\}$  is an unbounded increasing sequence (independent of x).

The Riemann integral of f on [a,b] exists if, by definition, there exists an increasing sequence  $\{\pi_n\}$  of partitions of [a,b] such that  $\lim_n l(\pi_n) = \lim_n [\overline{S}(f,\pi_n) - \underline{S}(f,\pi_n)] = 0$ , where  $l(\pi) := \max_k (x_k - x_{k-1})$ ,  $\overline{S}(f,\pi) := \sum_k M_k(\pi)(x_k - x_{k-1})$ ,  $\underline{S}(f,\pi) := \sum_k m_k(\pi)(x_k - x_{k-1})$ ,  $M_k(\pi) := \sup\{f(x) : x_{k-1} \le x \le x_k\}$  for a given partition  $\{a = x_0 < x_1 < \dots < x_n = b\}$ . Then  $\int_a^b f$  is defined to be the limit of the increasing sequence  $\{\underline{S}(f,\pi_n)\}$ . As an immediate consequence, |f| is integrable and  $|\int_a^b f| \le \int_a^b |f|$ , if  $\int_a^b f$  exists. Note. If  $\pi'_n$  is a refinement of  $\pi_n$  and if  $\{\pi_n\}$  is as in the definition

Note. If  $\pi'_n$  is a refinement of  $\pi'_n$  and if  $\{\pi_n\}$  is as in the definition of  $\int_a^b f$ , it is easy to see that  $\{\pi'_n\}$  also defines the integrability of f and yields the same value for  $\int_a^b f$ . Thus  $\int_a^b f$  is independent of the choice of  $\{\pi_n\}$ . For a partition  $\pi$  as above we further set  $\mu(\pi) = \max_k [M_k(\pi) - m_k(\pi)]$ .

The proof of the following results would be lengthy if one tries to stick to one type of definitions. The proofs given here seem to be very short.

III.1. Theorem. Let  $f:[a,b] \to R$  be bounded and let  $\{\pi_n\}$  be an increasing sequence of partitions of [a,b] such that  $\lim_{n\to\infty} l(\pi_n) = 0$ . Then  $\int_a^b f \text{ exists if either the sequence } v_n := \sum_k [M_k(\pi_n) - m_k(\pi_n)] \text{ is bounded or } \lim_n \mu(\pi_n) = 0$ . In particular, if f is monotone or continuous, then it is Riemann integrable.

**Proof.** If f is monotone or, more generally, if  $v_n \leq D$  for all  $n \in \mathbb{N}$ , then  $\overline{S}(f,\pi_n) - \underline{S}(f,\pi_n) \leq l(\pi_n)D$  for all  $n \in \mathbb{N}$  and hence  $\int_a^b f$  exists. If f is continuous or, more generally, if  $\lim_n \mu(\pi_n) = 0$ , then  $\overline{S}(f,\pi_n) - \underline{S}(f,\pi_n) \leq (b-a)\mu(\pi_n)$  for all  $n \in \mathbb{N}$  and hence  $\int_a^b f$  exists.

III.2. Theorem. Let f be the uniform limit of a sequence of Riemann integrable functions  $f_n : [a,b] \to \mathbb{R}$ . Then  $\int_a^b f$  exists.

Proof. Let  $\{\pi_n\}$  be an increasing sequence of partitions of [a,b] with

**Proof.** Let  $\{\pi_n\}$  be an increasing sequence of partitions of [a,b] with  $\lim_n l(\pi_n) = 0$ . Assume  $\int_a^b f$  does not exist. Then  $\overline{S}(f,\pi_n) - \underline{S}(f,\pi_n) \geq \varepsilon_0$  for all  $n \in \mathbb{N}$  and some  $\varepsilon_0 > 0$ . Let  $\{a_n\}$  be an unbounded increasing sequence establishing the uniform convergence of  $\{f_n\}$ . Since

$$M_k(f,\pi_n) - m_k(f,\pi_n) \le M_k(f_i,\pi_n) - m_k(f_i,\pi_n) + 2a_i^{-1},$$

it follows that

$$0 < \varepsilon_0 \le \overline{S}(f_i, \pi_n) - \underline{S}(f_i, \pi_n) + 2a_i^{-1}(b-a)$$
,  $(i = 1, 2, \cdots)$ .

Letting  $n \to \infty$ , it follows that  $0 < \varepsilon_0 \le 2a_i^{-1}(b-a)$  and hence  $a_i \le 2(b-a)/\varepsilon_0$ , (i=1,2,...); a contradiction.

We conclude this topic with another famous theorem.

III.3. Theorem. If  $\int_a^b f$  exists and if  $\varphi$  is continuous on some closed interval containing f([a,b]), then  $\int_a^b \varphi \circ f$  exists.

**Proof.** The  $\varepsilon$ - $\delta$  definition of continuity implies that  $\varphi$  is the uniform limit of a sequence  $\{\varphi_n\}$  of piecewise linear continuous functions which can be expressed in the form

$$\varphi_n(x) = a + bx + \sum_{k=0}^m c_k |x - x_k|,$$

where the constant  $m, a, b, c_0, c_1, \dots, c_m, x_0, x_1, \dots, x_m$  depend on n. The rest of the proof follows from Theorem III.2 and the fact that  $\varphi_n \circ f$  is the sum of integrable functions.

## Bibliography

For a history of the concepts discussed in this paper we may refer to [1]. The details of this work is given in [2]. For a similar approach we refer to [3].

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Proceedings of the Monthly Colloquium of the Iranian Mathematical Society Vol. 2, pp 155-177 (2004)

# Orthogonal Functions in the Calculus of Variations and Optimal Control

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Abstract: The solution of problems in the calculus of variations is obtained by using hybrid functions. The properties of the hybrid functions which consist of block-pulse functions plus Legendre polynomials and block-pulse functions plus Chebyshev polynomials are presented. Two examples are considered, in the first example the brachistochrone problem is formulated as a non-linear optimal control problem, and in the second example an application to a heat conduction problem is given. The operational matrix of integration in each case is introduced and is utilized to reduce the calculus of variations problems to the solution of algebraic equations. The method is general, easy to implement and yields very accurate results.

## 1. Introduction

There has been a considerable renewal of interest in the classical problems of the calculus of variations both from the point of view of mathematics and of applications in physics, engineering, and applied mathematics.

Finding the brachistochrone, or path of quickest decent, is a historically interesting problem that is discussed in virtually all textbooks dealing with the calculus of variations. In 1696, the brachistochrone problem was posed as a challenge to mathematicians by John Bernoulli. The solution of the brachistochrone problem is often cited as the origin of the calculus of variations as suggested in [1]. The classical brachistochrone problem deals with a mass moving along a smooth path in a uniform gravitational field. A mechanical analogy is the motion of a bead sliding down a frictionless wire. The solution to this problem has been obtained by various methods such as the gradient method [2], successive sweep algorithm [3-4], the classical Chebyshev method [5] and multistage Monte Carlo method [6].

Orthogonal functions (OF's) have received considerable attention in dealing with various problems of dynamic systems. The main characteristic of this technique is that it reduces these problems to those of solving a system of algebraic equations; thus greatly simplifying the problem. The approach is based on converting the underlying differential equations into an integral equation through integration, approximating various signals involved in the equation by truncated orthogonal series and using the operational matrix of integration P, to eliminate the integral operations. The form of P depends on the particular choice of the orthogonal functions. Special attention has been given to applications of Walsh functions [7], block-pulse functions [8], Laguerre series [9], Legendre polynomials [10] and Chebyshev polynomials [11].

There are three classes of sets of OF's which are widely used. The first includes sets of piecewise constant basis functions (PCBF'S) (e.g., Walsh, block-pulse, etc.). The second consists of sets of orthogonal polynomials (OP's) (e.g., Laguerre, Legendre, Chebyshev, etc.). The third is the widely used sets of sine-cosine functions (SCF's) in Fourier series. While OP's and SCF's together form a class of continuous basis functions, PCBF's have inherent discontinuities or jumps. The inherent features(continuity or discontinuity) of a set of OF's largely determine their merit for application in a given situation. References [12] and [13] have demonstrated the advantages of PCBF spectral methods over Fourier spectral techniques. If a continuous function is approximated by PCBF's, the resulting approximation is piecewise constant. On the other hand if a discontinuous function is approximated by continuous basis functions the discontinuities are not properly modeled. Signals frequently have mixed features of continuity and jumps. These signals are continuous over certain segments of time, with discontinuities or jump occuring at the transitions of the segments. In such situations, neither the CBF's nor PCBF's taken alone would form an efficient basis in the representation of such signals.

The direct method of Ritz and Galerkin in solving variational problems has been of considerable concern and is well covered in many textbooks [14], [15]. Chen and Hsiao [7] introduced the Walsh series method to variational problems. Due to the nature of the Walsh functions, the solutions obtained were piecewise constant. Hwang and Shih [9], Chang and Wang [10] and Horng and Chou [11], used Laguerre polynomials, Legendre polynomials and Chebyshev polynomials respectively to derive continuous solutions for the first example in [7]. Furthermore, Razzaghi and Razzaghi [16], [17] applied Fourier series and Taylor series respectively to derive continuous solution for the second example in [7] which

is an application to the heat conduction problem. It is shown in Razzaghi and Razzaghi [17] that, to obtain the Taylor series coefficient, an ill-conditioned matrix commonly known as the Hilbert matrix is used. Hence the Taylor series is not suitable for the solution of the second example in [7].

In the present paper we introduce a new direct computational method to solve problems of the calculus of variations. The method consists of reducing the variational problems into a set of algebraic equations by first expanding the candidate functions as hybrid functions with unknown coefficients. The hybrid functions, which consists of block-pulse functions plus

- a) Legendre polynomials and
- b) Chebyshev polynomials

are first introduced. The operational matrix of integration in each case is given and is used to evaluate the coefficients of hybrid functions in such a way that the necessary conditions for extremization are imposed. Two examples are considered. In example 1, the brachistochrone problem is formulated as an optimal control problem and in the second example we will demonstrate the application of operational matrix of integration for hybrid functions by considering the second example in [7]. It is shown that the hybrid functions of block-pulse and Legendre polynomials approach produces an exact solution for the heat conduction problem.

# 2. Properties of Hybrid Functions of Block-Pulse and Legendre Polynomials

Hybrid functions  $b(n, m, t), n = 1, 2, \dots, N, m = 0, 1, \dots, M - 1$ , have three arguments; n is the order of block-pulse functions, m is the order of Legendre polynomials, and t is the normalized time. They are defined

on the interval  $[0, t_f)$  as

$$b(n,m,t) = \begin{cases} P_m(\frac{2N}{t_f}t - 2n + 1), & t \in \left[\left(\frac{n-1}{N}\right)t_f, \frac{n}{N}t_f\right) \\ 0, & \text{otherwise.} \end{cases}$$
(1)

Here  $P_m(t)$  are the well-known Legendre polynomials of order m which are orthogonal with respect to the weight function w(t) = 1 and satisfy the following recursive formula.

$$P_{o}(t) = 1, P_{1}(t) = t$$

$$P_{m+1}(t) = \left(\frac{2m+1}{m+1}\right) t P_m(t) - \left(\frac{m}{m+1}\right) P_{m-1}(t) , \quad m = 1, 2, 3, \cdots$$

Since b(n, m, t) consists of block-pulse functions and Legendre polynomials, which are both complete and orthogonal, the set of hybrid functions of block-Pulse and Legendre polynomials is a complete orthogonal set.

### 2.1. Function Approximation

A function f(t), defined over the interval 0 to  $t_f$  may be expanded as

$$f(t) = \sum_{n=1}^{\infty} \sum_{m=0}^{\infty} c(n,m)b(n,m,t), \qquad (2)$$

where

$$c(n,m) = (f(t),b(n,m,t))$$

in which (.,.) denotes the inner product. If the infinite series in Eq. (2) is truncated, then Eq. (2) can be written as

$$f(t) \simeq \sum_{n=1}^{N} \sum_{m=0}^{M-1} c(n, m) b(n, m, t) = C^{T} B(t),$$
 (3)

where

$$C = [c(1,0), \cdots, c(1,M-1), c(2,0), \cdots, c(2,M-1), \cdots, c(N,0), \cdots, c(N,M-1)]^T, \quad (4)$$

and

$$B(t) = [b(1,0,t), \cdots, b(1,M-1,t)|b(2,0,t), \cdots, b(2,M-1,t)| \cdots |b(N,0,t), \cdots, b(N,M-1,t)]^{T}.$$
(5)

2.2. The Operational Matrix of the Hybrid of Block-pulse and Legendre Polynomials.

The integration of the vector B(t) defined in Eq. (5) can approximated by

$$\int_0^t B(t')dt' \simeq PB(t) \tag{6}$$

where P is the  $NM \times NM$  operational matrix for integration and is given by

$$P = \begin{pmatrix} E & H & H & \cdots & H \\ 0 & E & H & \cdots & H \\ 0 & 0 & E & \cdots & H \\ \vdots & \vdots & \vdots & & \vdots \\ 0 & 0 & 0 & \cdots & E \end{pmatrix}. \tag{7}$$

In Eq. (7)

$$H = rac{t_f}{N} egin{pmatrix} 1 & 0 & 0 & \cdots & 0 \ 0 & 0 & 0 & \cdots & 0 \ 0 & 0 & 0 & \cdots & 0 \ dots & dots & dots & dots \ 0 & 0 & 0 & \cdots & 0 \end{pmatrix},$$

and E is operational matrix of integration for Legendre polynomials on the interval  $\left[\left(\frac{n-1}{N}\right)t_f,\frac{n}{N}t_f\right]$  given in [18] by

$$E = \frac{t_f}{2N} \begin{pmatrix} 1 & 1 & 0 & 0 & \cdots & 0 & 0 & 0 \\ \frac{-1}{3} & 0 & \frac{1}{3} & 0 & \cdots & 0 & 0 & 0 \\ 0 & \frac{-1}{5} & 0 & \frac{1}{5} & \cdots & 0 & 0 & 0 \\ \vdots & \vdots & \vdots & \vdots & & \vdots & \vdots & \vdots \\ 0 & 0 & 0 & 0 & \cdots & \frac{-1}{2M-3} & 0 & \frac{1}{2M-3} \\ 0 & 0 & 0 & 0 & \cdots & 0 & \frac{-1}{2M-1} & 0 \end{pmatrix}.$$

## 2.3. The Approximation of $B(t)B^{T}(t)C$ .

The following property of the product of two Legendre polynomial vectors will also be used.

Let

$$P(t) = [P_0(t), P_1(t), \dots, P_{M-1}(t)]^T,$$

$$A = [a_0, a_1, \dots, a_{M-1}]^T.$$

Then we have

$$P(t)P^{T}(t)A = \tilde{A}P^{T}(t), \tag{8}$$

where  $\tilde{A}$  is an  $M \times M$  matrix given in [18].

Let

$$B_n(t) = [b(n,0,t),b(n,1,t),\ldots,b(n,M-1,t)]^T, \quad n = 1,2,\ldots,N,$$

$$\bar{C}_n = [c(n,0),c(n,1),\ldots,c(n,M-1)]^T, \quad n = 1,2,\ldots,N.$$

Then using Eqs. (4) and (5) we get

$$B(t) = [B_1(t), B_2(t), \cdots, B_N(t)]^T,$$
 (9)

$$C = [\bar{C}_1, \bar{C}_2, \cdots, \bar{C}_N]^T. \tag{10}$$

By using Eqs. (9) and (10) we obtain

$$B(t)B^{T}(t)C = \begin{pmatrix} B_{1}(t)B_{1}^{T}(t)\bar{C}_{1} & 0 & \cdots & 0 \\ 0 & B_{2}(t)B_{2}^{T}(t)\bar{C}_{2} & \cdots & 0 \\ \vdots & \vdots & \ddots & \vdots \\ 0 & 0 & \cdots & B_{N}(t)B_{N}^{T}(t)\bar{C}_{N} \end{pmatrix}.$$
(11)

Similarly to Eq. (8) we have

$$B_n(t)B_n^T(t)\bar{C}_n = \tilde{C}_n B_n(t), \qquad n = 1, 2, \cdots, N.$$
 (12)

Using Eqs. (11) and (12), we get

$$B(t)B^{T}(t)C = \tilde{C}B(t), \tag{13}$$

where  $\tilde{C}$  is an  $NM \times NM$  diagonal matrix given by

$$\tilde{C} = \begin{pmatrix} \tilde{C}_1 & 0 & \cdots & 0 \\ 0 & \tilde{C}_2 & \cdots & 0 \\ \vdots & \vdots & \ddots & \vdots \\ 0 & 0 & \cdots & \tilde{C}_N \end{pmatrix}.$$

# 2.4. Integration of $B(t)B^{T}(t)$ .

The integration of the cross product of two hybrid Legendre vectors can be obtained as

$$D = \int_0^{t_f} B(t)B^T(t)dt. \tag{14}$$

where D is a diagonal matrix, given by

$$D = \begin{pmatrix} L & 0 & \cdots & 0 \\ 0 & L & \cdots & 0 \\ \vdots & \vdots & \ddots & \vdots \\ 0 & 0 & \cdots & L \end{pmatrix}, \tag{15}$$

with L the  $M \times M$  diagonal matrix given by

$$L = \frac{t_f}{N} \begin{pmatrix} 1 & 0 & \cdots & 0 \\ 0 & \frac{1}{3} & \cdots & 0 \\ \vdots & \vdots & \ddots & \vdots \\ 0 & 0 & \cdots & \frac{1}{2M-1} \end{pmatrix}.$$

# 3. Properties of Hybrid Functions of Block-Pulse and Chebyshev Polynomials.

Hybrid functions  $\hat{b}(n, m, t)$ ,  $n = 1, 2, \dots, N, m = 0, 1, \dots, M - 1$ , have three arguments; n is the order of block-pulse functions, m is the order of Chebyshev polynomials, and t is the normalized time. They are defined on the interval  $[0, t_f)$  as

$$\hat{b}(n,m,t) = \begin{cases} T_m(\frac{2N}{t_f}t - 2n + 1), & t \in \left[\left(\frac{n-1}{N}\right)t_f, \frac{n}{N}t_f\right) \\ 0, & \text{otherwise.} \end{cases}$$
(16)

Here  $T_m(t)$  are the well-known Chebyshev polynomials of order m which are orthogonal with respect to the weight function  $w(t) = \frac{1}{\sqrt{1-t^2}}$  and satisfy the following recursive formula.

$$T_{o}(t) = 1$$
,  $T_{1}(t) = t$   
 $T_{m+1}(t) = 2tT_{m}(t) - T_{m-1}(t)$  ,  $m = 1, 2, 3, ...$ 

Since  $\hat{b}(n, m, t)$  consists of block-pulse functions and Chebyshev polynomials, which are both complete and orthogonal, the set of the hybrid functions of block-pulse and Chebyshev polynomials is a complete orthogonal set.

### 3.1. Function Approximation.

A function f(t), defined over the interval 0 to  $t_f$  may be expanded as

$$f(t) = \sum_{n=1}^{\infty} \sum_{m=0}^{\infty} c(n, m) \hat{b}(n, m, t),$$
 (17)

where

$$c(n,m) = (f(t), \hat{b}(n,m,t))$$

in which (.,.) denotes the inner product. If the infinite series in Eq. (17) is truncated then Eq. (17) can be written as

$$f(t) \simeq \sum_{n=1}^{N} \sum_{m=0}^{M-1} c(n,m)\hat{b}(n,m,t) = C^{T}\hat{B}(t),$$
 (18)

where

$$C = [c(1,0), \cdots, c(1,M-1)|c(2,0), \cdots, c(2,M-1) \\ |\cdots|c(N,0), \cdots, c(N,M-1)]^T,$$

$$\hat{B}(t) = [b(1,0,t), \dots, b(1,M-1,t)|b(2,0,t), \dots, b(2,M-1,t)| \dots |b(N,0,t), \dots, b(N,M-1,t)|^{T}.$$
(19)

3.2. The Operational Matrix of the Hybrid of Block-pulse and Chebyshev Polynomials

The integration of the vector B(t) defined in Eq. (19) can approximated by

$$\int_0^t B(t')dt' \simeq \hat{P}\hat{B}(t)$$

where  $\hat{P}$  is the  $NM \times NM$  operational matrix for integration and is

given by

$$\hat{P} = \begin{pmatrix} \hat{E} & \hat{H} & \hat{H} & \cdots & \hat{H} \\ 0 & \hat{E} & \hat{H} & \cdots & \hat{H} \\ 0 & 0 & \hat{E} & \cdots & \hat{H} \\ \vdots & \vdots & \vdots & & \vdots \\ 0 & 0 & 0 & \cdots & \hat{E} \end{pmatrix}.$$

In the above matrix

$$\hat{H} = \frac{t_f}{N} \begin{pmatrix} 1 & 0 & 0 & \cdots & 0 \\ 0 & 0 & 0 & \cdots & 0 \\ \frac{-1}{3} & 0 & 0 & \cdots & 0 \\ \vdots & \vdots & \vdots & & 0 \\ \frac{(-1)^{M-1}}{2M(M-2)} & 0 & 0 & \cdots & 0 \end{pmatrix},$$

and E is operational matrix of integration for Chebyshev polynomials on the interval  $\left[\left(\frac{n-1}{N}\right)t_f,\frac{n}{N}t_f\right]$  given in [11] by

$$\hat{E} = \frac{t_f}{N} \begin{pmatrix} \frac{1}{2} & \frac{1}{2} & 0 & 0 & \cdots & 0 & 0 & 0 \\ \frac{-1}{8} & 0 & \frac{1}{8} & 0 & \cdots & 0 & 0 & 0 \\ \frac{-1}{6} & \frac{-1}{4} & 0 & \frac{1}{12} & \cdots & 0 & 0 & 0 \\ \vdots & \vdots \\ \frac{(-1)^{M-1}}{2(M-1)(M-3)} & 0 & 0 & 0 & \cdots & \frac{-1}{4(M-3)} & 0 & \frac{1}{4(M-1)} \\ \frac{(-1)^M}{2M(M-2)} & 0 & 0 & 0 & \cdots & 0 & \frac{-1}{4(M-2)} & 0 \end{pmatrix}.$$

# 4. Hybrid Functions Direct Method

For now, we will use hybrid of block-pulse and Legendre polynomials, similar results can be obtained by using hybrid of block-pulse and Chebyshev polynomials. Consider the problem of finding the extremum of the functional

$$J(x) = \int_0^1 F[t, x(t), \dot{x}(t)] dt.$$
 (20)

The necessary condition for x(t) to extremize J(x) is that it should satisfy the Euler-Lagrange equation

$$\frac{\partial F}{\partial x} - \frac{d}{dt} \left( \frac{\partial F}{\partial \dot{x}} \right) = 0 \tag{21}$$

with appropriate boundary conditions. However, the above differential equation can be integrated easily only for simple cases. Thus numerical and direct methods such as the well-known Ritz and Galerkin methods have been developed to solve variational hybrid functions. problems. Here we consider a Ritz direct method for solving Eq. (21) using the

Suppose, the rate variable  $\dot{x}(t)$  can be expressed as

$$\dot{x}(t) = C^T B(t). \tag{22}$$

Using Eq. (6), x(t) can be represented as

$$x(t) = \int_0^t \dot{x}(t')dt' + x(0)$$
  
=  $C^T PB(t) + [x(0), 0, \dots, 0, x(0), 0, \dots, 0, \dots, x(0), 0, \dots, 0]^T B(t)$ . (23)

We can also express t in terms of B(t) as

$$t = \left[\frac{1}{2N}, \frac{1}{2N}, 0, \dots, 0, \frac{3}{2N}, \frac{1}{2N}, \dots, 0, \dots, 0, \dots, 0, \dots, \frac{2N-1}{2N}, \frac{1}{2N}, 0, \dots, 0\right] B(t) = d^T B(t) \quad (24)$$

Substituting Eqs. (22-24) in Eq. (20), the functional J(x) becomes a function of c(n,m),  $n=1,2,\cdots,N$ ,  $m=0,1,2,\cdots,M-1$ . Hence to find the extremum of J(x) we solve

$$\frac{\partial J}{\partial c(n,m)}=0, \qquad n=1,2,\cdots,N, \qquad m=0,1,\cdots,M-1. \tag{25}$$

The above procedure is now used to solve the following examples.

#### 5. Illustrative Examples

In this section two problems of the calculus of variations are considered. Example 1 is the classical brachistochrone problem, where as example 2 is an application to the heat conduction problem taken from [7].

#### 5.1. Example 1: The Brachistochrone Problem.

#### 5.1.1. The Brachistochrone Problem as an Optimal Control Problem

As an optimal control problem, the brachistochrone problem may be formulated as [5].

Minimize the performance index J,

$$J = \int_0^1 \left[ \frac{1 + U^2(t)}{1 - X(t)} \right]^{\frac{1}{2}} dt, \tag{26}$$

subject to

$$\dot{X}(t) = U(t),\tag{27}$$

with

$$X(0) = 0, \quad \dot{X}(1) = -0.5.$$
 (28)

Eqs. (26), (27) and (28), describe the motion of a bead sliding down a frictionless wire in a constant gravitational field. The minimal time transfer expression is obtained from the law of conservation of energy. Here X and t are dimensionless and they represent respectively the vertical and horizontal coordinates of the sliding bead.

As is well known the exact solution to the brachistochrone problem is the cycloid defined by the parametric equations

$$x = 1 - \frac{\beta}{2}(1 + \cos 2\alpha), \qquad t = \frac{t_0}{2} + \frac{\beta}{2}(2\alpha + \sin 2\alpha),$$
 (29)

where

$$\tan \alpha = \frac{dX}{dt} = U.$$

With the given boundary conditions, the integration constants are found to be

$$\beta = 1.6184891, \quad t_0 = 2.7300631.$$

## 5.1.2. The Numerical Method

Suppose, the rate variable  $\dot{X}(t)$  can be expressed approximately as

$$\dot{X}(t) = C^T B(t). \tag{30}$$

Using Eqs. (6) and (28), X(t) can be represented as

$$X(t) = \int_0^t \dot{X}(t')dt' + X(0)$$
  
=  $C^T PB(t)$ , (31)

and by using Eqs. (27) and (30) we have

$$U^{2}(t) = C^{T}B(t)B^{T}(t)C.$$
 (32)

Equation (32) can be simplified by using the property of the product of two hybrid Legendre function vectors given in Eq. (13).

# 5.1.3. The Performance Index Approximation

Using Eqs. (26), (31) and (32) the performance index J can be approximated as follows:

$$J = \int_0^1 \left( \frac{1 + C^T \tilde{C} B(t)}{1 - C^T P B(t)} \right)^{\frac{1}{2}} dt.$$
 (33)

Divide the interval [0,1] into N equal subintervals, we have

$$J = \sum_{n=1}^{N} \int_{\frac{n-1}{N}}^{\frac{n}{N}} \left( \frac{1 + C^T \tilde{C}B(t)}{1 - C^T PB(t)} \right)^{\frac{1}{2}} dt.$$
 (34)

In order to use the Gaussian integration formula we transform the tinterval  $(\frac{n-1}{N}, \frac{n}{N})$  into the  $\tau$ -interval (-1, 1) by means of the transformation

$$t = \frac{1}{2}(\frac{1}{N}\tau + \frac{2n-1}{N}). \tag{35}$$

The optimal control problem in Eqs. (26-28) is then restated as follows: Minimize

$$J = \frac{1}{2} \int_{-1}^{1} \left[ \frac{1 + u^{2}(\tau)}{1 - x(\tau)} \right]^{\frac{1}{2}} d\tau, \tag{36}$$

subject to

$$\frac{dx}{d\tau} = \frac{1}{2}u(\tau),\tag{37}$$

with

$$x(-1) = 0, \quad x(1) = -0.5.$$
 (38)

Using Eqs. (34) and (35) we get

$$J = \sum_{n=1}^{N} \frac{1}{2N} \int_{-1}^{1} \left( \frac{1 + C^{T} \tilde{C} B(\frac{1}{2}(\frac{1}{N}\tau + \frac{2n-1}{N}))}{1 - C^{T} P B(\frac{1}{2}(\frac{1}{N}\tau + \frac{2n-1}{N}))} \right)^{\frac{1}{2}} d\tau.$$
 (39)

Using the Gaussian integration formula, Eq.(39) can be approximated as

$$J \approx \sum_{n=1}^{N} \frac{1}{2N} \sum_{j=0}^{k} \left( \frac{1 + C^{T} \tilde{C} B(\frac{1}{2} (\frac{1}{N} \tau_{j} + \frac{2n-1}{N}))}{1 - C^{T} P B(\frac{1}{2} (\frac{1}{N} \tau_{j} + \frac{2n-1}{N}))} \right)^{\frac{1}{2}} w_{j}, \tag{40}$$

where  $\tau_j$ , j = 0, 1, ..., k are the k+1 zeros of Legendre polynomial  $P_{k+1}$ , and  $w_j$  are the corresponding weights, given in [19]. The idea behind the above approximation is the exactness of the Gaussian integration formula for polynomials of degree not exceeding 2k+1.

# 5.1.4. Evaluating the Vector C

The optimal control problem has now been reduced to a parameter optimization problem which can be stated as follows.

Find c(n, m),  $n = 1, 2, \dots, N$ ,  $m = 0, 1, \dots, M-1$  that minimizes Eq.(40) subject to

$$x(-1) = 0, \quad x(1) = -0.5.$$
 (41)

We now minimize Eq. (40) subject to Eq. (41) using the Lagrange multiplier technique. Suppose

$$J^* = J + \lambda_1 x(-1) + \lambda_2 [x(1) + 0.5].$$

The necessary conditions for a minimum are

$$\frac{\partial J^*}{\partial c(n,m)} = 0$$
  $n = 1, 2, \dots, N, m = 0, 1, \dots, M-1$  (42)

and

$$\frac{\partial J^*}{\partial \lambda_1} = 0, \qquad \frac{\partial J^*}{\partial \lambda_2} = 0. \tag{43}$$

Eqs. (42) and (43) give (NM+2) non-linear equations with (NM+2) unknowns which can be solved for c(n,m),  $\lambda_1$  and  $\lambda_2$  using Newton's iterative method. The initial values required to start Newton's iterative method have been chosen by taking  $x(\tau)$  as a linear function between x(-1) = 0 and x(1) = -0.5.

In Table 1 the results for hybrid Legendre approximation with N=2, k=5 and M=3, 4,5 together with N=2, k=8 and M=5 are listed, we compare the solution obtained using the proposed method with other solutions in the literature together with the exact solution.

Methods	x(1)	u(-1)	J
Dynamic programming gradient method[2]	-0.5	-0.7832273	0.9984988
Dynamic programming successive sweep method[3,4]	-0.5	-0.7834292	0.9984989
Chebyshev solutions[5]			
M=4	-0.5	-0.7844893	0.9984982
M=7	-0.5	-0.7864215	0.99849815
M = 10	-0.5	-0.7864406	0.9984981483
Hybrid Legendre, $N = 2, k = 5$			
M=3	-0.5	-0.7852418	0.9984989
M=4	-0.5	-0.7864397	0.9984983
M=5	-0.5	-0.7864402	0.9984981
Hybrid Legendre $N = 2, k = 8 \text{ and } M = 5$	-0.5	-0.7864408	0.99849814829
Exact Solution[4]	-0.5	-0.7864408	0.99849814829

Table 1. The hybrid Legendre and other solutions in the literature.

## 5.2. Example 2: Application to The Heat Conduction Problem

Consider the extremization of

$$J = \int_0^1 \left[ \frac{1}{2} \dot{x}^2 - x g(t) \right] dt = \int_0^1 F(t, x, \dot{x}) dt, \tag{44}$$

where g(t) is a known function satisfying

$$\int_0^1 g(t)dt = -1,$$

with the boundary conditions

$$\dot{x}(0) = 0$$
 ,  $\dot{x}(1) = 0$ . (45)

Schechter [20] gave a physical interpretation for this problem by noting an application in heat conduction and Chen and Hsiao [7] considered the case where g(t) is given by

$$g(t) = \begin{cases} -1, & 0 \le t < \frac{1}{4}, & \frac{1}{2} \le t < 1, \\ 3, & \frac{1}{4} \le t < \frac{1}{2}, \end{cases}$$
 (46)

and gave an approximate solution using Walsh functions. The exact solution is

$$x(t) = \begin{cases} \frac{1}{2}t^2, & 0 \le t < \frac{1}{4} \\ -\frac{3}{2}t^2 + t - \frac{1}{8}, & \frac{1}{4} \le t < \frac{1}{2} \\ \frac{1}{2}t^2 - t + \frac{3}{8}, & \frac{1}{2} \le t < 1. \end{cases}$$

Here of we solve the same problem using hybrid of Legendre and blockpulse functions with M=3 and N=4. First we assume

$$\dot{x}(t) = C^T B(t). \tag{47}$$

In view of Eq. (46), we write Eq. (44) as

$$J = \frac{1}{2} \int_0^1 \dot{x}^2(t)dt + 4 \int_0^{\frac{1}{4}} x(t)dt - 4 \int_0^{\frac{1}{2}} x(t)dt + \int_0^1 x(t)dt,$$

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$$\begin{split} J = & \frac{1}{2} \int_0^1 C^T B(t) B^T(t) C dt + 4 C^T P \int_0^{\frac{1}{4}} B(t) dt - \\ & 4 C^T P \int_0^{\frac{1}{2}} B(t) dt + C^T P \int_0^1 B(t) dt \,. \end{split}$$

Let

$$w(t) = \int_0^t B(t')dt',$$

then using Eq. (20), we have

$$J = \frac{1}{2}C^TDC + C^TP[4w(\frac{1}{4}) - 4w(\frac{1}{2}) + w(1)].$$
 (48)

where

$$D = \int_0^1 B(t)B^T(t)dt.$$

The boundary conditions in Eq.(45) can be expressed in terms of hybrid of Legendre and block-pulse functions as

$$C^T B(0) = 0$$
 ,  $C^T B(1) = 0$ . (49)

We now minimize Eq. (48) subject to Eq. (49) using the Lagrange multiplier technique. Suppose

$$J^* = J + \lambda_1 C^T B(0) + \lambda_2 C^T B(1), \tag{50}$$

where  $\lambda_1$  and  $\lambda_2$  are the two multipliers. Using Eq. (25) we obtain

$$\frac{\partial J^*}{\partial C} = DC + P[4w(\frac{1}{4}) - 4w(\frac{1}{2}) + w(1)] + \lambda_1 B(0) + \lambda_2 B(1) = 0.$$
(51)

We also have

$$w(1) = \frac{1}{4}[1,0,0,1,0,0,1,0,0,1,0,0]^{T},$$

$$w(\frac{1}{2}) = \frac{1}{4}[1,0,0,1,0,0,0,0,0,0,0]^{T},$$

$$w(\frac{1}{4}) = \frac{1}{4}[1,0,0,0,0,0,0,0,0,0,0]^{T},$$

$$B(0) = [1,-1,1,0,0,0,0,0,0,0,0]^{T},$$

$$B(1) = [0,0,0,0,0,0,0,0,1,1,1]^{T}.$$

Equations (49) and (51) define a set of 14 simultaneous linear algebraic equations from which the coefficient vector C and the multipliers  $\lambda_1$  and  $\lambda_2$  can be found. The vector  $C^T P$  is

$$C^{T}P = \frac{1}{64} \left[ \frac{2}{3}, 1, \frac{1}{3}, 2, -1, -1, -\frac{10}{3}, -3, \frac{1}{3}, -\frac{22}{3}, -1, \frac{1}{3} \right]^{T}$$
 (52)

Further, to define x(t) for t in the interval  $[0, \frac{1}{4}]$  we map  $[0, \frac{1}{4}]$  into [-1, 1] by mapping t into 8t-1 and similarly for the other intervals. Using the above equation and  $P_0 = 1$ ,  $P_1 = t$  and  $P_2 = \frac{3}{2}t^2 - \frac{1}{2}$ , we get

$$x(t) = \begin{cases} \frac{1}{64} \left[ \frac{2}{3} + (8t - 1) + \frac{1}{3} \left[ \frac{3}{2} (8t - 1)^2 - \frac{1}{2} \right] \right] = \frac{1}{2} t^2, & 0 \le t \le \frac{1}{4} \\ \frac{1}{64} \left[ 2 - (8t - 3) - \left[ \frac{3}{2} (8t - 3)^2 - \frac{1}{2} \right] \right] = -\frac{3}{2} t^2 + t - \frac{1}{8}, & \frac{1}{4} \le t \le \frac{1}{2} \\ \frac{1}{64} \left[ -\frac{10}{3} - 3(8t - 5) + \frac{1}{3} \left[ \frac{3}{2} (8t - 5)^2 - \frac{1}{2} \right] \right] = \frac{1}{2} t^2 - t + \frac{3}{8}, & \frac{1}{2} \le t \le \frac{3}{4} \\ \frac{1}{64} \left[ -\frac{22}{3} - (8t - 7) + \frac{1}{3} \left[ \frac{3}{2} (8t - 7)^2 - \frac{1}{2} \right] = \frac{1}{2} t^2 - t + \frac{3}{8}, & \frac{3}{4} \le t \le 1 \end{cases}$$

which is the exact solution. This exact solution can not be obtained either with CBF's or with PCBF's.

#### 6. Conclusion

The aim of present work is to develop an efficient and accurate method for solving problems of the calculus of variations. The problem has been reduced to solving a system of algebraic equations. Illustrative examples are included to demonstrate the validity and applicability of the technique. The advantages of using the hybrid Legendre method are:

- (1) The operational matrix P contains many zeros which plays an important role in simplifying the performance index.
- (2) The Gaussian integration formula is exact for polynomials of degree not exceeding 2k + 1.
- (3) Only small values of k, N and M are needed to obtain very satisfactory results for the brachistochrone problem.

(4) Hybrid functions approach provides an exact solution for the heat conduction problem.

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Proceedings of the Monthly Colloquium of the Iranian Mathematical Society Vol. 2, pp 179-200 (2004)

## The Algebra of Formal Power Series

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Abstract: Let K be a field. K[[X]] will denote the ring of formal power series in several commuting variables,  $X = (x_1, x_2, \dots, x_k)$  with coefficients in K. K((X)) will denote the field of fractions of K[[X]]. An element  $f \in K((X))$  is said to be an algebraic function over K if f is algebraic over the field of rational functions K(X). If further  $f \in K[[X]]$ , then f is said to be an algebraic series over K. A function which is not algebraic is called transcendental (over K).

For the case of one variable G. Christol et. al. (1980) have characterised the algebraic functions over a finite field in terms of automata. We generalise their argument and obtain the corresponding result for the case of several variables over a perfect field of positive characteristic. H. Furstenberg (1967) has shown that if K is a finite field and  $f = \sum_{n \geq 0} a_n x^n$ ,  $g = \sum_{n \geq 0} b_n x^n \in K[[X]]$  are aglebraic, then the Hadamard product of f and g,  $f * g = \sum_{n \geq 0} a_n b_n x^n$  is also algebraic. We apply the above characterisation of the algebraic series in several variables to generalise Furstenberg's result and prove that if K is an arbitrary field of positive characteristic and if  $f = \sum_{l} a_l X^l$ ,  $g = \sum_{l} b_l X^l \in K[[X]]$  are algebraic series, then  $f * g = \sum_{l} a_l b_l X^l$  is also an algebraic series. As an easy consequence of the above result we give a proof to

MSC (2000): 11J61, 12H25

Keywords: Differentially algebraic formal power series algebraically independent, transcendence basis, Hadamard product.

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Deligne's Theorem: If  $f \in K[[X]]$  and  $f = \sum_{\sigma} a_{\sigma} X^{\sigma}$  is an algebraic series in X over K, then  $D(f) = \sum_{n \geq 0} a_{n,1} t^n$  is an algebraic series in t over K. Then we consider the Hadamard product of rational formal power series and we show that if L is a field of characteristic zero and if  $f = \sum_{l} a_{l} X^{l}$ ,  $g = \sum_{l} b_{l} X^{l}$  are rational series in L[[X]], then  $f * g = \sum_{l} a_{l} b_{l} X^{l}$  (which is not in general a rational series) is always algebraic only if  $k \geq 2$ .

We introduce two methods for trying to decide whether or not a given formal power series is an algebraic series over a field. Our first method (in characteristic zero) is based on the reduction process modulo the prime p and our second method (in positive characteristic and so in characteristic zero using reduction) is based on the splitting process for functions.

In 1986, M. Mendes France and A. J. van der Poorten have shown that if  $f = \sum_{n=0}^{\infty} a_n x^n \in F[[X]]$  is algebraic, where F is a finite field of characteristic p > 0,  $a_0 = 1$  and  $f \neq 1$  and if  $\lambda$  is a p-adic integer, then  $f^{\lambda}$  is algebraic if and only if  $\lambda$  is rational. We generalise their result and prove the following theorem:

Suppose that K is a field of characteristic p > 0. Suppose that  $f = \sum_{n=0}^{\infty} a_n x^n \in K[[x]]$  is algebraic over K, where  $a_0 = 1$  and  $a_1 \neq 0$ . Let  $\lambda_1, \lambda_2, \dots, \lambda_n$  be p-adic integers. Then the following conditions are equivalent:

- (i)  $1, \lambda_1, \lambda_2, \dots, \lambda_n$  are linearly independent over Q.
- (ii)  $(1+x)^{\lambda_1}, (1+x)^{\lambda_2}, \cdots, (1+x)^{\lambda_n}$  are algebraically independent over Q.
- (iii)  $f^{\lambda_1}, f^{\lambda_2}, \dots, f^{\lambda_n}$  are algebraically independent over K(x).

A formal power series f is called D-Algebraic (respectively, D-finite) if it satisfies an algebraic (respectively, a linear) differential equation with polynomial coefficients. These notions are usually defined only over fields of characteristic zero and are not so significant over fields of characteristic p>0 as  $f^{(p)}\equiv 0$ . For a formal power series over a perfect field of positive characteristic we make the definition of E-algebraicity (respectively, E-finiteness) which is an analogue of the notion of D-algebraicity (respectively, D-finiteness).

It is slightly surprising that E-finite series are in fact algebraic series. However, the E-algebraic series (which are not all algebraic series) under ordinary addition and multiplication of series, form a field which is algebraically closed in K((x)) and has some other natural prperties. We also study the Hadamard product of two E-algebraic formal power series.

### 1. Preliminaries and Notations

Let K be a field and  $X = (x_1, x_2, \ldots, x_k)$ . We will denote the ring of formal power series by K[[X]], the field of fractions of K[[X]] by K((X)). An element  $f \in K((X))$  is said to be an algebraic function over K if f is algebraic over the field of rational functions K(X). If, further,  $f \in K[[X]]$ , then f is said to be an algebraic series over K. A function (or series) which is not algebraic is called transcendental over K.

Example 1.1. (i) The series

$$f = \sum_{n=0}^{\infty} {2n \choose n} x^n = \sum_{n=0}^{\infty} {-\frac{1}{2} \choose n} (-4)^n x^n = (1-4x)^{-\frac{1}{2}}$$

is algebraic over any field.

- (ii) The series  $f(x) = \sum_{n=0}^{\infty} x^{p^n}$  is algebraic over any field of characteristic p and transcendental over any field of characteristic q, where  $q \neq p$ .
- (iii) The series  $\sum_{n=1}^{\infty} x^{n!}$  is transcendental over any field (see Zariski and Samuel [25, p. 220]).
- (iv) Let p be a prime number and let  $S_p(n)$  be the sum of the digits of n is its p-adic expansion. That is, if

$$n = \sum_{i=0}^{\infty} n_i p^i, \qquad 0 \le n_i \le p - 1,$$

(actually a finite sum for  $n \in N$ ) is the *p*-adic expansion of n, then  $S_p(n) = \sum_{i=0}^{\infty} n_i$ . One can show that if

$$f(x) = \sum_{n=0}^{\infty} S_p(n) x^n \in F_p[[x]],$$

then  $f(x) = \frac{x^p-1}{x-1}f(x)^p + \frac{x}{(x-1)^2}$ . That is, f is algebraic over  $F_p$  (see [19]).

# 2. Some characterisations of algebraic series

The following theorem is a well-known result in the field of complex numbers.

Theorem (Biberbakh). Let  $f = \sum_{n=0}^{\infty} a_n z^n \in C[[z]]$ , where  $a_n$  just accepts the finite elements  $d_1, d_2, \ldots, d_r$ . Then f is rational or non-algebraic.

G. Christol et.al.[4] have the following characterisation of algebraic series in terms of p-automata.

**Theorem 2.1.** Suppose that F is a finite field of characteristic p. Suppose that  $f = \sum_{n=0}^{\infty} a_n x^n \in F[[x]]$ . Then the following are equivalent:

- i) f is algebraic over F.
- ii) The sequence  $(a_n)$  is generated by a p-automata.

When F is replaced by the ring of p-adic integers, Denef and Lipshitz in [5] have extended Theorem 2.1 to the case of several variables.

We gave a characterisation of algebraic series in several variables over perfect fields of positive characteristic in [15].

Recall that a field K is perfect if every irreducible polynomial over K is separable. Equivalently, every algebraic extension field of K is separable over K. For example, every finite field or every field of characteristic zero is perfect. If K is a field of characteristic p, then K is perfect if and only if the Frobenuis map  $\phi: K \longrightarrow K$  is an automorphism.

From now on, K will denote a perfect field of characteristic p > 0, unless explicity stated otherwise.

Firstly, we shall recall the splitting process for functions over a perfect field and the associated semilinear operators on the field of fractions of the ring of formal power series which we introduced in [15] and secondly, we shall employ such operators for characterising algebraic

functions (for the details of the proofs see [15] or [19]).

**Theorem 2.2.** Suppose that K is a field. If  $f \in K((X))$  is an algebraic function over L, where L is an extension of K, then f is algebraic over K.

The above theorem asserts that when we deal with algebraic functions over arbitrary fields of positive characteristic it is enough to consider such functions over perfect fields of positive characteristic, since every field L has a perfect extension, for example, the algebraic closure of L.

Let  $\iota$  be a non-negative vector, that is,  $\iota=(n_1,n_2,\cdots,n_k)$ , where  $i_j\in N, j=1,2,\cdots,k$ . Then  $X^\iota$  will denote the monomial  $x_1^{n_1}x_2^{n_2}\cdots x_k^{n_k}$ . We denote by  $\Lambda$  the set of all non-negative vectors, and by  $\Lambda_p$  the set  $Z_p^k$ , where  $Z_p=\{0,1,2,\cdots,p-1\}$ .

**Lemma 2.3** If  $f(X) \in K[[X]]$  (respectively K((X))), then f can be written uniquely as

$$f = \sum_{i \in \Lambda_n} X^i f_i^p \tag{2.1}$$

for some  $f_{\iota} \in K[[X]]$  (respectively K((X))).

For  $\iota \in \Lambda_p$  define  $E_{\iota} : K((X)) \longrightarrow K((X))$  by

$$E_{\iota}(f) = f_{\iota}. \tag{2.2}$$

Now for  $f \in K((X))$ , by Lemma 2.3 we have

$$f = \sum_{\iota \in \Lambda_p} X^{\iota} (E_{\iota}(f))^p. \tag{2.3}$$

The operator  $E_{\iota}$  is semilinear; that is, if  $f, g \in K((X))$  and  $\lambda \in K$ , then

$$E_{\iota}(\lambda f + g) = \lambda^{\frac{1}{p}} E_{\iota}(f) + E_{\iota}(g).$$

Moreover,  $E_{\iota}(g^p f) = g E_{\iota}(f)$ .

**Definition 2.4** Suppose that  $f, g \in K[[X]]$ , say

$$f = \sum_{t \in \Lambda} a_t X^t, \quad g = \sum_{t \in \Lambda} b_t X^t.$$

The Hadamard product of f and g, which will be denoted by  $f \ast g$ , is the series which is defined by

$$f * g = \sum_{t \in \Lambda} a_t b_t X^t.$$

**Lemma 2.5** If  $f, g \in K[[X]]$ , then for  $t \in \Lambda_p$ ,

$$E_t(f*g) = E_t(f)*E_t(g).$$

Let  $\Omega$  be the semigroup generated by the identity operator and the  $E_{\iota}$  for  $\iota \in \Lambda_p$ , with ordinary composition as multiplication. To each  $f \in K((X))$  we associate its orbit

$$\Omega(f) = \{ E(f) : E \in \Omega \}. \tag{2.4}$$

Then we have the following:

Theorem 2.6. Let  $f \in K((X))$ . Then  $< \Omega(f) >$ , the K-linear space spanned by  $\Omega(f)$ , is the smallest K-subspace of K((X)) containing f and which is invariant under each  $E_{\iota}$ ,  $\iota \in \Lambda_{\mathfrak{p}}$ .

**Theorem 2.7.** Let  $f \in K((X))$ . Then the following are equivalent:

- i) f is algebraic over K.
- ii) There exist elements  $a_o, a_1, \ldots, a_N$  in K[X] such that

$$\sum_{i=0}^{\infty} a_i f^{p^i} = 0,$$

where  $a_o \neq 0$ .

iii) There exists a finite dimensional K-subspace V of K((X)) such that  $f \in V$  and  $E_{\iota}(V) \subseteq V$ ,  $\iota \in \Lambda_p$ .

iv) 
$$dim_K < \Omega(f) > is finite.$$

**Example 2.8.** Let  $f = \sum_{n\geq 0} x^n y^{n^2}$ . We shall show that f is transcendental over  $F_2(x, y)$  and hence f will be transcendental with respect to any field of characteristic zero too.

Using the idea of Lemma 2.3 we split f and find the series

$$f_{1} = E_{(o,o)}(f) = \sum_{n\geq 0} x^{n} y^{2n^{2}},$$

$$f_{2} = E_{(o,o)(o,o)}(f) = \sum_{n\geq 0} x^{n} y^{2^{2}n^{2}},$$

$$\cdots,$$

$$f_{k} = E_{(o,o)(o,o)\dots(o,o)}(f) = \sum_{n\geq 0} x^{n} y^{2^{k}n^{2}}$$

Each series  $f_i$  corresponds to the set  $\{(n,m): m=2^in^2\}$ . These sets are all clearly distinct and hence the set  $\{f_1, f_2, f_3, \dots\}$  is infinite. Therefore, f is transcendental over  $F_2(x,y)$  by Theorem 2.7.

Corollary 2.9. Suppose that  $f,g \in K[[X]]$ . If f,g are algebraic series over K, then f \* g is again an algebraic series over K.

**Theorem 2.10.** Suppose that K is any field. If  $h \in K((X))$  is an algebraic function over L, where L is an extension field of K, then h is an algebraic function over K.

We are now in a position to have the main theorem.

**Theorem 2.11.** If K is a field of characteristic p > 0 and if f, g are algebraic series over K, then f \* g is again an algebraic series over K.

Deligne's theorem [4] can be proved directly from the main theorem.

Theorem 2.12. Suppose that K is a field of characteristic p > 0. If  $f \in K[[X]]$  and  $f = \sum_{\sigma \in \Lambda} a_{\sigma} X^{\sigma}$  is an algebraic series in X over K, then  $I(f) = \sum_{n \geq 0} a_{n,1} t^n$  is an algebraic series in t over K.

In Theorem 2.7 we considered  $< \Omega(f) >$ , the K-linear space spanned by  $\Omega(f)$ . Now we consider  $< \Omega(f) >$  as a K(X)-linear space spanned by  $\Omega(f)$ . We shall show that f is algebraic if and only if this space is a finite dimensional subspace of K((X)) too.

Theorem 2.13. Suppose that  $f \in K((X))$  and W is the K(X)-subspace of K((X)) spanned by  $\Omega(f)$ . Then f is an algebraic function if and only if  $\dim_{K(X)} W < \infty$ .

Recall that for a field extension  $L \subseteq F$ , a transcendence base of F over L is a subset S of F which is algebraically independent over L and is maximal (with respect to set-theoretic inclusion) in the set of all algebraically independent subsets of F.

From now on, for simplicity of notations, we just concentrate on series in one variable.

**Theorem 2.14.** Let  $f \in K((x))$ . Then f is algebraic over K if and only if  $K(x, \Omega(f))$  is a finitely generated field extension of K(x).

# 3. Some methods for transcendency

R.P. Stanley observed in [21] (via analytic techniques) that if  $f(x) = \sum_{n=0}^{\infty} \binom{2n}{n}^t x^n$ , t > 1, then f is transcendental over C(x) for even t, t > 1. He also stated that it is unknown for odd t > 1 whether or not

it is transcendental.

In this section apart from the method which was introduced in section 2, we introduce another method for deciding whether or not a given formal power series with integer coefficients is transcendental.

Our method (in characteristic zero) is based on the reduction process and looking at the degree of the corresponding explicit equation obtained when the reduction modulo the prime p of the formal power series is algebraic. If the degree of the corresponding polynomial equation is an unbounded function of the prime p, then there is no single polynomial for all primes p and therefore, the formal power series is transcendental.

We introduced our method by dealing with the above particular problem, which was raised by Stanley [21] and we could answer it in [23].

First we showed that (the reduction of) f is algebraic over any field of positive characteristic p and we then deduced (from the explicit equation obtained) that f is transcendental over any field of characteristic zero for any integer t > 1.

We also gave a generalisation of this result in the case of multinomial coefficients.

Note that if t=1, then f is algebraic of degree at most 2 over any field. In that case,  $f=(1-4x)^{-1/2}$  and for  $t\geq 1, f=1$  over any field of characteristic 2. Hence we may suppose that, in the case of positive characteristic p, p>2.

Throughout this section f will denote the series

$$f(x) = \sum_{n=0}^{\infty} \left( \begin{array}{c} 2n \\ n \end{array} \right)^{t} x^{n}$$

for  $t \in N$  and t > 1.

First we shall see that if a series h with integer coefficients is algebraic

over a field of characteristic zero, then the reduction of h is algebraic over  $F_p$ .

**Proposition 3.1** Suppose that K is any field of characteristic zero and

$$h(x) = \sum_{i=0}^{\infty} h_i x^i \in Z[[x]]$$

is algebraic over K of degree N. Then for any prime p,

$$\bar{h}(x) = \sum_{i=0}^{\infty} \bar{h_i} x^i \in F_p[[x]]$$

is algebraic over  $F_p$  of degree at most N, where  $\bar{a}$  is the image of a in  $F_p$ .

**Note.** Proposition 3.1 can be extended to the case of several variables.

Remark 3.2. Note that 
$$f(x) = \sum_{n=0}^{\infty} \binom{2n}{n}^t x^n = h * h * \cdots * h (t)$$
 times), where  $h(x) = \sum_{n=0}^{\infty} \binom{2n}{n} x^n$ , which is algebraic (with respect to any field) and \* denotes the Hadamard product operation.

Now, since over any field of positive characteristic the Hadamard product of two algebraic formal power series is again an algebraic formal power series (see [15, Collorary 5.5]) it follows that (the reduction of) f is algebraic over any field of positive characteristic. However, we shall now prove this directly by using Lucas' Theorem to find the corresponding explicit equation for (the reduction of) f, since we need this equation later.

**Theorem 3.3.** (Lucas' Theorem) For  $m, n \in N, p$  a prime,

$$\left(\begin{array}{c} m \\ n \end{array}\right)^p \equiv \left(\begin{array}{c} mp \\ np \end{array}\right) \equiv \left(\begin{array}{c} m \\ n \end{array}\right) \pmod{p}$$

and

$$\left(\begin{array}{c} mp+i \\ np+i \end{array}\right) \equiv \left(\begin{array}{c} m \\ n \end{array}\right) \left(\begin{array}{c} i \\ j \end{array}\right) \pmod{p}$$

for  $i, j \in N$  with  $0 \le i, j \le p - 1$ .

Proof. See, for example, Dickson [6, p. 271]. ///

**Proposition 3.4.** If p is any odd prime, then f is algebraic over  $\mathbb{Z}_p$ .

Theorem 3.5. Let  $t \in N, t > 1$ . If  $f(x) = \sum_{n=0}^{x} {2n \choose n}^{t} x^{n} \in \mathbb{Z}[[X]]$ , then f is transcendental over any field of characteristic zero.

**Theorem 3.6.** If  $g(x) = \sum_{m=0}^{\infty} {km \choose m, m, \dots, m}^t x^m$  where  $t, k \in N$  with  $t \ge 1, k \ge 3$ , then g is transcedental over any field of characteristic zero.

# 4. Rational power series

The following result, which is well known, is due to E. Borel.

Theorem 4.1. Let K be a field of characteristic zero. let  $f(x) = \sum_{n\geq 0} a_n x^n$ ,  $g(x) = \sum_{n\geq 0} b_n x^n \in K[[x]]$ . If f, g are rational, then  $f * g(x) = \sum_{n\geq 0} a_n x^n$  is again rational.

When the field K has positive characteristic the result is still true.

**Lemma 4.2.** Suppose that K is a field of characteristic p > 0 which is also algebraically closed. If f and g belong to K(x), then  $f * g \in K(x)$ .

**Proof.** see [16. Lemma 3.2].///

Lemma 4.3. Suppose that K is a field and L is an extension field of K. Suppose that  $f \in K[[x]]$ . If  $f \in L(x)$ , then  $f \in K(x)$ .

**Proof.** See [16, Lemma 3.3].///

**Theorem 4.4.** Suppose that K is a field of characteristic p > 0. If f and  $g \in K(x)$ , then  $f * g \in K(x)$ .

**Proof.** Suppose that L is an extension field of K which is algebraically closed. Then f and g belong to L(x). Therefore by Lemma 4.2,  $f * g \in L(x)$  and so by Lemma 4.3,  $f * g \in K(x)$ .

Theorems 4.1 and 4.4 do not hold in the case of several variables. For example if

$$f = \sum_{n,m \ge 0} \binom{n+m}{n} x^n y^m = \frac{1}{1-x-y},$$

which is rational, then (except in characteristic two)

$$f * f = \sum_{n,m \ge 0} {n+m \choose n}^2 x^n y^m = \{(1-x-y)^2 - 4xy\}^{-1/2},$$

which is not rational.

A question which arises here is

Suppose that K is a field and  $f, g \in K[[x_1, x_2, \dots, x_k]]$ . Suppose that f, g are rational series. Is the Hadamard product f \* g an algebraic series?

The answer is positive when the field K has positive characteristic [see [24, Corollary 5.5]]. If k > 2 and K has characteristic zero, then the following example shows that the answer is negative.

Example 4.5. If

$$f = \sum_{n_1, n_2, n_3 \ge 0} \left( \begin{array}{c} n_1 + n_2 + n_3 \\ n_1, n_2, n_3 \end{array} \right) x_1^{n_1} x_2^{n_2} x_3^{n_3} = \frac{1}{1 - x_1 - x_2 - x_3}$$

and

$$g = \sum_{n \ge 0} (x_1, x_2, x_3)^n = \frac{1}{1 - x_1 x_2 x_3},$$

which are rational series, then

$$f * g = \sum_{n \ge 0} \begin{pmatrix} 3n \\ n, n, n \end{pmatrix} t^n, \text{ where } t = x_1 x_2 x_3,$$

which is transcendental over a field of characteristic zero (see, for example, [23]). Note that, by [23], the same series is algebraic over a field of characteristic p > 0, but of degree equal to an unbounded function of p.

The only case which is left and still seems to be unknown is the case k=2 and when K has characteristic zero. In [24] we considered this case and proved the following theorem.

Theorem 4.6. The Hadamard product of two rational series of two variables over a field of characteristic zero is an algebraic series.

Remark 4.7. We cannot weaken the conditions of Theorem 4.6 to allow one of the series to be algebraic. For example if

$$R(x,y) = \sum_{n,m\geq 0} {n+m \choose n}^2 x^n y^m = \{(1-x-y)^2 - 4xy\}^{-1/2},$$

which is algebraic and

$$S(x,y) = \sum_{n\geq 0} (xy)^n = \frac{1}{1-xy},$$

which is rational, then

$$R * S = \sum_{n \ge 0} {2n \choose n}^2 t^n$$
, where  $t = xy$ ,

which is transcendental. (See [23].)

**Example 6.3.** Suppose that  $f = \sum_{n\geq 0} x^{n^2}$ , and  $g = \sum_{n\geq 0} \frac{1}{n!} x^n$ , which are D-algebraic. Then  $f * g = \sum_{n\geq 0} \frac{1}{(n^2)!} x^{n^2}$ , which is not D-algebraic. (See Lipshitz-Rubel [12, Proposition 7.3, p. 1209].)

A sub-algebra of  $D_L$  which is closed under the Hadamard product operation is the algebra of differentiably finite power series. A differentiably finite (D-finite, for short) power series is a series which satisfies a linear differential equation. The class of D-finite power series has been subject to extensive investigation, particularly whithin the theory of differential equations. A systematic exposition of their properties from a combinatorial point of view have been given by Stanley [21].

Theorem 6.4. Suppose that f, g are D-finite power series. Then f \* g is again a D-finite power series.

Proof. See Stanley [21].

Recently, Stanley's notion of D-finiteness has been of interest to several authors such as Gessel, Zeilberger, Lipshitz etc. Theorem 6.4 and many other results concerning D-finiteness have been generalised by Lipshitz [11] to the case of several variables.

Note that in Exampe 6.3, g is D-finite. Hence the Hadamard product of a D-algebraic and a D-finite power series is not D-algebraic. However, we have been able to prove that the Hadamard product of a rational formal power series and a D-algebraic formal power series is D-algebraic. First we need the following Lemma.

Lemma 6.5. Suppose that  $E \subseteq F$  is an algebraic extension of fields. Suppose that  $f \in E[[x]]$ . If  $tr.deg._FF(x, f, f', ..., f^{(n)}, ...) < \infty$ , then

$$tr.deg._{E}E(x,f,f^{'},...,f^{(n)},...)<\infty.$$

Theorem 6.6. Suppose that  $f, g \in L[[x]]$ . If f is rational and g is D-algebraic, then f \* g is D-algebraic.

### 7. E-algebraic functions

Differentially algebraic functions are usually defined only over fields of characteristic zero and are not so significant over fields of characteristic p > 0, as  $f^{(p)} \equiv 0$ . In this section we shall define an analogue of the concept of a D-algebraic function over a perfect field of characteristic p > 0.

From now on K will denote a perfect field of characteristic p > 0, unless explicitly stated otherwise. Moreover, all we discuss about can be generalised to the case of several variables.

**Definition.** 7.1 Suppose that  $f \in K((x))$ . We say that f is an E-algebraic function (over K(x)) if  $tr.deg._{K(x)}K(x,\Omega(f)) < \infty$ .

**Notation.** We shall denote by  $\Gamma_K$ , the set of all E-algebraic functions.

Example 7.2. Let  $K = F_2$  and  $\alpha$  be a 2-adic integer. Let  $f_{\alpha} = (1+x)^{\alpha} \in F_2[[x]]$ . If  $\alpha$  is rational, then  $f_{\alpha}$  is algebraic ([20]) and hence  $tr.deg._{F_2(x)}F_2(x,\Omega(f_{\alpha})) = 0$ . However, if  $\alpha$  is not rational, then  $f_{\alpha}$  is not algebraic over  $F_2$ . However, we show that  $f_{\alpha} \in \Gamma_K$  and hence the set  $\Gamma_K$  strictly contains the set of all algebraic functions.

Let  $\alpha = \sum_{i=0}^{\infty} \alpha_i 2^i$  be the 2-adic expansion of  $\alpha$ . Then

$$f_{\alpha} = (1+x)^{\alpha} = (1+x)^{\alpha_{\bullet}} \left[ (1+x)^{\frac{\alpha-\alpha_{\bullet}}{2}} \right]^{2}.$$

Hence  $E_o(f_\alpha) = (1+x)^{\frac{\alpha-\alpha_o}{2}}$  and  $E_1(f_\alpha) = \alpha_o E_o(f_\alpha)$ . Therefore,

$$E_o(f_\alpha)^2 = (1+x)^{\alpha-\alpha_o} = \frac{f_\alpha}{(1+x)^{\alpha_o}} \in F_2(x,f_\alpha).$$

Similarly,  $E_1(f_\alpha)^2 \in F_2(x, f_\alpha)$ . One can show that

$$[E_{i_1i_2...i_r}(f_{\alpha})]^{2^r} = \frac{cf_{\alpha}}{(1+x)^{\alpha_o+2\alpha_1+...+2^{r-1}\alpha_{r-1}}} \in F_2(x,f_{\alpha}),$$

where  $c \in F_2$  and  $E_{ij} = E_i o E_j$ . Therefore,

$$tr.deg._{F_2(x)}F_2(x,\Omega(f_\alpha))=1,$$

since  $f_{\alpha}$  is not algebraic. Thus  $f_{\alpha} \in \Gamma_K$ .

Note that if  $f \in K((x))$  is algebraic, then

$$tr.deg._{K(x)}K(x,\Omega(f))=0.$$

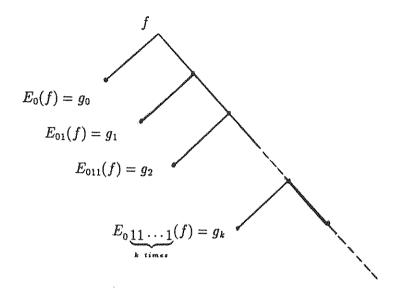
So as in the case of characteristic zero, every algebraic function is E-algebraic.

Now we have that  $\Gamma_K$  is a field.

Theorem 7.3.  $\Gamma_K$  with ordinary addition and multiplication of series is a field.

Let  $\overline{K(x)}$  be the algebraic closure of K(x) in K((x)). In Example 7.2 we showed that  $\overline{K(x)} \subset \Gamma_K$ . Now we show that  $\Gamma_K \subset K((x))$ . That is, we construct a power series  $f = \sum_{n=0}^{\infty} a_n x^n \in K[[x]]$  such that  $f \notin \Gamma_K$ .

Example 7.4. Let  $K=F_2$  and  $\theta$  be a 2-adic integer which is transcendental over Q. Then  $1, \theta, \theta^2, \cdots$  are linearly independent over Q. Let  $f_{\theta} = (1+x)^{\theta} \in F_2[[x]]$ . Then one can show that  $f_{\theta}, f_{\theta^2}, \cdots$  are algebraically independent over  $F_2(x)$  (see, for example, [17]). Let  $g_k = f_{\theta^k} = (1+x)^{\theta^k}$ , for k = 0, 1, 2, ... and define f as follows.



If we specify  $a_n$ , then, since  $F_2(g_o, g_1, ...)$  has infinite transcendence degree over  $F_2(x)$  and since  $F_2(g_o, g_1, ...) \subseteq F_2(x, \Omega(f))$ , we conclude that  $f \notin \Gamma_K$ .

In order to specify  $a_n$ , we write the binary expansion of n,  $n = \sum_{j=1}^{i} n_j 2^{j-1}$ , say

$$n = n_1 n_2 ... n_k n_{k+1} n_{k+2} ... n_i = 11 ... 10 n_{k+2} ... n_i$$

(where  $n_i = 1$ ) and if we use the definition of the  $E_i$  operator, then we can get

$$a_n = E_{n_i n_{i-1} \dots n_{k+2}}(g_k)(0) = \begin{pmatrix} \theta^k \\ m \end{pmatrix} \pmod{2}$$

where  $E_{ij} = E_i o E_j$  and  $m = \sum_{j=k+2}^i n_j 2^{j-k-2}$  as required. Therefore,

$$\overline{K(x)} \subset \Gamma_K \subset K((x)).$$

As we see from the above example  $tr.deg._{K(x)}K((x)) = \infty$ .

Barcanescu and Brezuleanu in [2] proved the following proposition.

**Proposition** 7.5. Suppose that K is a field of characteristic p > 0. Suppose that  $K \subseteq M$  is a separable field extension. Let  $u \in M \setminus K$ . Then the extension  $K(u) \subseteq M$  is a separable extension if and only if  $u \notin M^p K$ .

We use the above result and have that  $K(x) \subseteq \Gamma_K$  is a separable extension.

**Proposition** 7.6. i)  $\Gamma_K$  is closed under the E and D (derivative) operators.

- ii)  $K(x) \subseteq \Gamma_K$  is a separable extension.
- iii)  $tr.deg._{K(x)}\Gamma_K = \infty$ .
- iv)  $\Gamma_K$  is algebraically closed in K((x)).

Proof. See[20].

**Theorem** 7.7.  $\Gamma_K$  is not closed under the Hadamard product operation.

Proof. See [20].

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Proceedings of the Monthly Colloquium of the Iranian Mathematical Society Vol. 2, pp 201-225 (2004)

### Riemann Extension and Complete Lifts

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Abstract: After some fundamental definitions we define Homogeneous and symmetric spaces, both for Affine and Riemannian cases. Then we introduce  $s^k$ -manifolds where, k=2 gives the Cartan symmetric spaces. A generalization of  $s^k$ -manifolds are given by Ledger and Obata as s-manifolds (1968). These spaces are defined locally.

Ledger and Graham gave a general and globally defined sspaces called s-regular spaces. Further works has been done by Ledger, Gray, Sasaki, Vanake, Kowalski, Sekizawa, Razavi and myself, that is  $\Sigma$ -spaces are defined and lifted to the tangent spaces.

This lecture is restricted to the enlargement of homogeneous property to the tangent and cotangent bundles.

MSC (2000): 53022

Keywords: Homogeneous, Symmetric, Riemannian, Affine, sk-spaces, Σ-spaces

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#### 1. Introduction

A topological space M is locally Euclidean if, for every  $x \in M$  there exist an integer  $n \geq 0$ , an open set  $U \subseteq M$  of n, an open subset  $U' \subseteq \mathbb{R}^n$  and a homeomorphism  $\varphi: U \longrightarrow U'$ .

The integer n is uniquely determined by x and is called local dimension of M at x.

If M is Hausdorff and second countable, then by Brouwer theorem on invariance of domain (if  $U \subseteq \mathbb{R}^n$  and  $V \subseteq \mathbb{R}^n$  are open subsets such that U is homeomorphic to V, then m = n) then the local dimension is n at every point  $x \in M$ , and is called the dimension of M.

**Definition 1.1** A topological space M is a topological manifold of dimension n if

- 1. M is locally Euclidean of local dimension n.
- 2. M is Hausdorff.
- 3. M is second countable.

The locally Euclidean property allows us to choose local coordinates in any small region of M.

Definition 1.2 A coordinate chart on M is a pair  $(U,\varphi)$  where  $U\subseteq M$  is an open subset and  $\varphi:U\longrightarrow \mathbb{R}^n$  is a homeomorphism onto an open subset of  $\mathbb{R}^n$ . Let  $\varphi(p)=(x^1(p),\cdots,x^n(p))$  then the n-tuple  $(x^i(p))$  is taken as the coordinate of  $p\in M$  and  $x^i,s$  as coordinate functions. Relative to such a coordinatization, we can do calculus in the region U of M. The problem is that the point p will generally belong to infinitely many different coordinate charts and calculus in one of these coordinatizations about p might not agree with calculus in another. We

need the coordinate systems to be smoothly compatible in the following sense.

**Definition 1.3** Two coordinate charts  $(U, \varphi)$  and  $(V, \varphi)$  on M are said to be  $c^{\infty}$ -related if either  $U \cap V = \emptyset$  or  $\varphi \circ \psi^{-1} : \psi(U \cap V) \longrightarrow \varphi(U \cap V)$  is a diffeomorphism, between open sets of  $\mathbb{R}^n$ .  $\varphi \circ \psi^{-1}$  is called smooth changes of coordinates on  $U \cap V$ .

A  $c^{\infty}$  atlas on M is a collection  $\mathcal{A}=\{(U_{\alpha},\varphi_{\alpha})\}_{\alpha\in I}$  of coordinate charts such that

- (1)  $(qU_{\alpha}, \varphi_{\alpha})$  is  $c^{\infty}$ -related to  $(U_{\beta}, \varphi_{\beta})$   $\forall \alpha, \beta \in I$ .
- $(2) M = \bigcup_{\alpha \in I} U_{\alpha}$

Two  $c^{\infty}$  atlases  $\mathcal{A}$  and  $\mathcal{A}'$  are equivalent if  $\mathcal{A} \cup \mathcal{A}'$  is also a  $c^{\infty}$  atlas on M. Equivalence of  $c^{\infty}$  atlases is an equivalent relation. Each  $c^{\infty}$  atlas on M is equivalent to a unique maximal  $c^{\infty}$  atlas on M.

**Definition 1.4** A maximal  $c^{\infty}$  atlas  $\mathcal{A}$  on M is called smooth structure or differentiable structure or  $c^{\infty}$  structure on M. The pair  $(M, \mathcal{A})$  is called a smooth or differentiable or  $c^{\infty}$  manifold of dimension n or simply n-manifold.

By substituting  $c^k$  for  $c^{\infty}$  we obtain  $c^k$  manifold  $1 \leq k < \infty$ . The same is for real analytic or  $c^{\omega}$  manifold. In all these cases  $\mathbb{R}^n$  is called the model space. We can take  $\mathbb{C}^n$  as model space (complex manifold) or a Hilbert or Banach spaces, where the manifold will be infinite dimensional.

**Definition 1.5** A function  $f: M \longrightarrow R$  is said to be smooth if for each  $x \in M$ , there is a chart  $(U, \varphi) \in A$  such that  $x \in U$  and  $f \circ \varphi^{-1} : \varphi(U) \longrightarrow I \!\! R^n$  is smooth. The set of all smooth, real valued functions on M will be denoted by  $\mathcal{D}(M)$ .

**Lemma 1.6** The function  $f: M \longrightarrow \mathbb{R}^n$  is smooth iff

 $f \circ \varphi_{\alpha}^{-1} : \varphi_{\alpha}(U_{\alpha}) \longrightarrow \mathbb{R}^{n}$  is smooth for every  $(U_{\alpha}, U_{\alpha}) \in \mathcal{A}$ .

**Definition 1.7** Let M and N be  $c^{\infty}$  manifold with respective smooth structures  $\mathcal{A}$  and  $\mathcal{B}$ . A mapping  $f: M \longrightarrow N$  is said to be smooth if, for each  $x \in U_{\alpha}$ ,  $f(U_{\alpha}) \subseteq V_{\beta}$  and

$$\psi_{\beta} ofo \varphi_{\alpha}^{-1} : (\varphi_{\alpha}(U_{\alpha})) \longrightarrow \psi_{\beta}(V_{\beta})$$

is smooth.

Lemma 1.8 the map  $f: M \longrightarrow N$  is smooth if and only if, for all choices of  $(U_{\alpha}, \psi_{\alpha}) \in \mathcal{A}$  and  $(U_{\alpha}, \psi_{\alpha}) \in \mathcal{B}$  such that  $f(U_{\alpha}) \subseteq V_{\beta}$ , then the map

$$\psi_{\beta} ofo \varphi_{\alpha}^{-1} : (\varphi_{\alpha}(U_{\alpha})) \longrightarrow \psi_{\beta}(V_{\beta})$$

is smooth. [1;70]

**Definition 1.9** Let M be a differentiable manifold and  $p \in M$ , for a chart or local coordinate neighborhood  $(U, \varphi)$  of p, let

 $\alpha(-\varepsilon,\varepsilon)\subset I\!\!R\longrightarrow U\subset M$  be a differentiable curve such that  $\alpha(0)=p$ . Let C(U,p) be the set of all differentiable curves on M passing through  $p\in U$ . If  $\alpha,\beta\in C(U,p)$ , then  $\alpha$  and  $\beta$  are said to be infinitesimally equivalent at p or  $\alpha\stackrel{p}{\succeq}\beta$ , if and only if

$$\frac{d}{dt}(f(\alpha(t)))_{t=0} = \frac{d}{dt}(f(\beta(t)))_{t=0}$$

It is easy to check that  $\stackrel{p}{\simeq}$  is an equivalence relation. An equivalence class of  $\alpha \in C(U,p)$  is denoted by  $[\alpha]_p$ , and is called a tangent vector to U (or M) at p, and the set

$$T_pU = \{ [\alpha]_p; \alpha \in C(U,p) \}$$

is called the tangent space at p.

**Lemma 1.10** For each  $[\alpha]_p \in T_pU$  the operator

$$D[\alpha]_p:\mathcal{D}(U)\longrightarrow IR$$

is well defined, by choosing any representative  $\alpha \in [\alpha]_p$  and setting

$$D[\alpha]_p(f) = \frac{d}{dt}(f(\alpha(t)))_{t=0} \quad \forall f \in \mathcal{D}(U),$$

conversely  $[\alpha]_p$  is uniquely determined by the operator  $D[\alpha]_p$ , [1;28].

By the definition of infinitesimal curves there is a natural one to one correspondence between  $T_pU$  (or  $T_pM$ ) and  $I\!\!R^n$  introducing tangent vectors as a linear approximation of nonlinear curves. The key lemma for this, follows:

**Lemma 1.11** Let  $[\alpha]_p, [\beta]_p \in T_pU$  and  $a, b \in \mathbb{R}$  then there is a unique infinitesimal curve  $[\gamma]_p$  such that the associated derivatives on  $\mathcal{D}U$  satisfy

$$D[\gamma]_p = aD[\alpha]_p + bD[\beta]_p$$

**Proof:** Let  $\gamma:(-\varepsilon,\varepsilon)\longrightarrow U$  be a curve on M defined by

$$\gamma(t) = a\alpha(t) = b\beta(t) - (a+b-1)p$$

Since  $\alpha(0) = \beta(0) = p$  then  $\gamma(0) = p$ . Also

$$D[\gamma]_p(f) = \frac{d}{dt}(f(\gamma(t)))_{t=0}$$

$$= \frac{d}{dt}f(a\alpha(t) + b\beta(t) - (a+b-1)p)_{t=0}$$

$$= a\frac{d}{dt}(f(\alpha(t)))_{t=0} + b\frac{d}{dt}(f(\beta(t)))_{t=0}$$

$$= D[\alpha]_p(f) + D[\beta]_p(f)$$

$$D[\gamma]_p = aD[\alpha]_p + bD[\beta]_p$$

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Therefore the operator  $D[\alpha]_p$  makes  $T_pU$  or  $T_pM$  into an n-dimensional vector space over R. The zero vector is the infinitesimal curve represented by the constant curve  $\alpha: (\varepsilon, \varepsilon) \longrightarrow U$  by  $\alpha(t) = p$ . If  $[\alpha]_p \in T_pU$ . then  $-[\alpha]_p = [-\alpha]_p$  where  $(-\alpha)(t) = \alpha(-t)$ , defined for all sufficiently small value of t.

The operator  $D[\alpha]_p$  depends only the behavior of f in an arbitrary small neighborhood of p, that is  $D[\alpha]_p(f)$  depends only on the "germ" of the t at p.

**Definition 1.12** We say that  $f,g \in \mathcal{D}(u)$  are germinally equivalent at p and write  $f \stackrel{p}{\sim} g$ . if there is an open neighborhood W of p in U such that f/W = g/W. Germinally equivalent is an equivalence relation on  $\mathcal{D}(u)$ , and equivalence class  $[f]_p$  of  $f \in \mathcal{D}(u)$  is called the germ of f at p. The set  $\mathcal{D}(u)/\tilde{p}$  of germs at p is denoted by  $G_p$ .

**Definition 1.13** For each  $\alpha \in C(U,p)$  the operator  $D[\alpha]_p: G_p \longrightarrow \mathbb{R}$  is defined by  $D[\alpha]_p([f]_p) = \frac{d}{dt}(f(\alpha,t))_{t=0}$ .

**Definition 1.14**  $G_p$  together with the operations

- 1.  $a[f]_p = [af]_p$   $a \in \mathbb{R}$ ,  $[f]_p \in G_p$  (scalar multiplication)
- 2.  $[f]_p + [g]_p = [\frac{f}{w} + \frac{g}{w}]_p$   $[f]_p, [g]_p \in G_p(addition), W$  is open neighborhood of p in  $dom(f) \cap dom(g)$ .
- 3.  $[f]_p[g]_p = [(\frac{f}{w})(\frac{g}{w})]_p$   $[f]_p, [g]_p \in G_p$  (multiplication)

**Definition 1.15** The evaluation map  $e_p: G_p \longrightarrow \mathbb{R}$  is defined by  $e_p([f]_p) = f(p)$ .

A derivative operator (derivative) on  $G_p$  is an IR-linear map  $D:G_p\longrightarrow IR$  such that

$$D([f][g]) = D[f]e_p([g]) + c_p([p])D[g]$$
  
=  $g(p)D[f] + f(p)D[g].$ 

A derivative on  $G_p$  is also called a tangent vector to U at P. The set of all derivatives on  $G_p$  is denoted by  $T_p(U)$  or  $T_p(M)$  and is called the tangent space to M at p.

**Lemma 1.16** The tangent space  $T_pU$  or  $T_pM$  is a vector space over  $\mathbb{R}$  under the operations:

1. 
$$(aD)[f] = a(D[f])$$
  $a \in \mathbb{R}D \in T_pM, [f] \in G_p$ 

2. 
$$(D_1 + D_2)[f] = D_1[f] + D_2[f]$$
  $D_1, D_2 \in T_pM, [f] \in G_p$ 

The two definitions of  $T_PU$  give canonically isomorphic spaces [1;30] For example define  $(D_i)_pG_p \longrightarrow I\!\!R$  by

$$(D_i)_p[f]_p = (\frac{\partial t}{\partial x^i})(p)$$

It is proved that the set of tangent vectors  $(D_i)_p = (\frac{\partial t}{\partial x^i})(p)$  is a basis of the vector space  $T_pU$  or  $T_pM$  [1;32].

that is if V is a tangent vector in  $T_pU$ , then

$$V = \sum_{i=1}^{n} v^{i}(D_{i})_{p} = \sum_{i=1}^{n} v^{i}(\frac{\partial}{\partial x^{i}})_{p}$$

or simply by using the summation convention  $V=v^i(\frac{\partial}{\partial x^i})_p$ . Let  $TU=\bigcup_{x\in U}T_p(U)$  be the disjoint union, then there is a one to one correspondence  $TU\longleftrightarrow U\times I\!\!R$  given by

$$v^{i}(\frac{\partial}{\partial x^{i}})_{p} \longleftrightarrow (p; v^{1}, v^{2}, \dots, v^{n})$$
 (\*)

We use (\*) to transfer the topology of  $U \times \mathbb{R}$  to TU.

**Definition 1.17** Let  $f: M \longrightarrow N$  be a smooth map between differentiable manifolds M and n, where  $\dim M = m, \dim N = n$ . For any  $p \in U \subset M$  and  $f(p) \in V \subset N$  we define  $(df)_p = (f)_p : T_pU \longrightarrow T_{f(p)}V$  by

$$(df)_p[\alpha]_p = [fo\alpha]_{f(p)} \quad \forall [\alpha]_p \in T_pU.$$

This is called the differential of f at p.

**Lemma 1.18** The differential  $(df)_p$  is well defined linear map. Relative to the basis  $(\frac{\partial}{\partial x^i})_p$  of  $T_pU$  and  $(\frac{\partial}{\partial y^j})_{f(p)}$  of  $T_{f(p)}V$  the matrix of  $(df)_p$  is the Jacobian matrix of f at p. That is  $Jf = (\frac{\partial y^j}{\partial x^i})_p$   $1 \le i \le n$   $1 \le j \le m$ 

The differential  $(f_{\alpha})_p$  computed at all  $p \in U$ , assemble to a mapping

$$f_* = df : TU \longrightarrow TV$$

given by

$$f_{lpha}\left(p,\left[egin{array}{c} v' \ dots \ v^m \end{array}
ight]
ight)=\left(f(p),J_f\left[egin{array}{c} v' \ dots \ v^m \end{array}
ight]
ight).$$

Here we have identified TU with  $U \times \mathbb{R}^m$  and TV with  $V \times \mathbb{R}^n$  [1;35]

# 2. Tangent and cotangent bundles

Let M be a  $c^{\infty}$ , m-manifold with structure  $\{(U_{\alpha}, \varphi_{\alpha})\}_{\alpha \in I}$ . Consider the set  $TM = \bigcup_{p \in M} T_p M$  as disjoint union, for each  $U_{\alpha}$   $\alpha \in I$ , define

$$T(U_{\alpha}) = \bigcup_{p \in U_{\alpha}} T_p(U) \subseteq TM$$

. Then the individual linear map  $(d\varphi_{\alpha})_p$   $p\in U_{\alpha}$  unite to define a set map, considering (\*)

$$d\varphi_\alpha: T(U_\alpha) \longrightarrow T(\varphi_\alpha(U_\alpha)) = \varphi_\alpha(U_\alpha) \times I\!\!R^n \subseteq I\!\!R^{2n}$$

More precisely, if  $V_p$  denotes a tangent vector to M at  $p \in U_{\alpha}$ .

$$(d\varphi_{\alpha})(V_p) = (\varphi_{\alpha})(p), (d\varphi_{\alpha})_p V_p$$

and this defines a bijection of  $TU_{\alpha}$  onto an open subset of  $\mathbb{R}^{2n}$ . If  $U_{\alpha} \cap U_{\beta} \neq \phi$  consider

$$d\varphi_{\alpha}.d\varphi_{\beta}^{-1}:T(\varphi_{\beta}(U_{\alpha}\cap U_{\beta}))\longrightarrow T(\varphi_{\alpha}(U_{\alpha}\cap U_{\beta}))$$

is a  $c^{\infty}$  diffeomorphism between open sets of  $\mathbb{R}^{2n}$ . We Topologies the set TM such that if

$$W \subseteq d\varphi_{\alpha}(T(U_{\alpha})) = T(\varphi_{\alpha}(U_{\alpha})) \subseteq IR^{2n}$$

is an open set, then  $d\varphi_{\alpha}^{-1}(w)$  is to be an open subset of TM.

Lemma 2.1 The above sets from the base of a topology on TM and, in this topology, TM is a topological manifold of dimension 2m. Furthermore, the system  $\{T(U_{\alpha}), d\varphi_{\alpha}lpha\}_{\alpha\in I}$  is a  $c^{\infty}$  atlas on TM, determining a maximal atlas A.

 $(TM, \mathcal{A})$  is called the tangent bundle of M. The map  $\phi: TM \longrightarrow M$  determined by  $\pi(p, v) = p$   $p \in M, v \in T_pM$  is smooth. TM is locally a Cartesian product of  $U_{\alpha} \subset M$  and  $\mathbb{R}^m$  and  $\pi^{-1}(p) = T_pM$  is the tangent space of M at p.

**Definition 2.2** A vector field on M is a smooth map

 $X: M \longrightarrow M$ ;  $P \longmapsto X_p$  such that  $\pi o X = id_M$ , the set of all vector fields on M is denoted by  $D^1(M)$ . If  $(U, x^1, x^2, \dots, x^n)$  is a coordinate chart on M, then  $X = X^i \frac{\partial}{\partial x^i}$  where  $X^i: U \longrightarrow IR$  are smooth function, on M for  $1 \le i \le n$ .

In general Let  $(U, x^1, x^2, \ldots, x^n)$  be a coordinate chart on M, denote by  $(X^i)$  the system of Cartesian coordinates on each tangent space  $T_p(M)$ . Then in any open subset  $\pi^{-1}(U)$  of TM we introduce local coordinates  $(x^i, X^i)$   $1 \le i \le n$ , which are called induced coordinate on  $\pi^{-1}(u)$ . We denote

$$\partial_i = \frac{\partial}{\partial x^i}$$
 and  $\bar{\partial}_i = \frac{\partial}{\partial X^i}$ 

# 3. Covectors and 1-forms

**Definition 3.1** Let M be a differentiable manifold  $p \in M$ . The dual space  $T_p^*M = Hom(T_pM, I\!\!R)$  is called the cotangent space of M at p. Each element  $w \in T_p^*M$  is called a cotangent vector to M at p.

A typical cotangent vector is the differential map. Let  $U\subseteq M$  be open,  $p\in U$ , and let  $f\in c^\infty(U)$ . Since  $T(I\!\!R)=I\!\!R$ , we obtain a linear functional  $(df)_p:T_pM\longrightarrow T(R)=R$ , so  $(df)_p\in T_p^*M$ . It is evident that  $(df_p)$  depends only on the germ  $[f]_p\in G_p$ , so we obtain a surjective  $I\!\!R$ -linear map  $d:G_p\longrightarrow T_p^*M$ 

For each  $X_p \in T_pM$ ;  $(df)_pX_p = X_p(f)$ .

**Definition 3.2** If  $\varphi: M \longrightarrow N$  is a smooth map between manifolds, if  $p \in M$ , and if  $w \in T_p^*M$ , then  $\varphi_p^*(w) \in T_p^*(M)$  is defined by

$$\varphi_p^*(w)X_p = w(\varphi_{*p}(X_p)).$$

The linear map  $\varphi_p^*$  is called the adjoint of  $\varphi_{*p}$ .

**Lemma 3.3** Relative to local coordinates  $x^1, \ldots, x^n$  about  $p \in M$ , The differentials  $dx^1, \ldots, dx^n$  at p from a basis of  $T_p^*M$  and

$$(dx^i)_p(rac{\partial}{\partial x^j})_p=rac{\partial x^i}{\partial x^j}(p)=\delta_{ij}$$

hence  $\dim T_p^* M = m$  [1;160]

Let  $\varphi: M \longrightarrow N$  be smooth map between smooth manifolds. If  $x^1, \ldots, x^m$  are coordinates about  $p \in M$  and  $y^1, \ldots, y^n$  coordinates about  $\varphi(p) \in N$ , we have basis  $(\frac{\partial}{\partial x^i})_p \qquad 1 \leq i \leq m$  for  $T_pM$  and  $(\frac{\partial}{\partial y^j})_{\varphi(p)}$  for  $T_{\varphi(p)}N$  and the dual basis  $(dx^i)_p$  for  $T_p^*M$  and  $(dy^j)_{\varphi(p)}$  for  $T_\varphi^*N$ . The relations are as follows.

$$(\frac{\partial}{\partial y^{i}})_{\varphi(p)} = (\frac{\partial x^{i}}{\partial y^{i}})_{p}(\frac{\partial}{\partial x^{i}})_{p} \quad 1 \leq i \leq m \quad 1 \leq j \leq n$$

Riemann extension and complete lifts

$$(dy^i)_{\varphi(p)} = (\frac{\partial y^j}{\partial x^i})_p (dx^i)_p \quad 1 \le i \le m \quad 1 \le j \le n$$

Now let  $A = \{(U_{\alpha}, \varphi_{\alpha})\}_{\alpha \in I}$  be a maximal smooth atlas on M. As a set let

$$T^*M = \bigcup_{p \in M} T_p^*M$$

. We topologies each  $T^*(U_{\alpha})$  via the bijection

$$\psi_{\alpha}: T^{\alpha}(U_{\alpha}) \longrightarrow U_{\alpha} \times I \!\! R^n$$

defined by

$$\psi_{lpha}(w_i(dx^i)_p) = (p, \left[egin{array}{c} w_1 \ dots \ w_m \end{array}
ight])$$

repeating the case for tangent bundle,  $T^*M$  becomes a 2m dimensional smooth manifold, and for any  $w \in T_p^*M$  we have  $w = w_i dx^i$  for every tangent vector  $X_p = (X^i(\frac{\partial}{\partial x^i})_p \in T_pM$  and 1-form  $w = w_i (dx^i)_p \in T_p^*M$  we have

$$\begin{split} w_p(X_p) &= (w_i(dx^i)_p)(X^j(\frac{\partial}{\partial x^j})_p) = w_i X^j(dx^i(\frac{\partial}{\partial x^j}))_p \\ &= w_i X^j \partial_{ij} = w_i X^i. \qquad 1 \le i, j \le n \end{split}$$

Lie bracket is defined as

$$[ , ]: D^{1}(M) \times D^{1}(M) \longrightarrow D^{1}(M)$$
$$[ , ](X,Y) = XY - YX$$

and  $(D^1(M), [ , ])$  is called the Lie algebra of vectorfields on M, or simply Lie algebra of M. In general let  $(U; x^1, \ldots, x^n)$  be a coordinate chart on M, then every 1-form w at  $p \in U$  is written as  $w = \xi_i dx^i$ . Then  $(\xi_i)$  are the system of Cartesian coordinate in any open subset,  $\pi_1^{-1}/U$ 

on  $T^*M$ , we can introduce local coordinate  $(x^i, \xi_i)$   $1 \le i \le n$ , which are called induced coordinates. We denote

$$\partial_i = \frac{\partial}{\partial x^i}$$
 ,  $\partial^i = \frac{\partial}{\partial \xi_i}$ 

# 4. Lie groups and Lie algebra

A Lie group G is a smooth manifold without boundary which is also a group such that the group operation  $\varphi: G \times G \longrightarrow G$ ,  $\varphi(x,y) = xy$  and the inversion  $\psi: G \longrightarrow G$ ;  $\psi(x) = x^{-1}$  are boat smooth.

For example every finite dimensional vector space over  $I\!\!R$  or  $I\!\!C$  is a Lie group under vector addition.

$$(M(n,\mathcal{C})+), (Gl(n,\mathcal{C}), \times)(M(n,R), +),$$
  
 $(Gl(n,R), \times).U(n), O(n), Sl(n,R), \dots, S^7, S^3, S^1$ 

are all Lie groups.

**Definition 4.1** Let G be a Lie group,  $g \in G$ . Let translation by g is a smooth map  $L_g: G \longrightarrow G$  defined by  $L_a(x) = ax \quad \forall x \in G$ . A vectorfield  $X \in D^1(G)$  is left invariant if, for each  $g \in G$   $(dl_g)X = X$ . The set of all left invariant vectorfields on G is denoted by L(G).

**Proposition 4.2** The subset  $L(G) \subset D^1(G)$  is the Lie algebra of the Lie group G and the evaluation map  $E: L(G) \longrightarrow T_eG$  by  $E(X) = X_e$  is an isomorphism of vector spaces (e is the unit of G). Hence, dimL(G) = dimG.[1,130]

Let A be  $m \times m$  matrix and  $\alpha(t) = \exp A$   $t \in R$  such that  $\alpha(0) = I$ . Then  $\alpha(t+s) = \alpha(t)\alpha(s)$   $t,s \in R$  and  $\alpha^{-1}(t) = \exp(-tA)$  that is  $\alpha(t)$  is a 1-parameter subgroup of Gl(n,R) of G is a Lie group then the exponential map  $\exp: L(G) \equiv T_e G \longrightarrow G$  is defined by  $\exp(X) = \alpha_X(1)$  where  $\alpha_X(t) = \exp tX$ .

# 5. Homogeneous space

**Definition 5.1** Let M be a smooth manifold and G a Lie group. A smooth map  $\varphi: G \times M \longrightarrow M$  written  $\varphi(g,x) = gx$ , is said to be an action of G on M, and G is called a Lie transformation group on M if

1. 
$$\varphi(g_2, \varphi(g_1, x)) = \varphi(g_1g_2, x)$$
 or  $g_1(g_2x) = (g_1g_2)x$   
 $\forall g_1, g_2 \in G \text{ and } \forall x \in M$ 

2. 
$$\varphi(e,x) = x \ e \in G \forall x \in M$$

**Definition 5.2** An orbit of the action  $G \times M \longrightarrow M$  is a set of points of the form  $\{gx_0; g \in G\}$  where  $x_0 \in M$ . The action is transitive if M itself is an orbit, in which case M is said to be homogeneous space of G.

The orbits of a group action are equivalence classes, two points  $x, y \in M$  being equivalent under the action if  $\exists g \in G$  such that gx = y.

**Definition 5.3** Let M be a homogeneous space of G and let  $x_0 \in M$ . The isotropy of  $x_0$  is the set

$$G_{x_0} = \{g \in G; gx_0 = x_0\}$$

which is a property Lie subgroup of G.

**Proposition 5.4** There is a smooth structure on quotient space where  $G/G_{x_0}$  and  $M = G/G_{x_0}$ , is a homogeneous space where  $G \times M \longrightarrow M$  is given by  $g(hG_{x_0}) = (gh)G_{x_0}$  [1,144]

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**Corollary 5.5** If G is a Lie group and  $H \subseteq G$  is a closed normal subgroup, then the group G/H has a smooth structure in which it is a Lie group.[1,148]

For example 
$$S^{n-1} = o(n)/o(n-1)$$
 and  $S^{2n-1} = U(n)/U(n-1)$ .

**Definition 5.6** Let  $\varphi: \mathbb{R} \times n \longrightarrow M$  be a smooth map such that  $\varphi(o,p) = p$  and  $\varphi(s,\varphi(t,p)) = \varphi(s+t,p)$ , That is the additive group  $\mathbb{R}$  is a transformation group on M. For each  $t \in \mathbb{R}$  we define  $\varphi_t M \longrightarrow M$  such that  $\varphi_t(p) = \varphi(t,p)$  then it is easy to show that  $\varphi_0 = Id$  and  $\varphi_{s+t} = \varphi_s o \varphi_t$  also  $\varphi_t^{-1} = \varphi_{-t}$ .

 $\{\varphi_t; t \in I\!\!R\}$  is a group. Called 1-parameter group of transformations on M.

#### 6. Connections

Let  $U \subseteq \mathbb{R}^n$  be open. Given  $X, Y \in D^1(U)$ , define  $D_X(Y) \in D^1(M)$  as follows.

Write  $X = X^i \frac{\partial}{\partial x^j}$   $Y = Y^j \frac{\partial}{\partial x^j}$  and define

$$D_X(Y) = X(Y^j) \frac{\partial}{\partial x^j} = X^i \frac{\partial x^j}{\partial x^i} \frac{\partial}{\partial x^i}$$

We can view D as an R-bilinear map

$$D: D^1(U) \times D^1(U) \longrightarrow D^1(U)$$

It has the following properties:

1. 
$$D_{fX}(Y) = fD_X(Y)$$
.  $\forall f \in D(U) \forall X, Y \in D^1(U)$ 

2. 
$$D_X(fY) = X(f)Y + fD_X(Y)$$

 $D_X(Y)$  is called derivative of the vectorfield Y in X direction. The operation D is called Euclidean connection.

**Definition 6.1** Let M be a smooth manifold. An Affine connection or M is an R-bilinear map

$$\nabla: D^1(M) \times D^1(M) \longrightarrow D^1(M)$$

written as  $\nabla(X,Y) = \nabla_X(Y)$  with the following properties, for each  $X \in D^1(M); \nabla_X : D^1(M) \longrightarrow D^1(M)$  is a linear mapping, that is

1. 
$$\nabla_{fX+gY}(Z) = f(\nabla_X Z) + g(\nabla_Y Z)$$
  $f, g \in c^{\infty}(M), Y, Z \in D^1(M)$ 

2. 
$$\nabla_X(fY) = X(f)Y + f(\nabla_X Y)$$

The operator  $\nabla_X$  is called covariant differentiation with respect to X.

If  $\nabla$  is an offine connection on M and  $(U_1x^1,\ldots,x^n)$  is a coordinate chart, set  $X=\frac{\partial}{\partial x^i}$  then

$$\nabla_{\frac{\partial}{\partial x^i}}(\frac{\partial}{\partial x^j}) = \Gamma_{ij}^k(\frac{\partial}{\partial x^i})$$

The smooth functions  $\Gamma_{ij}^k$  are called the Christoffel symbols.[3,27]

**Definition 6.2** Let M be a manifold with an offine connection  $\nabla$ . Let  $X \in D^1(M)$  and  $\alpha: (-\epsilon, \epsilon) \longrightarrow M$  be a differentiable curve on M. X is called parallel along  $\alpha$  if  $\nabla_{\alpha(t)} X = 0$ 

**Definition 6.3** Every manifold M has a connection.[1,301]

Let V be a vector space over  $\mathbb{R}$ . A bilinear form on V is defined to be a map  $\varphi: V \times V \longrightarrow \mathbb{R}$  which is linear in each variable.

**Definition 6.4** Suppose  $\nabla$  is an offine connection on M and  $\varphi: M \longrightarrow M$  be smooth and

$$d\varphi(\nabla_X Y) = \nabla_{d\varphi X}(d\varphi Y) \quad \forall X, Y \in D^1(M)$$

Then  $\varphi$  is called offine transformation of  $(M, \nabla)$ .

**Definition 6.5** Let  $(M, \nabla)$  be an offine manifolds, The curve  $\alpha: (-\epsilon, \epsilon) \longrightarrow \alpha(t)$  is called a geodesic if the family of tangent vector  $\alpha(t)$  are parallel along  $\alpha$ . That is  $\nabla^{\alpha(t)} = 0$ .

In a coordinate neighborhood  $(U, x^1, \ldots, x^n)$ ;  $\alpha(t) = (x^i(t))$  is geodesic if it satisfies the system of differential equation

$$\frac{d^2x^k}{dt^2} + \Gamma^k_{ij}\frac{dx^i}{dt}\frac{dx^j}{dt} = 0 \qquad [3,30]$$

 $\varphi$  is called symmetric if  $\varphi(x,y) = \varphi(y,x)$ . a symmetric form is called positive definite if  $\varphi(x,y) \geq 0$  and  $\varphi(x,x) = 0 \iff x = 0$   $\varphi$  is also called an inner product on V or  $\varphi$  is a tensor of type (0,2).

**Definition 6.6** A tensorfield g of smooth bilinear forms on a smooth manifold M which assigns to each  $p \in M$  a symmetric, positive definite, bilinear form  $g_p$  of type (0,2) at  $T_pM$ , that is

$$g_p: T_pM \times T_pM \longrightarrow R$$
 (inner product)

is called a Riemannian metric on M and (M,g) is a Riemannian manifold. If  $g_p$  fails to be positive definite then g is called pseudo-Riemannian and (M,g) is pseudo-Riemannian. [2,155]

For any coordinate neighborhood  $(U, x^1, \ldots, x^n)$ ,

$$g(\frac{\partial}{\partial x^i}, \frac{\partial}{\partial x^j}) = g_{ij}$$

are called the components of g, that is  $g = (g_{ij})$ 

For 
$$M = \mathbb{R}^n$$
  $\left(\frac{\partial}{\partial x^i}, \frac{\partial}{\partial x^j}\right) = \delta_{ij}$ .

**Definition 6.7** A connection  $\nabla$  on the Riemannian manifold (M,g) is a Riemannian connection if for all  $X,Y,Z\in D^1(M)$ .

$$Xg(X,Y) = g(\nabla_X Y, Z) + g(Y, \nabla_X Z).$$

If  $\nabla$  is symmetric then it is called Livi-Civita connection on (M,g).

**Proposition 6.8** A Riemannian manifold M has a unique Livi-Civita connection [1,305]

In any local coordinate chart, the matrix  $(g_{ij})$  of metric coefficients is nonsingular, so we can define  $(g^{kl}) = (g_{ij})^{-1}$ . The coefficient  $g^{kl}$  are rational functions of the metric coefficients  $g_{ij}$  and satisfy  $g_{ik}g^{kl} = \delta_i^l$ .

**Definition 6.9** Let (M,g) and (N,g') be two smooth Riemannian or (pseudo-Riemannian) manifolds and  $\varphi: M \longrightarrow N$  be smooth.  $\varphi$  is called a local isometry if for each  $p \in M$ , there exist open neighborhoods U of p and V of  $\varphi(p)$  such that  $\varphi_p^*(h) = g$ . In other words  $\varphi$  is an tangent space at p and  $\varphi(p)$ . That is  $g_p(X_p, Y_p) = h_{\varphi(p)}(d\varphi X, d\varphi Y)$ 

**Definition 6.10** A vector X on (M,g) is called an isometry or killing vectorfield if local 1-parameter group of local transformations generated by X in a neighborhood of each  $p \in M$ , consists of local isometrics. In the same manner, a vectorfield X on  $(M,\nabla)$  is called an infinitesimal offine transformation of M, if for each  $p \in M$  a local 1-parameter group of local transformation  $\varphi_t$  of a neighborhood U of p into M, preserves the connection  $\nabla$ , more precisely if  $\varphi_t: U \longrightarrow M$  is an offine mapping, where U is provided with offine connection.

An infinitesimal isometry is necessarily an infinitesimal offine transformation.

The Riemannian manifolds (M, g) is geodesically complete if  $\exp_p(X_p)$  is defined for all  $p \in M$  and for all  $X_p \in T_p(M)$ . Equivalently, every local geodesic extends (uniquely) to a geodesic  $\alpha(t)$  for  $-\infty < t < \infty$ .

**Proposition 6.11** (Hopt-Rinow) if M is geodesically complete then  $\exp_p: T_pM \longrightarrow M$  is surjective.[1,318].

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# 7. Riemannian homogeneous space

Let (M, g) be a connected Riemannian manifold. The group of all isometries  $\varphi: M \longrightarrow M$ , will be denoted by I(M). The action of I(M) on M preserves all intrinsic properties.

**proposition 7.1** (Myers, Steenrod). If (M,g) is a Riemannian manifold, the group I(M), with compact -open topology, is isomorphic, as a topological group, to a Lie group such that the natural action  $I(M) \times M \longrightarrow M$  denoted by  $(\varphi, x) \longrightarrow \varphi(x)$  is smooth and M = I(M)/k, where k is a closed subgroup, hence M is homogeneous manifold. [1,345]

**Proposition 7.2** If M is a homogeneous manifold, then it is a complete Riemannian manifold.

**Definition 7.3** A Riemannian symmetric space is a Riemannian manifold (M,g) with the property that, for each  $p \in M$ , is  $\varphi_p \in I(M)$  such that  $\varphi_p(p) = p$  and  $d\varphi_p = -I$ , where I is the identity transformation on  $T_p(M)$ .

Remark that,  $\varphi_p$  reserves every geodesic s(t) through x. That is, if s(0) = x, then  $\varphi_p(s(t)) = s(-t)$ .

**Proposition 7.4** If (M,g) is a symmetric space, then M is complete. If (M,g) is connected symmetric space, then M is s homogeneous Riemannian manifold.

#### 8. Lifts

Let M be a manifold of dimension n. For any local coordinates  $(U, x^1, \ldots, x^n)$  on M, we have  $X = X^i \frac{\partial}{\partial x^i}, w = w_i dx^i$ , and local coordinates  $(\pi^{-1}(U), x^i, X^i) 1 \le i \le n$  for TM,  $\pi_1^{-1}(U)$ ;  $x^i, w_i$ ) for  $T^*M$ . Also

 $\partial_i = \frac{\partial}{\partial x^i}$ ,  $\bar{\partial}_i = \frac{\partial}{\partial X^i}$  basis for tangent spaces of TM and  $\partial_i = \frac{\partial}{\partial x^i}$ ,  $\partial^i = \frac{\partial}{\partial w^i}$  basis for cotangent spaces of  $T^*M$ .

Let  $w \in D_1(M)$  be 1-form on M, we may regard w as a function on TM and denoted by  $\tau w$ .

For a function  $\varphi \in c^{\infty}(M)$  we denote  $\varphi c = \tau(d\varphi) \in TM$  and called complete lift of  $\varphi$  to TM. For any vectorfield X on M we define the complete lift  $X^c \in D^1(TM)$  by  $X^c \varphi^c = (X\varphi)^c \quad \forall \varphi \in c^{\infty}(M)$ .

For any 1-form  $w \in D_1(M)$  we define 1-form  $w^c \in D_1(TM)$  by  $w^c(X^c) = (w(X))^c \quad \forall X \in D^1(M)$ , and  $w^c$  is called the complete lift of w to TM.

For  $X=X^i\partial_i\in D^1(M), \ X^\nu=X^i\frac{\partial}{\partial X^i}=\xi^i\bar\partial_i$  is called the vertical lift of X to TM. Also  $X^h=X^i\partial_i-X^iX^j\Gamma^k_{ij}\bar\partial_k$  is called the horizontal lift of X on TM.

For a metric tensorfield  $g = (g_{ij})$  on M we define

$$\left[\begin{array}{cc} X^k \partial_k g_{ij} & g_{ij} \\ g_{ij} & 0 \end{array}\right]$$

as the complete lift of g to TM. In local coordinates if  $g = g_{ij}dx^idx^j$  then  $g^c = \Gamma g_{ij}\delta \xi^i dx^j$ , where  $\delta \xi^i = dx^i + \Gamma^k_{jk}dx^j$ .

For a vectorfield  $X \in D^1(M)$ , the vertical lift  $X^{\nu}$  on  $T^*M$  is a function on  $T^*M$  defined by  $X^{\nu}(\ )=w(X_p)$ .

For lift to  $T^*M$  we have vertical lift a 1-form w to be a vertical vectorfield  $w^{\nu}$  given by

$$w^{\nu}(X^{\nu}) = (w(X))^{\nu} \qquad X \in D^{1}(M)$$

For a function  $X^{\nu} \in c^{\infty}(T^*M)$ ,  $dx^{\nu}$  is a 1-form given by

$$dx^{\nu} = (w_i \partial_i X^i) dx^j + X^i dw_i$$

For a vectorfield  $X \in D^1(M)$  we define a vectorfield  $X^c \in D^1(T^*M)$  by  $X^c = (dx^{\nu})$ 

in, where  $\in$  is a tensorfield of type (2,0) on  $T^*M$  with components

$$\left[\begin{array}{cc} 0 & I \\ I & 0 \end{array}\right]$$

where, I is the Kronrcker tensorfield.

For a vectorfield  $X \in D^1(M)$  the horizontal lift is defined by  $hX(Y^{\nu}) = (\nabla_X Y)^{\nu}$ .

In coordinates:

$$\begin{split} X^{\nu} &= w_i X^i & X = X^i \partial_i \\ w^{\nu} &= w_i \frac{\partial}{\partial w_i} = w_i \partial^i & w = w_i dx^i \\ X^c &= X^i \partial_i - w_i (\partial_j X^i) \partial^j & X = X^i \partial_i \\ hX &= X^i \partial_i + w_i X^j \Gamma^i_{kj} \partial^k & X = X^i \partial_i \end{split}$$

**Proposition 8.1** Let g be a pseudo-Riemannian metric on M. Then  $g^c$  is a pseudo-Riemannian metric on TM. (Yauo, Kobayashi).

**Proposition 8.2** If X is a killing vectorfield of pseudo-Riemannian manifold (M,g), then  $X^{\nu}$  and  $X^{c}$  are killing vectorfields of pseudo-Riemannian manifold  $(TM,g^{c})$ .

Let  $\nabla$  be a offine connection on M, there exists a unique offine connection  $\nabla^c$  on TM, defined by

$$\nabla_{X^c}^c y^c = (\nabla_X Y)^c. \quad \forall X, Y \in D^1(M).$$

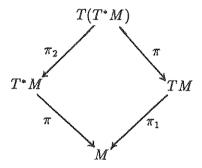
If  $\nabla$  is a Riemannian connection on M, with respect to the pseudo-Riemannian metric g, then  $\nabla^c$  is the Riemannian connection of TM with respect to the pseudo-Riemannian metric  $g^c$ . Also if X is an infinitesimal offine transformation of M, then both  $X^{\nu}$  and  $X^c$  are infinitesimal offine transformation of TM with respect to  $\nabla^c$ .[3]

**Proposition 8.3** If the group of offine transformation of M with respect to  $\nabla^c$  is transitive on TM. Also if M is a pseudo-Riemannian

(offine) symmetric space with metric g (connection  $\nabla$ ), then TM is also a pseudo-Riemannian (offine) symmetric space with metric  $g^c$  (connection  $\nabla^c$ ).

#### 9. Riemann extension

Define a canonical 1-form  $\theta$  on  $T^*M$  via the following diagram



that is, if  $\bar{X}$  is a vectorfield on  $T^*M$ , define  $\theta(\bar{X}) = \pi_2(\bar{X})(\pi_*(\bar{X}))$ . In terms of coordinates.

$$\theta(x^1,\ldots,x^n,w_1,\ldots,w_n)=w_idx^i$$

we call  $\theta$ , the canonical 1-form on  $T^*M$ . Denote  $\Omega = d\theta$ .  $\Omega$  is called a canonical 2-form on  $T^*M$ . In terms of coordinates:

$$\Omega(x^1,\ldots,x^n,w_1,\ldots,w_n)=dw_i\wedge dx^i$$

In matrix form

$$\Omega = \left[ egin{array}{cc} 0 & I \\ -I & 0 \end{array} 
ight].$$

This implies that  $\Omega$  has maximal rank, and consequently defines a bundle isomorphism between  $T(T^*M)$  and  $T^*(T^*M)$ .

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Let H be the horizontal distribution on  $T^*M$  determined by the connection  $\nabla$  on M. Then  $\pi_*$  defines an isomorphism of  $H_{\bar{p}}$  and the tangent space at  $\pi(\bar{p}) \in M$ . That is

$$\pi_*(hX)_{\bar{p}} = X_{\pi(\bar{p})} \qquad \bar{p} \in T^*M.$$

The tangent space  $T_{\bar{p}}(T^*M)$  splits into horizontal and vertical subspaces,

$$T_{\bar{p}}(T^*M) = H_{\bar{p}} + V_{\bar{p}}$$

where  $\stackrel{\cdot}{V}_{\bar{p}}=\{\bar{X}\in T_{\bar{p}}(T^*M);\pi_*\bar{X}=0\}$  and

$$H_{\bar{p}} = \{hX \in T_{\bar{p}}(T^*M); X \in T_{p}M\}.$$

In general every vectorfield  $\bar{X} \in D^1(T^*M)$  can be written as:

$$\begin{split} \bar{X} &= \bar{X}^i \partial_i + \bar{w}_i \partial^i = (\bar{X} \partial_i + \Gamma^l_{ik}) w_c X^{-k} \partial^i) + (\bar{w}_i - \Gamma^l_{ik} w x^{-k}) \partial^i \\ &= \bar{X}^i e_i + \bar{V}_i \partial^i = \bar{X}^h + \bar{X}^v. \end{split}$$

where  $\bar{X}^h \in H$ ,  $\bar{X}^v \in V$ .  $\bar{X}^i$  and  $V_i$  are generally functions of x and w.

We define the Riemann extension metric  $\bar{g}$  on  $T^*M$  by

$$\bar{g}(\bar{X},\bar{Y}) = \Omega(\bar{X}^v,\bar{Y}^h) + \Omega(\bar{Y}^v,\bar{X}^h) \quad \bar{X},\bar{Y} \in T(T^*M)$$

In terms of coordinates  $\bar{X} = \bar{X}^i e_i + \bar{Y}_i \partial^j, \bar{Y} = \bar{Y}^i e_i + \bar{U}_i \partial^i$ 

$$\bar{g}(\bar{X}, \bar{Y}) = (\Gamma^l_{ki} + \Gamma^l_{ik}) w_l \bar{X}^i Y^k + \bar{X}^i \bar{U}_i + \bar{Y}^i \bar{V}_i$$

or

$$g: \left[egin{array}{ccc} -(\Gamma_{ik}^l+\Gamma_{ki}^l)w_l & I \ I & 0 \end{array}
ight]$$

This shows that  $\bar{g}$  depends only on the symmetric part of the connection  $\nabla$  on M. Hence we shall consider M to be an offine manifold with

symmetric connection  $\nabla$ .  $(T^*M, \bar{g})$  is called the Riemann extension of  $(M, \nabla)$ .

We have calculated the Lie algebra of vectorfields on  $(T^*M, \bar{g})$  and the connection  $\bar{\nabla}$  arising from  $\bar{g}$  on  $T^*M$ .

We have proved that the covariant derivative of  $\bar{g}$  is zero that is  $\bar{\nabla}\bar{g}=0$ , which is the case for any Riemannian connection  $\nabla$  on a Riemannian manifold (M,g). i.e.  $\nabla g=o$ .

We have modified the metric  $\bar{g}$  for a pseudo-Riemannian manifold (M,g) so that  $T^*M$  becomes a natural Riemann extension. That is the metric defined on  $T^*M$  on the zero section coinsides with the original metric of M.

Let g be the pseudo-Riemannian metric on M and define

$$\bar{g}_c(\bar{X},\bar{Y}) = \Omega(\bar{X}^v,\bar{Y}^h) + \Omega(\bar{Y}^v,\bar{X}^h) + cg(\pi_*\bar{X},\pi_*\bar{Y})$$

where  $\bar{X}, \bar{Y} \in D^1(T^*M)$  and  $c \in \mathbb{R}$ . For c = 0 it is the metric  $\bar{g}$ , and for c = 1 the additional term is just the horizontal lift of g, and for c > 0 the additional terms is hypothetic to the horizontal lift of g.

We have proved that  $\nabla g = 0 \Longrightarrow \bar{\nabla} \bar{g}_c = 0$ .

**Definition 9.1**  $(T^*M, \bar{g}_c)$  is called a generalized Riemann extension for the pseudo-Riemannian manifold (M, g).

We have calculated the covariant derivatives of all vectorfields on  $T^*M$  and hence get the Riemann extension connection  $\bar{\nabla}$  on  $T^*M$ .

We have calculated the curvature tensorfield  $\bar{R}$  on  $T^*M$  and prove the following theorem:

**Theorem 9.2** Riemann extension  $(T^*M, \bar{g})$  of a manifold  $(M, \nabla)$  is locally symmetric if and only if  $(M, \nabla)$  is locally symmetric.[4]

Now we show that complete lift and Riemann extension of a pseudo-Riemannian manifolds are isomorphic but for offinly connected manifolds they are not necessary isomorphic. M. Toomanian

Let (M,g) be a pseudo-Riemannian manifold. For any local coordinate  $(x^i), 1 \leq i \leq n$  in M we have associated the local coordinates  $(x^i, X^i)$  and  $(x^i, w_i)$  to TM and  $T^*M$  respectively, also  $g^c$  and  $\bar{g}$  the complete lifts and the generalized Riemann extension metrics:

Let  $\varphi: TM \longrightarrow T^*M$  be mapping defined by

$$\varphi(x^i, X^i) = (x^i, g_{ij}X^j)$$
 ,  $g = g_{ij}$ 

It is obviously differentiable and its inverse  $\varphi^{-1} = (x^i, w_i) = (x^i, g^{ij}w_j)$  is also differentiable, hence  $\varphi$  is a diffeomorphism. Furthermore,

$$g^c = A^t \bar{g} A$$

where

$$A = \begin{pmatrix} \frac{\partial(x^i, w_i)}{\partial(x^j, w_j)} \end{pmatrix} = \begin{pmatrix} \delta_{ij} & 0 \\ X^k \partial_j g_{ik} & g_{ij} \end{pmatrix}.$$

then  $\varphi$  is an isometry. This together with the proposition 8.3 gives the following theorem.

**Theorem 9.3** If M is a pseudo-Riemannian symmetric space, then its Riemann extension is a pseudo-Riemannian symmetric space.

Now let  $M = \mathbb{R}^2$  with coordinates  $(x^1, x^2)$  and  $\nabla$  with coordinates  $\Gamma_{11}^2 = x^2$   $\Gamma_{ij}^k = 0$  i, j, k = 1, 2 we have proved that  $T = 0, \nabla R = 0$  and M is an offine locally symmetric space.

We have construct the Riemann extension  $(T^*M, \bar{g})$  and proved that it is locally symmetric pseudo-Riemannian manifold.

We also construct the complete lift  $(TM, \nabla^c)$  and proved that  $\nabla^c$  does not arise from any pseudo-Riemannian metric.

We have proved that following theorem:

Theorem 9.4 The generalized Riemann-extension of a pseudo-Riemannian homogeneous manifold (M,g) is a pseudo-Riemannian homogeneous manifold  $(T^*M, \bar{g}_e)$ .[4]

We have calculated general killing vector fields, curvature tensorfield, general infinitesimal vectorfields and their Lie algebra of  $(T^*M, \bar{g}_c)$ .[5]

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Proceedings of the Monthly Colloquium of the Iranian Mathematical Society Vol. 2, pp 227-270 (2004)

# The Stable Iterative Algorithms with the CADNA Library for Solving Sparse Linear Systems

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Abstract: This paper shows the efficiency of iterative methods for solving large sparse systems of linear algebraic equations when they are performed with an efficient round-off error analysis method. In a first part the most robust iterative methods such as GMRES, hybrid GMRES, A<sup>T</sup>A-orthogonal s-step Orthomin(k), and QMR algorithms are briefly presented and some of their properties are recalled. Particularly, it is shown that in the floating-point arithmetic there exist some cases in which the properties of these algorithms are lost, e.g., the result is false, or the coefficients of the GMRES residual polynomial are non significant and lead to serious round-off errors. The subject of this

MSC (2000): 65F10, 65F25, 65Go5

Keywords: Iterative methods, Hybrid methods, CESTAC method, stochastic arithmetic, CADNA library.

paper is to show how by using the CADNA library, it is possible during the run of the codes of these algorithms to detect the numerical instabilities, to stop correctly the process, and to evaluate the accuracy of the results provided by the computer. Numerical examples are used to show the good numerical properties.

#### 1. Introduction

In recent years there has been significant progress in development of iterative methods for solving sparse real linear systems of the form

$$Ax = b, (53)$$

where A is a nonsymmetric matrix of order n. The GMRES, hybrid GMRES,  $A^TA$ -orthogonal s-step Orthomin(k), and QMR algorithms are the most popular iterative methods for solving such systems. From the mathematical standpoint, these methods, based on a given initial point  $x^{(0)}$  considered as an approximation of the solution to the problem to be solved, consist in computing a sequence  $x^{(0)}, x^{(1)}, \ldots, x^{(n)}$  such that, if the method converge,  $x^{(n)}$  tends towards a limit  $x_s$ , when  $n \to \infty$ . Obviously it is impossible to reach this limit, and consequently termination criteria are proposed to stop the iterative process, such as

if 
$$||x^{(q)} - x^{(q-1)}|| \le \epsilon$$
 then stop,  
if  $||x^{(q)} - x^{(q-1)}|| \le \epsilon ||x^{(q)}||$  then stop,

where  $\epsilon$  is an arbitrary positive value. It is clear that these termination are not satisfactory because

$$||x^{(q)} - x^{(q-1)}|| \le \epsilon \iff ||x^{(q)} - x_s|| \le \epsilon,$$

$$||x^{(q)} - x^{(q-1)}|| \le \epsilon ||x^{(q)}|| \iff ||x^{(q)} - x_s|| \le \epsilon ||x_s||.$$

From the informatical standpoint, the situation is even more serious. In fact, if an iterative method is implemented on a computer, each  $X^{(q)} \in F$  only has a certain number of decimal significant digits. If the  $\epsilon$  selected is less than accuracy of  $X^{(q)}$ , these termination criteria are no longer meaningful. Thus two problems are raised.

- •How can the iteration process be stopped correctly?
- •What is the accuracy of the informatical solution given by the machine?

In addition, in the floating-point arithmetic there exist some cases in which the properties of these algorithms are lost, e.g., the coefficients of the GMRES residual polynomial are non-significant and lead to serious round-off errors, or the  $A^TA$ -orthogonal method has slow convergence, because of round-off error propagation.

In section 4, we observe that, the stochastic arithmetic allows the development of a termination criterion, called the "optimal termination criterion", which stops the iterative process as soon as a satisfactory informatical solution is reached, and the use of other appropriate criteria for overcoming the instabilities which exist in the algorithms.

In section 2 we briefly describe the GMRES, hybrid GMRES,  $A^TA$ -orthogonal s-step Orthomin(k) and QMR algorithms and discuss about the problems which exist in the implementation of these algorithms on a computer.

In section 3 we give a brief description of stochastic round-off analysis, the CESTAC method, and the CADNA software [25, 4]. Section 4 is devoted to the use of the CESTAC method and CADNA library for overcoming the problems which exist in the implementation of the mentioned algorithms on a computer using the floating-point arithmetic. Moreover, we will observe that by using the CADNA library and introducing the appropriate stopping criteria, it is possible, during the run of

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the code of these algorithms to detect the numerical instabilities and to stop correctly the iterative process, and to restart it in order to improve the computed solution. Some numerical results are given to show the good numerical properties.

# 2. Hybrid GMRES, $A^TA$ -orthogonal s-step Orthomin(k), and QMR algorithms

#### 2.1 GMRES and Hybrid GMRES algorithms

The GMRES method by Saad and Schultz [22], is one of the most popular methods for solving linear systems with a nonsymmetric matrix. The idea of GMRES is to construct an approximate solution of the form  $x_m = x_0 + z_m$  where  $z_m$  is an element of the Krylov subspace  $K^m(A; r_0) = \operatorname{span}\{r_o, Ar_0, \ldots, A^{m-1}r_0\}$  with the following property:

$$||r_m||_2 = ||b - Ax_m||_2 = \min_{z \in K^m(A;r_0)} ||r_0 - Az||_2,$$
 (54)

where  $r_0 = b - Ax_0$ . The basic structure of the GMRES algorithm is as follows.

#### Algorithm 2.1. GMRES

- 1. Compute  $r_0 = b Ax_0, \beta = ||r_0||_2$ , and  $v_1 = r_0/\beta$
- 2. Define the  $(m+1) \times m$  matrix  $\bar{H}_m = \{h_{ij}\}_{1 \le i \le m+1, 1 \le j \le m}$ . Set  $\bar{H}_m = 0$
- 3. For j = 1, 2, ..., m Do:
- 4. Compute  $w_i = Av_i$
- 5. For i = 1, 2, ..., j Do:
- $6. h_{ij} = (w_j, v_i)$

The stable iterative algorithms with the CADNA ...

$$7. w_i = w_i - h_{ij}v_i$$

- 8. EndDo
- 9.  $h_{j+1,j} = ||w_j||_2$ . If  $h_{j+1,j} = 0$  set m = j and goto 12
- 10.  $v_{j+1} = w_j/h_{j+1,j}$
- 11. EndDo
- 12. Compute  $y_m$  the minimizer of  $||\beta e_1 \bar{H}_m y||_2$  and  $x_m = x_0 + V_m y$

The GMRES iteration constructs a sequence of residual polynomials that minimize the norm of the residual

$$||r_m||_2 = ||p_m(A)r_0||_2 = \min_{\substack{p \in P_m \\ p(0) = 1}} ||p(A)r_0||_2, \ m = 1, 2, \dots.$$

With these GMRES polynomials the following hybrid GMRES is proposed in [20]:

Start with a random initial guess  $x_0$ .

Phase I: Run GMRES until  $||r_m||_2$  drops by a suitable amount. Set  $\nu = m$ .

Phase II: Re-apply the GMRES residual polynomial  $p_{\nu}(z)$  cyclically until convergence.

The main idea of this algorithm is to suppose that at the  $\nu th$  GMRES step the relation

$$\frac{\|r_{\nu}\|_{2}}{\|r_{0}\|_{2}} = \frac{\|p_{\nu}(A)r_{0}\|_{2}}{\|r_{0}\|_{2}} = \tau$$

holds for some  $\tau < 1$  , and moreover we have

$$||p_{\nu}(A)||_2 \simeq \tau.$$

So that by re-applying the GMRES polynomial  $p_{\nu}(z)$  cyclically we can reduce the residual norm. Of course, these assumptions do not always hold and it must modify the algorithm in order to cope with failure. By assumption that storage is not limited, we propose the following safeguarding procedure which differs from Nachtigal, Reichel, and Trefethen algorithm [20] in process 2:

- 1. If any cycle of  $\nu$  steps of phase II reduces  $||r_m||_2$  by a factor less than  $\sqrt{\tau}$  that is, if the convergence is more than twice as slow as expected return to phase I.
- 2. Carry out additional GMRES steps  $\nu+1, \nu+2, \ldots, \nu'$  of phase I until  $||r_{\nu'}||_2/|r_{\nu}||_2 < \frac{1}{2}$ , and calculate a new polynomial  $p_{\nu'}(z)$ .
- 3. Begin a new phase II iteration with the new polynomial  $p_{\nu'}(z)$ , starting from the previous best value  $x_m$ , which will come either from the previous phase II if the convergence there was slow but positive, or from the new phase I if there was actual divergence in the previous phase II.

The residual polynomial  $p_{\nu}(z)$  can be implicitly constructed by GMRES. The details of calculating the coefficients of  $p_{\nu}(z)$  can be found in [20].

When the hybrid GMRES is implemented on a computer, at each iteration q of this algorithm we have one approximate solution from phase I, denoted  $x_{0\nu}$ , and some approximate solutions from phase II, denoted  $x_{k\nu}, k=1,2,\ldots$  During the run of the program if any of these approximate solution is a satisfactory solution, in the sense that its residual norm is reduced by a factor  $\epsilon$  (where  $\epsilon$  is an arbitrary positive value), then the program can be stopped. So, for stopping the process at iteration q with  $\nu < n, k = 0, 1, \ldots$  we can use the following termination

criterion:

if 
$$||r_{k\nu}||_2 \le \epsilon ||r_0||_2$$
 then stop.

In the case  $\nu = n$ , the program must be stopped because the GMRES method converges, in the absence of rounding errors, in at most n iterations.

Note that in the above termination criterion  $\epsilon$  is an arbitrary value, the results of example 1 section 4 (Table 2, in which only the decimal significant digits are printed) show that when  $\epsilon$  is chosen too large ( $\epsilon = 10^{-6}$ ) the process is broken-off too early and consequently the solution obtained has a poor accuracy. On the contrary when  $\epsilon$  is chosen too small ( $\epsilon = 10^{-16}$ ) the iterative process is stopped too late and many useless iterations are performed, without improving the accuracy of the solution obtained with  $\epsilon = 10^{-15}$ . In practice it is absolutely impossible to choose correctly the value of the convergence tolerance  $\epsilon$ . But as explained in section 4 with CADNA library it is possible to break-off the iteration of an iterative process as soon as a satisfactory computed solution is reached, and this without using any arbitrary  $\epsilon$ .

Let us now consider the linear system

$$\begin{bmatrix} 21 & 130 & 0 & 2.1 \\ 13 & 80 & 4.74 + E8 & 752 \\ 0 & -0.4 & 3.9816 + E8 & 4.2 \\ 0 & 0 & 1.7 & 9E - 9 \end{bmatrix} .x = \begin{bmatrix} 153.1 \\ 849.74 \\ 7.7816 \\ 2.6E - 8 \end{bmatrix}, (55)$$

which was described in [25]. The exact solution is  $x = [1.0, 1.0, 10^{-8}, 1.0]^T$ . The approximate solution obtained with the initial guess  $x_0 = [0, 0, \dots, 0]^T$  and the FORTRAN code of the hybrid GMRES algorithm, performed on a SUN4 computer, in double precision is as follows

$$x(1) = -89.8760751972095591,$$
  $x(2) = 15.6799813793030225,$   $x(3) = 2.4747823346160658E - 08,$   $x(4) = 1.0000000003115719.$ 

It is necessary to say that this approximate solution has been obtained by phase I with  $\nu=4$ . This solution is false and hybrid GMRES algorithm is not able, because of the propagation of the round-off errors, to provide a satisfactory solution for this example. However, nothing in the software nor in the solution allows the user to be aware that the computed solution is false. In section 3 we will show that the CADNA library is able to estimate the propagation of round-off errors during the run of the code and to furnish the accuracy of the computed solution.

Let us consider another linear system with the  $n \times n$  block diagonal matrix

$$A = \left[ egin{array}{cccc} M_1 & & & & \\ & M_2 & & & \\ & & \ddots & & \\ & & & M_{n/2} \end{array} 
ight], \quad ext{where } M_j = \left[ egin{array}{cccc} 1 & j-1+lpha \\ 0 & -1 & \\ 56 \end{array} 
ight],$$

which with  $\alpha=0$  corresponds to the example  $B_{\pm 1}$  of [21], and the second member  $b=[5,-3,4,-4,1,\ldots,1]^T$ . The exact solution is given by  $x_{2j-1}=b_{2j-1}+(j-1+\alpha)b_{2j}$  and  $x_{2j}=-b_{2j}$ . With  $\alpha=1.11,\,n=150$ , and the initial guess  $x_0=[0,0,\ldots,0]^T$  the computed coefficients  $\alpha_i$  of the GMRES polynomial  $p_6(z)$  with single and double floating-point arithmetic rounded to nearest mode are presented in Table 1. In section 4, thanks to the CADNA library, we consider that these coefficients are non significants. It is clear that, phase II is not able here to improve the approximate solution which has been obtained by phase I. In general, when  $\nu$  has a large value, it is possible to have an unstable solution  $x_{k\nu}, k=1,2,\ldots$  and also a GMRES polynomial with unsignificant coefficients. In this situation the use of phase II has obviously no sense. So, in order to avoid the performation of many useless operations of phase II and to prevent an overflow which may occur, it is better the program perform only phase I as soon as such an instability occurs during the

run. We now face the question of how can detect these kind of instabilities? The CADNA library is a precious tool for obtaining an answer to this question. In section 4 we show that with CADNA library it is possible, by including a simple test to detect it.

	single precision	double precision
$\alpha_0$	-2.9381578E+05	1.6409978460272019E+14
$\alpha_1$	3.0272866E+05	-1.6409978460272156E+14
$\alpha_2$	5.8763362E+05	-3.2819956920543981E+14
$\alpha_3$	-6.0545688E+05	3.2819956920544550E+14
$\alpha_4$	-2.9381784E+05	1.6409978460271962E+14
$lpha_5$	3.0272922E+05	-1.6409978460272297E+14

Table 1

### 2.2 $A^TA$ -orthogonal s-step Orthomin(k) algorithm

In [7], Chronopoulos developes  $A^TA$ -orthogonal s-step Orthomin(k) algorithm for nonsymmetric matrices with symmetric part  $M=(A+A^T)/2$  positive definite or indefinite. In this method the s directions  $\{r_i,\ldots,A^{s-1}r_i\}$  are formed and are  $A^TA$ -orthogonalized simultaneously to k of the preceding directions  $\{p_j^1,\ldots,p_j^s\}, j=j_i,\ldots,i$  where  $j_i=max(0,i-k+1)$ . The norm of residual  $||r_{i+1}||_2$  is minimized simultaneously in all s new directions in order to obtain  $x_{i+1}$ . More details of the s-step Orthomin(k) algorithm can be found in [7]. The following notation facilitates the description of the algorithm.

$$W_i = [(Ap_i^j, Ap_i^l)], \text{ where } 1 \leq j, l \leq s$$

$$\underline{a_i} = [a_i^1, \dots, a_i^s]^T \text{ (the steplengths in updating } x_i \text{ )}$$

$$\underline{m_i} = [(r_i, Ap_i^1), \dots, (r_i, Ap_i^s)]^T$$

$$\underline{c_i^l} = [(A^l r_{i+1}, Ap_i^1), \dots, (A^l r_{i+1}, Ap_i^s)]^T$$

$$\underline{b}_j^l = \{b_j^{l,m}\}_{m=1}^s$$
 for  $j=j_i,\ldots,i$  and  $l=1,\ldots,s$ , where  $j_i=max(0,i-k+1)$ 

(the coefficients to  $A^TA$ -orthogonalize to the previous directions)

$$P_i = [p_i^1, \dots, p_i^s]$$
 (the direction vectors)

$$R_i = [r_i, Ar_i, \dots, A^{s-1}r_i]$$
 (the residuals).

A description of s-step Orthomin(k) method can be given as follows:

Algorithm 2.2. s-step Orthomin(k)

Select  $x_0$ 

$$P_0 = [r_0 = b - Ax_0, Ar_0, \dots, A^{s-1}r_0]$$

For i = 0 Until Convergence Do

Compute mi, Wi

Call Scalar1

$$x_{i+1} = x_i + P_i \underline{a}_i$$

$$r_{i+1} = r_i - AP_i\underline{a}_i$$

Compute  $\underline{c}_i^i, j = j_i, \dots, i$ 

Call Scalar2

Compute 
$$R_{i+1} = [r_{i+1}, Ar_{i+1}, \dots, A^{s-1}r_{i+1}]$$

$$P_{i+1} = R_{i+1} + \sum_{j=i}^{i} P_{j} [\underline{b}_{i}^{l}]_{l=1}^{s}$$

Compute  $AP_{i+1}$  or,

$$AP_{i+1} = AR_{i+1} + \sum_{j=j_i}^{i} AP_j[\underline{b}_j^l]_{l=1}^{s}$$

#### EndFor

Scalar1: Decomposes  $W_i$  and solves  $W_i\underline{a}_i = \underline{m}_i$ .

Scalar2: Solves 
$$W_j \underline{b}_j^l = -\underline{c}_j^l$$
 for  $j = j_i, \ldots, i$  and  $l = 1, \ldots, s$ , where  $j_i = max(0, i - k + 1)$ 

The solution of the linear systems may cause a quick loss of orthogonality of the s-dimensional direction subspaces  $P_i$  because the matrix  $W_i$  may have a very large condition number. Numerical tests [8, 9, 10] have shown that the condition number of  $W_i$  is small for  $s \leq 5$ . One way to alleviate the orthogonality loss which can occur for large s > 5 is to  $A^TA$ -orthogonalize the s direction vectors in each iteration. In [23],  $A^TA$ -orthogonal s-step Orthomin(k) was developed and shown to be stable for large values of s(up to s = 16). In this method the direction vectors within each subspace  $P_i$  are  $A^TA$ -orthogonalized using the Modified Gram-Schmidt method. The linear systems need not be solved at each iteration since the  $W_i$  matrix is the identity matrix if  $P_i$  is perfectly  $A^TA$ -orthogonalized. By using the notation  $j_i = max(0, i - k + 1)$  the algorithm can be described as follows:

Algorithm 2.3.  $A^T A$ -orthogonal s-step Orthomin(k)

Select  $x_0$ 

Compute  $r_0 = b - Ax_0$ 

For i = 0 Until Convergence Do

Compute 
$$AP_i = [Ar_i, A^2r_i, \dots, A^sr_i],$$
  
and set  $P_i = [r_i, Ar_i, \dots, A^{s-1}r_i]$ 

If 
$$(0 < i)$$
 Then

Compute 
$$\underline{b}_{j}^{l} = [-(A^{l}r_{i}, Ap_{j}^{1}), \dots, -(A^{l}r_{i}, Ap_{j}^{s})]^{T}$$
, for  $l = 1, \dots, s$  and  $j = j_{i-1}, \dots, i-1$ 

Compute  $P_{i} = P_{i} + \sum_{j=j_{(i-1)}}^{i-1} P_{j}[\underline{b}_{j}^{l}]_{l=1}^{s}$ 

Compute  $AP_{i} = AP_{i} + \sum_{j=j_{(i-1)}}^{i-1} AP_{j}[\underline{b}_{j}^{l}]_{l=1}^{s}$ 

#### EndIf

Apply the Modified Gram-Schmidt method to the matrix  $AP_i$  to obtain final  $AP_i$  and  $P_i$ 

Compute 
$$\underline{a}_i = [(r_i, Ap_i^1), \dots, (r_i, Ap_i^s)]^T$$

$$x_{i+1} = x_i + P_i \underline{a}_i$$

$$r_{i+1} = r_i - AP_i \underline{a}_i$$

#### EndFor.

The main problem in the use of  $A^TA$ -orthogonal s-step Orthomin(k) method, with floating-point arithmetic, is the choice of s. Let us consider the results of this method with different values of s and k for the examples 4-6 of section 4.2 (Tables 6, 8, and 10, in which the number of iterations to convergence are printed). These results clearly show that when s has a small or large value the method has slow convergence, and for each problem and each k there exists an s which minimize the number of iterations to convergence. However, as above mentioned, the slow convergence of the method with large value of s is due to the round-off errors propagation. Hence, it is not possible to determine a good value of s without estimating the round-off errors propagation. In section 4, it is shown that by using the CADNA library, which is an efficient tool for doing so, we will be able to determine a good value of s.

Here, another problem is also the choice of the value  $\epsilon$  for stopping criterion  $||r_i||_2 \leq \epsilon$ . When  $\epsilon$  is chosen too large, the iterative process is stopped too soon, and consequently the solution obtained has a poor accuracy. On the contrary, when  $\epsilon$  is chosen small, it is possible, due to the numerical instabilities, many useless iterations are performed without improving the accuracy of the solution. How can the iterative processs be stopped correctly, and restarted in order to improve the computed solution? The CADNA library is a precious tool for obtaining an answer to this question. In section 4 we will show that with CADNA library, it is possible, by including the simple tests to stop and to restart correctly the iterative process.

#### 2.3 The QMR algorithm

The quasi-minimal residual method (QMR) is based on the look-ahead Lanczos algorithm proposed in [13]. In the following,  $A \in \mathbb{C}^{N \times N}$  is always assumed to be a given  $N \times N$  matrix. Let  $v_1, w_1 \in \mathbb{C}^N$  be any two vectors different from the zero vector. Starting with  $v_1, w_1$ , the lookahead Lanczos algorithm generate two sequences of vectors  $v_1, v_2, \dots, v_n$  and  $w_1, w_2, \dots, w_n, n = 1, 2, \dots$ , that satisfy

$$span\{v_1, v_2, \dots, v_n\} = K_n(v_1, A), 
span\{w_1, w_2, \dots, w_n\} = K_n(w_1, A^T),$$
(57)

and can be grouped into k = k(n) blocks

$$V_{l} = [v_{n_{l}}v_{n_{l+1}} \cdots v_{n_{l+1}-1}], \qquad W_{l} = [w_{n_{l}}w_{n_{l+1}} \cdots w_{n_{l+1}-1}],$$

$$l = 1, 2, \cdots, k-1,$$

$$V_{k} = [v_{n_{k}}v_{n_{k}+1} \cdots v_{n}], \qquad W_{k} = [w_{n_{k}}w_{n_{k}+1} \cdots w_{n}],$$

where

$$1 = n_1 < n_2 < \cdots < n_l < \cdots < n_k < n < n_{k+1}.$$

The blocks are constructed such that we have

$$W_j^T V_l = \left\{ egin{array}{ll} 0 & ext{if } j 
eq l, \ D_l & ext{if } j = l, \end{array} 
ight. \quad j,l = 1,2,\cdots,k,$$

where

$$D_l$$
 is nonsingular,  $l=1,2,\cdots,k-1$ , and  $D_k$  is nonsingular if  $n=n_{k+1}-1$ .

The first vectors  $v_{n_i}$  and  $w_{n_i}$  in each block are called *regular*, the remaining vectors are called *inner*. The kth block is called *complete* if  $n = n_{k+1} - 1$ ; in this case, at the next step n + 1, a new block is started with the regular vectors  $v_{n_{k+1}}$  and  $w_{n_{k+1}}$ . Otherwise, if  $n < n_{k+1} - 1$ , the kth block is *incomplete* and at the next step, the Lanczos vectors  $v_{n+1}$  and  $w_{n+1}$  are added to the kth block as inner vectors.

With these preliminaries, the basic structure of the look-ahead Lanczos algorithm is as follows.

Algorithm 2.4 sketch of the look-ahead Lanczos algorithm [13].

0) Choose 
$$v_1, w_1 \in \mathbb{C}^N$$
 with  $||v_1|| = ||w_1|| = 1$ ;  
Set  $V_1 = v_1, W_1 = w_1, D_1 = W_1^T V_1$ ;  
Set  $n_1 = 1, k = 1, v_0 = w_0 = 0, V_0 = W_0 = \emptyset, \rho_1 = \xi_1 = 1$ ;  
For  $n = 1, 2, \cdots$  do:

- 1) Decide whether to construct  $v_{n+1}$  and  $w_{n+1}$  as regular or inner vectors and go to 2) or 3), respectively;
- $2)(Regular\ step.)\ Compute$

$$\begin{split} \tilde{v}_{n+1} &= A v_n - V_k D_k^{-1} W_k^T A v_n - V_{k-1} D_{k-1}^{-1} W_{k-1}^T A v_n, \\ \tilde{w}_{n+1} &= A^T w_n - W_k D_k^{-T} V_k^T A^T w_n - W_{k-1} D_{k-1}^{-T} V_{k-1}^T A^T w_n, \end{split}$$
 (58)

set 
$$n_{k+1} = n + 1, k = k + 1, V_k = W_k = \emptyset$$
, and go to 4);

3)(Inner step.) Compute

$$\begin{split} \tilde{v}_{n+1} &= A v_n - \zeta_{n-n_k} v_n - (\eta_{n-n_k}/\rho_n) v_{n-1} - V_{k-1} D_{k-1}^{-1} W_{k-1}^T A v_n, \\ \tilde{w}_{n+1} &= A^T w_n - \zeta_{n-n_k} w_n - (\eta_{n-n_k}/\xi_n) w_{n-1} - W_{k-1} D_{k-1}^{-T} V_{k-1}^T A^T w_n; \\ (59) \end{split}$$

4) Compute  $\rho_{n+1} = ||\tilde{v}_{n+1}||$  and  $\xi_{n+1} = ||\tilde{w}_{n+1}||$ ;

If 
$$\rho_{n+1} = 0$$
 or  $\xi_{n+1} = 0$ , stop;

Otherwise, set

$$v_{n+1} = \tilde{v}_{n+1}/\rho_{n+1}, \qquad w_{n+1} = \tilde{w}_{n+1}/\xi_{n+1}, V_k = [V_k \ v_{n+1}], \quad W_k = [W_k \ w_{n+1}], \quad D_k = W_k^T V_k.$$
(60)

Now, we list some properties of Algorithm 2.4 which will be used in the sequel. First, in view of (8), we have

$$||v_n|| = ||w_n|| = 1, \quad n = 1, 2, \dots$$
 (61)

It is convenient to introduce the notation

$$V^{(n)} = [v_1 \ v_2 \dots v_n] \quad (= [V_1 \ V_2 \dots V_k]),$$
  

$$W^{(n)} = [w_1 w_2 \dots w_n] \quad (= [W_1 W_2 \dots W_k]).$$
(62)

Hence, by (5),

$$K_n(v_1, A) = \{ V^{(n)}z \mid z \in \mathbb{C}^N \}, K_n(w_1, A^T) = \{ W^{(n)}z \mid z \in \mathbb{C}^N \}.$$
 (63)

Moreover, the recursions for the v's in (6) and (7) can be rewritten in matrix formulation as follows:

$$AV^{(n)} = V^{(n)}H_n + [0 \dots 0 \tilde{v}_{n+1}]. \tag{64}$$

Here,

$$H_{n} := \begin{bmatrix} \alpha_{1} & \beta_{2} & 0 & \cdots & 0 \\ \gamma_{2} & \alpha_{2} & \beta_{3} & \ddots & \vdots \\ 0 & \gamma_{3} & \ddots & \ddots & 0 \\ \vdots & \ddots & \ddots & \ddots & \beta_{k} \\ 0 & \cdots & 0 & \gamma_{k} & \alpha_{k} \end{bmatrix}$$

$$(65)$$

is a block-tridiagonal matrix of size  $n \times n$  that also is upper Hessenberg. Furthermore, the diagonal blocks  $\alpha_1, \alpha_2, \dots, \alpha_k$  of the matrix  $H_n$  are all squares, and their sizes are just the lengths of the look-ahead steps.

Let  $x_0 \in \mathbb{C}^N$  be an arbitrary initial guess for the solution of (1), and set  $r_0 = b - Ax_0$ . At the n<sup>th</sup> iteration QMR computes an approximate solutions of (1) of the form

$$x_n \in x_0 + K_n(r_0, A).$$
 (66)

If we choose  $v_1 = r_0/\|\rho_0\|$  and any  $w_1 \in \mathbb{C}^N$ ,  $\|w_1\|_2 = 1$ , as the starting vectors for the look-ahead Lanczos algorithm, then, by (11), the right Lanczos vectors  $v_1, v_2, \dots, v_n$  span Krylov subspace  $K_n(r_0, A)$  in (14). Therefore, any iterate (14) can be presented in the form

$$x_n = x_0 + V^{(n)} z_n, \quad z_n \in \mathbb{C}^n, \tag{67}$$

and  $V^{(n)}$  is the matrix defined in (10). For this approximate solution  $x_n$  the residual satisfies

$$r_n = V^{(n+1)}(f_{n+1} - H_n^{(e)} z_n).$$
(68)

where  $f_{n+1} = [||r_0||_2 \ 0 \ \cdots \ 0]^T \in IR^{n+1}$ , and  $H_n^{(e)}$ , defined by

$$H_n^{(e)} = \begin{bmatrix} H_n \\ \rho_{n+1}(e_n^{(n)})^T \end{bmatrix}, \quad (e_n^{(n)})^T = [0 \dots 0 \ 1]^T \in \mathbb{R}^n. \quad (69)$$

Freund and Nachigal [14] suggested to choose  $z_n$  as the solution of the least-squares problem

$$|| f_{n+1} - H_n^{(e)} z_n ||_2 = \min_{z \in \mathbb{C}^n} || f_{n+1} - H_n^{(e)} z ||_2.$$
 (70)

The motivation of this choice of  $z_n$  is as follows. The vector z which minimizes (18) can be found with considerably less work than would be needed to minimize the residual norm, since the matrix  $V^{n+1}$  will not usually be unitary. More details of the QMR method can be found in [14].

We observe that in each step of a look-ahead Lanczos process, it is necessary to decide whether to construct the Lanczos vectors  $v_{n+1}$  and  $w_{n+1}$  as regular or inner vectors. As we know [14], for a regular step it is necessary that  $D_k = W_k^T V_k$  is nonsingular. Therefore, in implementation of QMR in the floating-point arithmetic, the smallest singular value of matrix  $D_k$  is computed and the criterion

$$\sigma_{min}(D_k) \leq Tol$$

is used to check whether this matrix is singular or close to singular, and to decide whether to construct the Lanczos vectors  $v_{n+1}$  and  $w_{n+1}$  as regular or inner vectors. Here Tol is a suitable chosen tolerance. The efficiency of the algorithm depends on a good choice of the Tol and to construct correctly the Lanczos vectors  $v_{n+1}$  and  $w_{n+1}$ . In addition, if the quantity  $\sigma_{min}(D_k)$  is badly computed, propagation of round-off errors will affect drastically all the computation. In section 4, we show that QMR algorithm with the CADNA library, and using the appropriate test, and optimal terminations, it is possible in constructing the Lanczos vectors  $v_{n+1}$  and  $w_{n+1}$  to decide correctly, to restart the QMR algorithm if it is necessary, to stop the program as soon as a satisfactory solution is reached, to estimate the accuracy of the solution, and to save computer time, because many useless operations and iterations are not performed.

In the following section we give a brief description of CESTAC method which is an efficient method for solving the numerical problems such as those described above.

# 3. The CESTAC method

# 3.1 Basic ideas of the CESTAC method

Any result R provided by a computer always contains an error resulting from round-off errors propagation. It has been proved [2] that a computed result R is modelized to the first order in  $2^{-p}$  as the equation

$$R = r + \sum_{i=1}^{n} u_i(d) 2^{-p} \alpha_i,$$

where r is the exact result,  $\alpha_i$  is the round-off error, and  $u_i(d)$  are quantities depending exclusively on the data. The integer n is the number of arithmetical operations involved in the computation of R, and the integer p is the number of bits in the mantissa.

The CESTAC method (Contrôle et Estimation Stochastique des Arrondis de Calculs) was developed by La Porte and Vignes, and was then generalized by the latter. It is based on a probabilistic approach of the round-off errors propagation, it has been presented in [12, 16, 17], this method allows to estimate the round-off error on each result and consequently provides the accuracy of this result.

The basic idea of this method consists in performing the same code several times in order to propagate the round-off error differently each time. Several samples of R containing different round-off error are then obtained. The first digits common to all the samples are significant and the others are not significant and represent the round-off error propagation. The aim is then to obtain these samples of R. They are obtained by the use of random arithmetic.

Indeed, each result r of any floating-point (FP) arithmetical operator is always bounded by two consecutive FP values  $R^-$  and  $R^+$ . The random arithmetic consists in randomly choosing either  $R^-$  or  $R^+$  with a probability 0.5. Then when a same code is executed N times with a computer using this random arithmetic, for each result of any floating-point arithmetic, N different results  $R_i$ ,  $i = 1, \ldots, N$  will be provided. It has been proved [2, 5] that, under certain hypothesis, this N results belong to a quasi-Gaussian distribution centered on the exact result r. So, in practice, the use of the CESTAC method consists in :

- (i) Running in parallel N times (N=2 or 3) the program with this new arithmetic. Consequently, for each result R of any floating-point arithmetic operation, a set of N computed results  $R_i, i=1,\ldots,N$  is obtained.
- (ii) Taking the mean value  $\bar{R} = \frac{1}{N} \sum_{i=1}^{N} R_i$  of the  $R_i$  as the computed result.
- (iii) Using Student distribution to estimate a confidence interval for R, and then compute the number  $C_{\bar{R}}$  of significant digits of  $\bar{R}$ , defined by  $C_{\bar{R}} = \log_{10}(\sqrt{N}|\bar{R}|/\tau_{\beta}\sigma)$ , with  $\sigma^2 = \frac{1}{N-1}\sum_{i=1}^{N}(R_i \bar{R})^2, \tau_{\beta}$  is the value of the Student distribution for N-1 degrees of freedom and a probability level  $1-\beta$ .

#### 3.2 Stochastic arithmetic

By using the CESTAC method so that the N runs of the computer program take place in parallel, the N results of each arithmetic operation can be considered as realisations of Gaussian random variable centered on the exact result. We can therefore define a new number, called stochastic number, and a new arithmetic, called stochastic arith-

metic, applied to these numbers. We present below the main definitions and properties of this arithmetic. For more details see [6]

**Definition 1.** We define the set S of stochastic numbers as the set of Gaussian random variables. We denote an element  $X \in S$  by  $X = (\mu, \sigma^2)$ , where  $\mu$  is the mean value of X and  $\sigma$  its standard deviation. If  $X \in S$  and  $X = (\mu, \sigma^2)$ , there exists  $\lambda_{\beta}$ , depending only on  $\beta$ , such that:

$$P(X \in [\mu - \lambda_{\beta}.\sigma, \ \mu + \lambda_{\beta}.\sigma]) = 1 - \beta.$$

 $I_{\beta,X} = [\mu - \lambda_{\beta}.\sigma, \ \mu + \lambda_{\beta}.\sigma]$  is a confidence interval of  $\mu$  at  $(1 - \beta)$ . An upper bound to the number of significant digits common to  $\mu$  and each element of  $I_{\beta,X}$  is

$$C_{\beta,X} = \log_{10}(\frac{|\mu|}{\lambda_{\beta} \cdot \sigma}).$$

The following definition is the modelling of the concept of informatical zero proposed in [24].

**Definition 2.**  $X \in S$  is a stochastic zero, denoted  $\underline{0}$ , if and only if:

$$C_{\beta,X} \leq 0$$
 or  $X = (0,0)$ .

**Definition 3.** Let  $X_1 = (\mu_1, \sigma_1^2)$ ,  $X_2 = (\mu_2, \sigma_2^2)$  be two elements of S. We define the four elementary stochastic operations denoted (s+, s-, s\*, s/) on the stochastic numbers by

$$\begin{array}{lll} X_1 \; s + \; X_2 & \stackrel{def}{=} & (\mu_1 + \mu_2, \; \sigma_1^2 + \sigma_2^2), \\ X_1 \; s - \; X_2 & \stackrel{def}{=} & (\mu_1 - \mu_2, \; \sigma_1^2 + \sigma_2^2), \\ X_1 \; s * \; X_2 & \stackrel{def}{=} & (\mu_1 * \mu_2, \; \mu_2^2 \sigma_1^2 + \mu_1^2 \sigma_2^2), \\ X_1 \; s / \; X_2 & \stackrel{def}{=} & (\mu_1 / \mu_2, \; (\frac{\sigma_1}{\mu_2})^2 + (\frac{\mu_1 \sigma_2}{\mu_2^2})^2), & \text{with } \mu_2 \neq 0. \end{array}$$

**Definition 4.** Let  $X_1$  and  $X_2$  be two elements of S,  $X_1$  is stochastically equal to  $X_2$ , denoted  $X_1$   $s = X_2$ , if and only if

$$X_1 s - X_2 = \underline{0}.$$

**Definition 5.** Let  $X_1 = (\mu_1, \sigma_1^2)$  and  $X_2 = (\mu_2, \sigma_2^2)$  be two elements of S.  $X_1$  is stochastically strictly greater than  $X_2$ , denoted  $X_1$   $s > X_2$ , if and only if:

$$\mu_1 - \mu_2 > \lambda_\beta \sqrt{\sigma_1^2 + \sigma_2^2}$$

**Definition 6.** Let  $X_1 = (\mu_1, \sigma_1^2)$ ,  $X_2 = (\mu_2, \sigma_2^2)$  be elements of S.  $X_1$  is stochastically greater than  $X_2$ , denoted  $X_1$   $s \ge X_2$ , if and only if:

$$X_1 \quad s > X_2 \quad or \quad X_1 \quad s = X_2$$

Based on these definitions, the following properties of stochastic arithmetic have been proved:

- s = is reflexive and symmetric but is not transitive,
- s > is transitive
- $s \ge$  is reflexive, anti-symmetric, but is not transitive;

$$a \ s \le b \Rightarrow a \ s = b \ or \ a \ s < b$$
 $a \ s = b \ and \ b \ s < c \Rightarrow a \ s \le c$ 
 $a \ s \le b \ and \ b \ s < c \Rightarrow a \ s \le c$ 

 $\underline{0}$  is absorbent for operation s\* and is the neutral element for operation s+. Let  $x,y\in I\!\!R$  and  $X,Y\in S$  respectively be their representative. If X s < Y  $\Rightarrow$  x < y. Thus stochastic arithmetic retrieves properties of exact arithmetic, lost by usual floating-point arithmetic such as associativity, distributivity, the concept of remarkable identities.

On computer, by using the synchronous implementation of the CESTAC method and by identifying the notions of informatical zero and stochastic zero, it is possible to use the stochastic arithmetic with its definitions. The use of stochastic arithmetic in scientific codes allows:

 during the run of a scientific code, to estimate the accuracy of any numerical result, to detect the numerical instabilities, and to check the branchings; 248 Faezeh Toutounian

2. to eliminate the programming expedients that are absolutely unfounded, such as those used, for example, in termination criteria of iterative methods, and replace them by criteria that directly reflect the mathematical condition that must be satisfied at the solution.

## 3.3 The CADNA library

CADNA (Control of Accuracy and Debugging for Numerical Applications) is a library for programs written in FORTRAN 77, FORTRAN 90, or in ADA which allows the computation using stochastic arithmetic by automatically implementing the CESTAC method [3, 4]. CADNA is able to estimate the accuracy of the computed results, and to detect numerical instabilities occuring during the run. To use the CADNA library, it suffices to place the instruction USE CADNA at the top of the initial FORTRAN or ADA source code and to replace the declarations of the real type by the stochastic type and to change some statements as printing statements.

During the run, as soon as a numerical anomaly (for example, appearance of informat ical zero in a computation or a criterion) occurs, a message is written in a special file called Cadna\_stability\_f90.lst. The user must consult this file after the program has run. If it is empty, this means the program has been run without any problem, that it has accordingly been validated, and that the results have been given with their associated accuracy. If it contains messages, the user, using the debugger associated with the compiler, will find the instructions that are the cause of these numerical anomalies, and must reflect in order to correct them if necessary. The program execution time using the CADNA library is only multiplied by a factor 3, which is perfectly acceptable in view of the major advantage offered, i.e., the validation of

programs. CADNA is also able to estimate the influence of data errors on the result provided by the computer.

## 4. Using the CADNA library in iterative methods

### 4.1 Using the CADNA library in hybrid GMRES method

As we have seen in section 2.1, in implementation of hybrid GMRES method three difficulties arise. Let us first consider the third problem which concerns to the numerical instabilities which may occur in phase II. By remarking that a GMRES polynomial with unsignificant coefficients and also an unstable solution  $x_{k\nu}$ ,  $k=1,2,\ldots$  always provide an unsignificant residual norm with large magnitude, we can define with CADNA library the test

if 
$$||r_{k\nu}||_2 = \underline{0}$$
 and  $||r_{k\nu}||_2 \ge 1.0$  then  $index = 1$ , (71)

with initial value index = 0, which allows us by checking the value of parameter index to detect the presence of an unstable solution. As consequence, the program, during the run, can decide to perform the both phase I and II, or only the phase I according to the value of this parameter index.

Now, we consider the first problem which concerns to choose a stopping criterion. As explained in [24], from mathematical point of view, once we know a solution  $x_m$ , we can validate its validity by checking the value of the residual norm

$$||r_m||_2 = ||b - Ax_m||_2 = 0.$$

Obviously with usual floating arithmetic this equality is never satisfied even when the solution  $x_m$  is the exact solution, because of the round-

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off errors propagation. However, with CADNA library in view of its properties, the result will be

$$||r_m||_2 = ||b - Ax_m||_2 = \underline{0}.$$

So, as soon as the residual norm of a stable computed solution is equal to the informatical zero, a satisfactory informatical solution is reached and the iterative process must be stopped. Now, by noting that it is possible, in phase II, to have an unstable solution  $x_{k\nu}$ ,  $k=1,2,\ldots$  and we can detect it by the test (19), we define the termination criterion

$$if ||r_{k\nu}||_2 = \underline{0} \quad then stop, \tag{72}$$

for checking the value of the residual norm of a stable solution  $x_{k\nu}, k=0,1,\ldots$ . This termination criterion stops the iterative hybrid GMRES process as soon as a satisfactory solution is reached either by phase I or by phase II. It is necessary to say that the instability of the solution  $x_{k\nu}, k=1,2,\ldots$  must be checked before using the termination criterion (20).

Finally the second problem, which concerns the accuracy of the computed solution, will be solved by using the CADNA library. Since, as it was explained in section 3.3, the CADNA library is able to estimate the accuracy of the computed results and to furnish the results with their exact decimal figures.

Let us now, to present the examples and the results which we obtained by the FORTRAN code of hybrid GMRES method, combined with CADNA library and the above tests. Computation have been performed on a SUN4 computer in double or simple precision with the stochastic arithmetic using the CADNA library.

Example 1. We consider the linear system with

$$A = \begin{bmatrix} 2 & 1 & & & & \\ 0 & 2 & 1 & & & \\ 1 & 0 & 2 & \ddots & & \\ & \ddots & \ddots & \ddots & & \\ & & 1 & 0 & 2 & 1 \\ & & & 1 & 0 & 2 \end{bmatrix}, \quad b = \begin{bmatrix} 3 \\ 3 \\ 4 \\ \vdots \\ 4 \\ 3 \end{bmatrix},$$

which was described in [15], and the dimension equal to 400. The exact solution is given by  $x = [1, 1, ..., 1]^T$ . The results obtained by using floating-point arithmetic with  $\epsilon = 10^{-6}, 10^{-15}, 10^{-16}$ , and CADNA library are presented in Table 2. Only the decimal significant digits are printed. At the last lines of this table we can find the corresponding value of indices  $\nu, k$  of the presented solution  $x_{k\nu}$ .

For this example the initial guess has been  $x_0 = [0,0,\dots,0]^T$ . From the results of Table 2, it appears that when  $\epsilon$  has too large value, the iterative process is stopped too soon (at  $\nu=4,k=11$ ), and the solution furnished is not the best that the computer may provide. When  $\epsilon$  is chosen too small ( $\epsilon=10^{-16}$ ) the iterative process is stopped too late (at  $\nu=6,k=1$ ), many useless iterations are performed without improving the accuracy of the solution with  $\epsilon=10^{-15}$ . By using the CADNA library, the optimal termination criterion (20) has stopped the iterative process at  $\nu=4,k=48$ , and solution is reached with about 15 exact significant digits on all the elements.

	$\epsilon = 10^{-6}$	$\epsilon = 10^{-15}$	$\epsilon = 10^{-16}$	CADNA library
x(1)	1.0000000	1.00000000000000000	1.00000000000000000	0.1000000000000000E+1
x(2)	0.999999	1.00000000000000000	1.00000000000000000	0.9999999999999E+0
x(3)	1.000000	1.0000000000000000	1.0000000000000000	0.1000000000000000E+1
x(4)	1.000000	0.99999999999999	0.999999999999999	0.100000000000000E+1
x(5)	0.999999	1.000000000000000000	1.00000000000000000	0.9999999999999E+0
:		:		•
x(396)	1.000000	1.000000000000000000	1.000000000000000000	0.999999999999E+0
x(397)	0.9999999	1.000000000000000000	1.00000000000000000	0.100000000000000E+1
x(398)	0.9999999	1.0000000000000000	1.000000000000000	0.100000000000000E+1
x(399)	1.000000	1.00000000000000000	1.000000000000000000	0.9999999999999E+0
x(400)	0.9999999	1.00000000000000000	1.00000000000000000	0.100000000000000E+1
	u = 4	$\nu = 4$	$\nu = 6$	$\nu = 4$
	k = 11	k = 46	k = 1	k = 48

Table 2

Remark: For controling the quality of a computed solution X of a linear system Ax = b we can use the normalized residuals test. As explained in [18, 19], this test consists in computing the normalized residuals

$$ho_i^* = rac{|
ho_i|}{2^{-p} \sqrt{m_i^q (\Sigma_{j=1}^n (AijX_j)^2 + B_i^2)}}, \quad i = 1, \ldots, n,$$

where

$$\rho_i = B_i - \sum_{j=1}^n A_{ij} X_j$$

and  $A_{ij}$ ,  $B_i$  are the normalized floating-point representations of  $a_{ij}$  and  $b_i$ , respectively,  $X_j$  is the jth element of computed solution X, and q=1 for rounding to the nearest mode, and q=2 for other rounding modes. The integer  $m_i$  is the number of nonzero elements of row i, and the integer p is the number of bits in the mantissa.

The three following cases can occur:

case I: All the n normalized residuals are of the order of mag-

nitude 1:

$$\rho_i^* \sim 1, \quad \forall i \in [1, 2, \dots, n],$$

thus the computed solution X is a satisfactory informatical solution.

case II: At least one of the normalized residuals is strictly strictly greater than one, but strictly strictly smaller than  $2^p$ :

$$1 \ll \rho_i^* \ll 2^p,$$

thus the computed solution X is not a satisfactory informatical solution, but it is possible to improve it by an increment vector  $\Delta X$  which may be obtained by solving the linear system  $A\Delta X = R$ , where R is the residual vector with ith element  $\rho_i$ .

case III: At least one of the normalized residuals is of the order of magnitude  $2^p$ :

$$\rho_i^* \sim 2^p$$

thus X is a bad solution and in general, we can not improve it. If this situation occurs, this means that the using method is not adapted to the proposed system.

By applying this test to the computed solutions of the above example, we discover that all the above computed solutions are the satisfactory informatical solutions except the one which obtained by floating-point arithmetic with  $\epsilon=10^{-6}$ . By improving this solution two times, we could obtain a satisfactory informatical solution. We observe that without using CADNA library, it is very difficult to obtain a satisfactory informatical solution.

Example 2. Let us again consider the linear system (3). The solution obtained with CADNA library is as follows

$$x(1) = \underline{0},$$
  $x(2) = \underline{0},$   $x(3) = \underline{0},$   $x(4) = 0.9999999E + 000.$ 

These results show that the first three elements of computed solution are non significant and the last one has 7 significant digits. We observe that only by using the CADNA library it is possible to conclude that the results obtained for the first three elements of computed solution must be due to the round-off errors propagation.

**Example 3.** Let us again consider the linear system (4). The solutions furnished by using single floating-point arithmetic with  $\epsilon =$  $10^{-5}, \epsilon = 10^{-6}$ , and CADNA library are presented in the Table 3. With  $\epsilon = 10^{-7}$  any solution has not been obtained, because an overflow occured during the run of code. The CADNA library detected the numerical instabilities and showed that all the coefficients of GMRES polynomial presented in Table 1 are 0, i.e., they have no significant digit. The test of normalized residuals showed that the solution obtained by CADNA library is a satisfactory informatical solution, but those which obtained by floating-point arithmetic are not. By solving the corresponding linear systems  $A\Delta X = R$  we could improve the solutions obtained with  $\epsilon = 10^{-5}, \epsilon = 10^{-6}$  and obtain the satisfactory informatical solutions which are presented in Table 4. Finally, with many difficulties, we obtained the satisfactory informatical solutions, but what is the accuracy of each element of these solutions? As we observe that the CADNA library not only has obtained a satisfactory informatical solution, but also has furnished the elements of the computed solution with their exact significant digits.

	$\epsilon = 10^{-5}$	$\epsilon = 10^{-6}$	CADNA library
x(1)	1.6700622	1.6700689	0.167001E+01
x(2)	2.9999490	2.9998803	0.299997E+01
x(3)	-4.4399438	-4.4398289	-0.443998E+01
x(4)	3.9999347	3.9998538	0.399999E+01
x(5)	4.1099434	4.1098661	0.41099E+01
	•	:	•
x(146)	-0.9999837	-0.9999635	-0.99999E+00
x(147)	75.1088562	75.1074753	0.75109E+02
x(148)	-0.9999837	-0.9999635	-0.99999E+00
x(149)	76.1087952	76.1073151	0.76109E+02
x(150)	-0.9999837	-0.9999635	-0.99999E+00
	$\nu = 7$	$\nu = 11$	$\nu = 12$
	k = 0	k = 0	k=0

Table 3

	$\epsilon = 10^{-5}$	$\epsilon = 10^{-6}$
x(1)	1.6699998	1.6700000
x(2)	3.00000000	3.00000000
x(3)	-4.4400001	-4.4400010
x(4)	4.0000000	4.0000000
x(5)	4.1100001	4.1100001
	•	b
x(146)	-1.0000000	-1.0000000
x(147)	75.1100006	75.1100006
x(148)	-1.0000000	-1.0000000
x(149)	76.1100006	76.1100006
x(150)	-1.0000000	-1.0000000
	$\nu = 9$	$\nu = 5$
	k=0	k = 0

Table 4

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## 4.2 Using the CADNA library in $A^TA$ -orthogonal s-step Orthomin(k) method

As we have seen in section 2.2, in implementation of  $A^TA$ -orthogonal s-step Orthomin(k) method two problems arise. The first one is:

• How to determine a good value of s?

Let us first consider Tables 5, 7, and 9 of the examples 4-6, respectively. These Tables present the minimum number of significant digits of the norm of orthogonal direction vectors of  $P_0$  (the first s-dimensional subspace) which are furnished by the CADNA library for differents values of s. It emerges from these results that, for each problem, this number begins to decrease from certain s. When it has a small value for some s, the large error exists at the beginning of the iterative process and can lead to the serious round-off errors, and then the slow convergence (see the results of Tables 6, 8, and 10 of the mentioned examples which represent the number of iterations to convergence for different values of s). By noting this remark, it has been observed in experiments that, for double precision, we can obtain a good value of s by taking the highest value of s for which all the orthogonal direction vectors of  $P_0$  have the norm with at least 10 significant digits. By using the CADNA library and increasing the value of s (for example, 4 by 4), it is very easy to determine a such value of s, because the number of significant digits of the norm of orthogonal direction vectors of  $P_0$ , for each s, can be furnished by the CESTAC function which exists in this library, and returns the number of significant digits of every stochastic variable.

Now, we consider the second problem which is:

• How can the iterative process be stopped correctly?

As we mentioned in section 2.2, when we use the stopping criterion

$$||r_i||_2 \le \epsilon, \tag{73}$$

it is possible, due to numerical instabilities or/and the stationarity, this stopping criterion is never satisfied. So, we need to use the additional termination criteria for stopping the process in the cases:

- (i) The algorithm is stationary and can not converge.
- (ii) The computer is not able to distinguish the vector  $r_i$  from the null vector and to improve the computed solution, because of the round-off errors propagation.

As explained in [24, 25], the stochastic arithmetic allows the development of two termination criteria for these cases.

In stochastic arithmetic, when the iterative process becomes stationary, that is, the difference between two iterates is nonsignificant, the components of the vector  $x_i - x_{i-1}$  are stochastic zeros. So, with CADNA library which allows using stochastic arithmetic by automatically implementing the CESTAC method, and using the stopping criterion

$$||x_i - x_{i-1}||_1 = \underline{0}. \tag{74}$$

it is possible to stop the iterative process as soon as it becomes stationary.

In stochastic arithmetic, when the computer is not able to distinguish the vector  $r_i$  from the null vector and to improve the computed solution, because of the round-off errors propagation, the components of  $r_i$  are stochastic zeros and a satisfactory informatical solution is available. So, with CADNA library, and using the stopping criterion

$$||r_i||_2 = \underline{0}. \tag{75}$$

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it is possible to stop the iterative process as soon as the case (ii) occurs and a satisfactory informatical solution is reached.

It is clear that, in the above cases which the iterative process is stopped by the criterion (22) or (23) before the criterion (21) is satisfied, the computed solution will not be a solution with desired accuracy ( $||r_i||_2 \le \epsilon$ ) and it is necessary to improve it by an increment vector  $\Delta x_i$ . For doing this, we need the classical type value of the residual vector  $r_i$  of the computed solution  $x_i$  for solving the linear system  $A\Delta x_i = r_i$  by restarting the iterative process. Fortunately, with CADNA library, it suffices for obtaining the classical type value of  $r_i$  to use the **old\_type** function which exists in this library, and returns the corresponding classical type value of every stochastic variable.

We observe that, with CADNA library, the criteria (22) and (23) stop the iterative process as soon as the cases (i) and (ii) occur, and make it possible to save computation time, because many useless iterations are avoided, to restart the iterative process in order to improve the satisfactory informatical solution which is furnished, and to obtain the solution with the desired accuracy. Consequently, with CADNA library and using the termination criteria (21)-(23), and including the test for restarting the process in the cases in which the process is stopped by the stopping criterion (22) or (23), we can have a stable and efficient  $A^TA$ -orthogonal s-step Orthomin(k) algorithm with the value of s furnished by the method discussed above for solving the linear system and obtaining the desired approximate solution (with  $||r_i||_2 \le \epsilon$ ).

Let us now, to present the examples and the results which we obtained by the FORTRAN code of  $A^TA$ -orthogonal s-step Orthomin(k) method, with floating-point arithmetic for different values of s, and this code with the CADNA library, and the above tests, for the value of

s furnished by the computer. Computation have been performed on a SUN4 computer in double precision. For floating-point arithmetic the stopping criterion was  $||r_i||_2 \le \epsilon$  and the maximum number of iterations allowed set to 1000.

#### Example 4.

We consider the constant-coefficient elliptic equation

$$-\Delta u + 2P_1 u_x + 2P_2 u_y = f, (76)$$

which was described in [11], on the unit square  $\Omega = \{(x,y)|0 \le x, y \le 1\}$  with Dirichlet boundary conditions. Discretizing (24) on  $n_1 \times n_1$  grid gives rise to a sparse linear system of equations of order  $n = n_1^2$ . By using the second order centered differences for the first derivatives and the Laplacian, the coefficient matrix has the form

After scaling the matrix and right-hand side by  $h^2(h = 1/(n_1 + 1))$ , the matrix entries are given by

$$a = 4$$
,  $b = -(1 + p_1)$ ,  $c = -(1 + p_2)$ ,  
 $d = -1 + p_1$ ,  $e = -1 + p_2$ ,

where  $p_1 = P_1h$ ,  $p_2 = P_2h$ . In our test we take  $P_1 = 0$ ,  $P_2 = 50$ . The grid size is h = 1/21, leading to a problem of size 400. The right hand side is determined so that the solution x to the discrete system is 1 everywhere. This allows an easy verification of the results. With  $\epsilon = 10^{-9}$  and  $x_0 = [0, \dots, 0]^T$  the results obtained are presented in Tables 5 and 6. Table 5 contains the minimum number of significant digits of the norm of orthogonal direction vectors of  $P_0$  for different value of s. Table 6 contains the number of iterations needed to satisfy the stopping criterion (21). For CADNA library there are two numbers in this Table, the first one, denoted by TN, presents the total number of iterations needed in the different runs of the iterative process. The second number, denoted by NR, presents the number of restarting of the iterative process.

s	4	8	12	16	20	24	28	32	36	40
MIN	14	14	14	13	11	9	7	5	3	3

Table 5: The minimum number of significant digits of the norm of orthogonal direction vectors of  $P_0$ .

	-	dout	ole f	loat	ing-	poir	ıt aı	ithi	neti	c	CADNA	library
	<del></del>									40		NR
k = 1	35	27	19	11	7	6	7	9	18	22	5	1
k=2	46	23	10	7	6	17	16	19	27	34	5	1
k = 4	52	12	8	8	8	53	57	65	88	90	5	1

Table 6: The number of iterations to convergence

The results presented in Table 5 show that the highest value of s for which all the orthogonal direction vectors of  $P_0$  have the norm with at

least 10 significant digits is s=20. Table 6 shows that the value s=20 is a good value for this example. With computed value s=20, and the CADNA library, the solution was reached with only 5 iterations for all the values k=1,2,4 which is less than those needed with floating-point arithmetic for different values of s. It must be noted that the process was stopped by the stopping criterion (23) at 3th iteration and restarted for improving the computed solution for which the norm of residual was  $||r_3||_2 = 0.484E - 5$ ,  $||r_3||_2 = 0.755E - 6$ ,  $||r_3||_2 = 0.755E - 6$ , for k=1,2,4, respectively. We observe that, for this example, by using CADNA library we could determine a good value of s (s=20), and obtain the desired approximate solution (with  $||r_i||_2 \le 10^{-9}$ ) with the minimum number of iterations. So, for this example, the algorithm using CADNA library is more efficient than that using floating-point arithmetic.

#### Example 5.

We consider the linear system with

$$A = \left[ \begin{array}{cccc} \alpha & 1 & & & \\ -1 & \alpha & 1 & & \\ & \ddots & \ddots & \ddots & \\ & & & 1 & \\ & & & -1 & \alpha \end{array} \right], \quad b = \left[ \begin{array}{c} 1+\alpha \\ \alpha \\ \vdots \\ \alpha \\ \alpha-1 \end{array} \right],$$

which was described in [1], and the dimension equal 400. With  $\alpha = 10^{-8}$ ,  $\epsilon = 10^{-5}$  and the initial vector  $x_0 = [0, \dots, 0]^T$  the results are given in the Tables 7 and 8.

ક		4	8	12	16	20	24	28	32	36	40
M	N	14	14	14	14	14	13	10	7	4	2

Table 7: The minimum number of significant digits of the norm of orthogonal direction vectors of  $P_0$ .

	d	oub	le fl	oati	ng-p	oin	t ari	thm	etic	:	CADNA	library
8	4	8	12	16	20	24	28	32	36	40	TN	NR
k=1	100	50	36	25	20	19	17	15	24	29	18	1
k=2	100	50	37	25	20	20	18	17	23	39	17	1
k=4	100	50	39	25	20	22	20	21	37	58	17	1

Table 8: The number of iterations to convergence

For this example the computed value of s is 28. From the results presented in Table 8 it clearly appears that s = 28 is a good value for this example. With the CADNA library, the solution was reached with 18 iterations for k = 1, and 17 iterations for k = 2, 4. The iterative process was stopped by the stopping citerion (22) at 16th iteration, and restarted in order to improve the computed solution for which the norm of residual was  $||r_{16}||_2 = 0.285E - 3$ ,  $||r_{16}||_2 = 0.163E - 3$ ,  $||r_{16}||_2 =$ 0.124E-3, for k=1,2,4, respectively. We observe that, with CADNA library for k=1 the total number of iterations to convergence is slightly greater than the smallest which corresponds to s = 32, and for k =2,4 this number which is equal to 17 is less than or equal to those needed for different value of s. So, for this example, with CADNA library the program is able to determine a good value of s (s = 28) and to furnish the desired approximate solution (with  $||r_i||_2 \leq 10^{-5}$ ) with a reasonable number of iterations. Consequently,  $A^TA$ -orthogonal s-step Orthomin(k) performed with CADNA library is an efficient tool for solving the linear system of this example.

#### Example 6.

We consider the linear system with

$$A = \begin{bmatrix} 1 & & & & \alpha \\ & 2 & & & \\ & & \ddots & & \\ & & n-1 & \\ & & & n \end{bmatrix}, \quad b = \begin{bmatrix} 1 \\ 1 \\ \vdots \\ 1 \end{bmatrix},$$

which was described in [23, 26] and the dimension equal 100. With  $\alpha = 2 \times 10^6$ ,  $\epsilon = 10^{-10}$  and the initial vector  $x_0 = [0, ..., 0]^T$  the results are listed in the Tables 9, 10.

S	4	8	12	16	20	24	28	32	36	40
MIN	15	11	5	2	2	1	0	0	0	0

Table 9: The minimum number of significant digits of the norm of orthogonal direction vectors of  $P_0$ .

	double fl	oatin	g-po	oint	arithi	netic	CADNA	librai	гy
s	4	8	12	16	20	•,•	TN	NR	
k=1	29	14	10	92	*		13	2	
k=2	27	15	19	85	*		14	3	
k=4	22	74	68	131	*		15	3	

Table 10: The number of iterations to convergence.

\* = problem reached iteration count limit.

For this example, the highest value of s for which all the orthogonal direction vectors of  $P_0$  have the norm with at least 10 significat digits is s = 8. The results of Table 10 show that, with CADNA library, the behavior is similar as in the example 2. It must be noted that the process has been stopped by the stopping criterion (23) two times for k = 1, and three times for k = 2, 4.

## 4.3 Using the CADNA library in the QMR method

As we have seen in section 2.3, in implementation of QMR in the floatingpoint arithmetic, the smallest singular value of matrix  $D_k$  is computed and the criterion

if 
$$(\sigma_{min}(D_k) \leq Tol)$$
 then go to step 3)

is used to check whether this matrix is singular or close to singular, and to decide whether to construct the Lanczos vectors  $v_{n+1}$  and  $w_{n+1}$  as regular or inner vectors. Here Tol is a suitable chosen tolerance. The efficiency of the algorithm depends on a good choice of the Tol and to construct correctly the Lanczos vectors  $v_{n+1}$  and  $w_{n+1}$ , and if the quantity  $\sigma_{min}(D_k)$  is badly computed, propagation of round-off errors will affect drastically all the computation. In addition, It has been shown that [14], in exact arithmetic, the stopping criterion in step 4) of algorithm 2.4 will be satisfied after at most N step -except in a very special situation. If  $\rho_{n+1} = 0$  or  $\xi_{n+1} = 0$  then  $K_n(v_1, A)$  is A-invariant subspace or  $K_n(w_1, A^T)$  is  $A^T$ -invariant subspace, respectively. In the first case  $x_n$  is the exact solution and the QMR algorithm must be stopped. In the second case, by restarting the QMR method and using the last available QMR iterate  $x_{n-1}$  (which is a good choice [14]) as the new initial guess, it is possible to improve the approximate solution. So, in floating-point arithmetic, we have to use the following criterions for stopping or restarting the QMR algorithm,

if 
$$(\rho_{n+1} \le \epsilon_1)$$
 then stop, (77)

if 
$$(\xi_{n+1} \le \epsilon_1)$$
 then restart, (78)

where  $\epsilon_1$  is a suitable chosen tolerance. It is necessary to mention that, it is possible, due to round-off errors propagation,  $\rho_{n+1}$  becomes non-significant and in the same time it has a large value  $(\rho_{n+1} > \epsilon_1)$ . In this

situation, it is clear that, the iterations of QMR method are not able to improve the approximate solution and the QMR method must be restarted with the last available QMR iterate  $x_{n-1}$ . Finally, for QMR Algorithm a convergence criterion is needed. We can stop the QMR algorithm as soon as the criterion

$$||r_n||_2 \le \epsilon_2 ||r_0||_2,\tag{79}$$

is satisfied, where  $\epsilon_2$  is also a suitable chosen tolerance. As we observe another problem is how to choose  $\epsilon_1$  and  $\epsilon_2$ . We now face the following questions:

- How can we detect the informatical singularity of the matrix  $D_k$  and the numerical instabilities which may occur in the program?
- How can the iterative process be stopped or restarted correctly?

In order to overcome these drawbacks, we propose to introduce the stochastic arithmetic in the QMR algorithm, which is able to estimate the round-off error propagation, and to detect the informatical singularity of the matrix  $D_k$  by means the following simple test,

if 
$$(\sigma_{min}(D_k) = \underline{0})$$
 then go to step 3) (80)

and to restart or to stop the process by the following tests,

if 
$$(\rho_{n+1} \le \epsilon_1 \text{ or } C_{\rho_{n+1}} < 1)$$
 then (if  $||r_n|| = \underline{0}$ ) then stop else restart, (81)

if 
$$(\xi_{n+1} \le \epsilon_1 \text{ or } C_{\xi_{n+1}} < 1)$$
 then (if  $||r_n|| = \underline{0}$ ) then stop else restart, (82)

if 
$$(||r_n||_2 = \underline{0})$$
 then stop. (83)

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It has been observed in experiment that, for double precision,  $\epsilon_1 = 10^{-8}$  is a good choice for tests (29)-(30) and we will have a stable algorithm.

Let us now, to present an example and the results which we obtained by the FORTRAN code of the QMR algorithm, with floating-point arithmetic for  $Tol = 10^{-4}$  and difference value of  $\epsilon_2$ ; and this code with stochastic arithmetic and the test (28) and the termination criteria (29)-(31). Computation have been performed in double precision and the number of iterations allowed set to 200000.

**Example 7.** We consider the ill-conditioned linear system Hx = bwith the Hilbert matrix  $H(h_{ij} = 1/(i+j-1))$ , and dimension equal to 100. The right hand side is determined so that the exact solution x is 1 everywhere. This allows an easy verification of the results. With  $x_0$  =  $[0,\ 0,\cdots,\ 0]^T,\,\epsilon=10^{-7},\,\epsilon=10^{-8},$  and stochastic arithmetic the results obtained are presented in Table 11. These results show that the results obtained with  $\epsilon = 10^{-8}$  have not more significant digits than those obtained with  $\epsilon = 10^{-7}$ , but the iterative process is stopped too late (at iteration 138312). So many useless iterations are performed without improving the accuracy of the solution. It is necessary to mention that with  $\epsilon = 10^{-9}$  the iterative process has not been converged. As we observe, in stochastic arithmetic, the better solution is furnished with only 6841 iterations which is very smaller than those of floating-point arithmetic with  $\epsilon = 10^{-7}$  and  $\epsilon = 10^{-8}$ . So, we can conclude that the use of criterion (28) which detect the informatical singularity of matrix  $D_k$  allows to stabilize the algorithm, and the use of the optimal criteria (29)-(31) allow to continue the iterations, to restart QMR method if it is necessary, and to stop the program as soon as a satisfactory solution is reached.

	$\epsilon = 10^{-7}$	$\epsilon = 10^{-8}$	Stochastic
			arithmetic
x(1)	0.9999542	1.0000064	1.0000
x(2)	1.0007666	0.9998362	1.00
x(3)	0.9976219	1.0008133	0.999
•	•		•
x(98)	0.9968086	0.9986009	0.995
x(99)	0.9963570	0.9983446	0.995
x(100)	0.9958868	0.9980741	0.994
iter	19579	138312	6841

Table 11

#### 5. Conclusion

In this paper we have seen in the implementation of iterative methods the following problems arise

- How can the iterative process be stopped correctly?
- What is the accuracy of the computed solution given by computer?
- How can detect the numerial instabilities which may occur in the program?

We observed that the use of CADNA library allows us to solve these problems. It has been shown that the CADNA library with the optimal termination criterion and the appropriate test is able to stop the program as soon as a satisfactory solution is reached, to estimate the accuracy of the solution, to detect the numerical instabilities, to prevent an overflow which may occur, and to save computer time, because many useless operations and iterations are not performed. Consequently, the mentioned iterative methods with CADNA library are the robust and efficient tools for solving large nonsymmetric systems of linear equations.

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Proceedings of the Monthly Colloquium of the Iranian Mathematical Society Vol. 2, pp 271-281 (2004)

## Local Defining Ideals of Ordinary Singularities

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Dedicated to the late Professor Karim Seddighi

Abstract: We review some basics on the theory of generic singularities in algebraic geometry. We state certain approaches towards the explicit local equations of such singularities. The main point of view is the constructive approach in commutative algebra and algebraic geometry.

#### 1. Introduction

One of the goals in algebraic geometry is the "local classification" of algebraic varieties. Namely, if X is an algebraic variety over an algebraically closed field k and  $\mathcal{O}_{X,x}$  is the local ring of regular functions at x (i.e., the ring of "germs" of rational functions on a neighborhood

MSC (2000): 14B05, 14B25

Keywords: Generic singularities, Gröbner basis, local defining ideal.

of x), how can one classify these rings up to isomorphism? This is a far-reaching problem. The "local analytic classification" is the classification of  $\widehat{\mathcal{O}}_{X,x}$ , the completion of  $\mathcal{O}_{X,x}$ . The Cohen structure theorem states that if X is smooth at x, then  $\widehat{\mathcal{O}}_{X,x} = k[[x_1, \cdots, x_r]]$ , the ring of formal power series in r variables where  $r = \dim X$ . Such a classification for  $\mathcal{O}_{X,x}$  with X smooth at x, does not exist.

The theory of ordinary singularities is an approach toward the classification of  $\widehat{\mathcal{O}}_{X,x}$  when X is singular at x. Ordinary singularities arise from "generic" projection of smooth varieties. More precisely, let  $Y \subset \mathbb{P}^n$  be a projective smooth variety of dimension r in the n-dimensional projective space. Let E be a point outside Y. Let  $\mathbb{P}^{n-1} \subset \mathbb{P}^n$  be the projective space of dimension n-1 and let  $\pi:Y \longrightarrow \mathbb{P}^{n-1}$  be the projection from E. If E is in general position, then  $Y_1 = \pi(Y)$  has "nice" singularities. We may repeat this procedure of projecting for  $Y_1$ , and continue as long as the projection is a birational map onto its image. The singularities of the resulting variety X are called "ordinary singularities". A "node" is the only ordinary singularity of curves, and in this case

$$\widehat{\mathcal{O}}_{X,x} = k[[x_1, x_2]]/(x_1x_2).$$

The ordinary singularities of surfaces are also fairly simple. They consist of ordinary double points, normal (triple) crossings, and pinch points, where the completion of the local rings are isomorphic to

$$k[[x, y, z]]/(xy), k[[x, y, z]]/(xyz), k[[x, y, z]]/(x^2z - y^2),$$

respectively. However, the ordinary singularities of higher dimensional varieties are more complicated. Intuitively, ordinary singularities are "stable", and hence they are appropriate models for approximations of critical points of natural phenomena.

Many authors such as K. Mount, O. Villamayor, J. Roberts, J. Lluis and A. Holme have given substantial contributions to the subject. A reasonably complete "parametric" representation of ordinary singularities is now available. However, to apply results from "the ideal theory of the ring of polynomials over a field", one needs the explicit equations of these ordinary singularities. For ordinary "double points", this is rather simple. Indeed, if x is an ordinary double point, then

$$\widehat{\mathcal{O}}_{X,x} = k[[x_1, \cdots, x_p]]/(x_1, \cdots, x_p)(x_{p+1}, \cdots, x_{2p}),$$

or,

$$\widehat{\mathcal{O}}_{X,x}=k[[z,z_1,\cdots,z_n,u_1,\cdots,u_n]]/I,$$

where,

$$I = \text{the ideal of } 2 \times 2 \text{-minors of} \begin{pmatrix} z_1 & u_1 & z_2 & u_2 & \dots & z_n & u_n \\ zu_1 & z_1 & zu_2 & z_2 & \dots & zu_n & z_n \end{pmatrix},$$
(See [Z] and [ZZ]).

In the first case, the normalization is the product of two complete regular local rings generalizing a node, while in the second case, the normalization is itself a complete regular local ring generalizing a pinch point.

Using the above classification of the ordinary double points, the quotient forms for the classification of triple points follow from [R1] and [SZ]. Here the goal is to see some picture of the progress on the quotient forms of the completions of local rings at ordinary singularities of higher "multiplicities". We will omit the proofs.

## 2. Ordinary Singularities with Higher Multiplicities

For some time, in collaboration with P. Salmon, we have been trying to provide a nice presentation for ordinary singularities of higher multiplicities as a quotient of the polynomial ring by an explicit ideal. Let A be the completion of the local ring at a point x on a variety X. Let B be the normalization of A. It is well-known that B is a finite A-module (e.g., see[S]). Let q be the length of B as an A-module. The positive integer q is called the multiplicity of X at x or simply the multiplicity of x. Roughly speaking, a point of multiplicity q is a point at which q "simple branches" of the variety intersect, or, it is a "limiting position" of such points. If B is a domain, the point x is called a "unibranched" point of multiplicity q. Using these notions, we concentrate on unibranched ordinary singularities of multiplicity q. The parametric equations of these points are given in [R2] as:

$$z = yt + xt^2 + \dots + vt^{q-2} + t^q,$$

$$z_i = y_i t + x_i t^2 + \dots + u_i t^{q-1}; i = 1, \dots, m.$$

In order to obtain the local defining ideal of the variety at such a point, we need all relations obtained by "eliminating" t in the above equations. Consider the "resultant matrix":

$$N = \begin{bmatrix} N_0 \\ N_1 \\ \vdots \\ N_n \end{bmatrix},$$

where.

is a  $(q-1) \times (2q-1)$  matrix, and

is a  $q \times (2q-1)$  matrix.

Conjecture 1. (January 1989) The local defining ideal of the unibranched ordinary singularity with multiplicity q is generated by the maximal minors of N.

**Lemma 1.** To prove Conjecture 1, it is enough to prove it for the case  $y = x = \cdots v = 0$ . In this case the maximal minors of N are the same as the maximal minors of

$$M = egin{bmatrix} M_1 \\ M_2 \\ \vdots \\ M_m \end{bmatrix}$$

where

$$M_i = egin{bmatrix} z_i & zu_i & zv_i & \ldots & zy_i \ y_i & z_i & zu_i & \ldots & zxi \ \vdots & \ddots & \ddots & \ddots & \ddots \ \vdots & \ddots & \ddots & \ddots & \ddots & \ddots \ u_i & v_i & \ldots & y_i & z_i \end{bmatrix}$$

for  $i = 1, \dots, m$  (we are using  $z_i$  for  $-z_i$  for simplicity).

Remark 1. The concept of resultant matrix first appeared in [SZ]. This generalizes the usual Sylvester resultant of two polynomials in one variable and is in general a larger ideal compared to the classical "uresultant" (e.g., see [W1]), and hence is more useful from the scheme theoretical point of view.

Remark 2. The set of maximal minors of M is not a minimal generating set for this ideal. Indeed, the number of maximal minors of M is  $\binom{mq}{q}$ , which is by far larger than  $m^q$ , the expected number for a minimal generating set.

Remark 3. Putting z = 1, the matrix  $M_i$  becomes a circulant matrix (see [D]). In the sense that each row is a permutation of the first row by a cycle of length q. Assigning certain degrees to the variables as in the next section,  $M_i$  will be the homogenization of the this circulant matrix.

## 3. An Approach via Gröbner Algebra.

Gröbner bases are certain tools for computations in the ring of polynomials. To distinguish the "initial monomial" of a polynomial as the largest monomial appearing in the polynomial, a total order is given on

the monomials. A generating set of an ideal whose initials generate the ideal of initials of the given ideal, is called a Gröbner basis of the ideal. The initials of a Grobner basis is a set of monomials which can best reflect the given ideal.

Let's reorder the matrix M as

$$M = egin{bmatrix} z_{m} & zu_{m} & zv_{m} & \ldots & zy_{m} \\ \vdots & \vdots & \ddots & \ddots & \ddots & \vdots \\ \vdots & \vdots & \ddots & \ddots & \ddots & \ddots & \vdots \\ z_{1} & zu_{1} & zv_{1} & \ldots & zy_{1} \\ y_{m} & z_{m} & zu_{m} & \ldots & zx_{m} \\ \vdots & \ddots & \ddots & \ddots & \ddots & \vdots \\ \vdots & \ddots & \ddots & \ddots & \ddots & \vdots \\ \vdots & \ddots & \ddots & \ddots & \ddots & \ddots & \vdots \\ u_{m} & v_{m} & \ldots & v_{m} & z_{m} \\ \vdots & \ddots & \ddots & \ddots & \ddots & \ddots & \vdots \\ u_{m} & v_{m} & \ldots & \ddots & \ddots & \ddots & \vdots \\ \vdots & \ddots & \ddots & \ddots & \ddots & \ddots & \vdots \\ u_{1} & v_{1} & \ldots & v_{1} & z_{1} \\ \end{bmatrix}$$

Let  $\deg u_i = 1$ ,  $\deg v_i = 2$ ,  $\cdots$ ,  $\deg y_i = q-1$ ,  $\deg z_i = \deg z = q$ . Consider the "degree reverse lexicographic" order on the monomials induced by

$$z_m > z_{m-1} > \cdots > z_1 > z > y_m > \cdots > y_1 > x_m > \cdots > x_1 > u_m > \cdots > u_1.$$

Definition 1. A maximal minor of M is called principal if no entry on

its main diagonal is a multiple of z.

**Lemma 2.** The ideal of maximal minors of M is generated by the set of principal maximal minors of M.

Lemma 3. The initial monomial of a principal maximal minor of M with respect o the above monomila order, is the product of the entries on the main diagonal of the minor.

Conjecture 2. (August 1997) The set of principal minors of M forms a Gröbner basis for the ideal of maximal minors of M.

Theorem 1. Conjecture 2 implies Conjecture 1.

**Definition 2.** A principal minor is called a distinguished principal minor (DP), if the entries on the main diagonal are non-decreasing from the top to the bottom.

Conjecture 3. The set of DP minors of M generates the ideal of maximal minors of M.

Theorem 2. Conjecture 3 implies Conjecture 2.

Lemma 4. The number of DP minors of M is  $m^q$ .

Conjecture 3 has lead to certain basic determinantal identities which do not seem to be well-known. Let's state just one of such identities.

Theorem 3. Let

$$D = \begin{bmatrix} L_1 \\ L_2 \\ \cdot \\ \cdot \\ \cdot \\ L_n \end{bmatrix}$$

be a  $n \times n$  matrix,  $\theta \in S_n$  any permutation with r fixed points. Then

#### 4. The "episode" of Elimination Theory!

The concept of resultant has been a central one in the elimination theory heavily studied in the 18th and the 19th centuries mostly with constructive methods. In the early part of the 20th century, many of the leading mathematicians in algebraic geometry and commutative algebra. became less interested in elimination theory and its classical methods. A. Weil in his influential book "Foundation of Algebraic Geometry, 1946" ([W], page 31), right before quoting an important result from C. Chevalley, says, "The device that follows · · · it may be hoped, finally eliminates from algebraic geometry the last traces of Elimination Theory ...". A section of the famous book by B.L. van der Wearden, "Modern Algebra 1953" [W1], is devoted to elimination theory. In the later editions (1959-) [W2], the name "Algebra" is adopted for the book, and van der Wearden eliminates the section on elimination theory! As a result, a simple constructive proof of the existence of a resultant system essentially due to Kronecker, is replaced by a non-constructive proof using the nontrivial Hilbert's Nullstellensatz in the second volume of the book. In the seventies of this century, the tide starts to turn. This is reflected in S. Abhyankar's famous poem of 1970's:

Eliminate, eliminate, eliminate,

Eliminate the eliminators of elimination theory

Today, with the enormous influence of computers in mathematics, the use of constructive methods has been inevitable [E]. In commutative algebra and algebraic geometry, the basic notion for such constructive methods, has been the notion of "Gröbner basis". We have been trying to put some of the constructive methods of elimination theory in the frame work of the theory of Gröbner bases.

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زندگی علمی و اجتماعی انجمن ریاضی ایران از همایش ریاضیدانان در سال ۱۳۴۹ آغاز شده و رشد و تکامل آن با تشکیل و ادامهٔ همایشهای گوناگون دمساز بوده است: کنفرانس سالانهٔ ریاضی ایران که سی و چهارمین آن شهر بورماه ۱۳۸۲ در شاهرود برگزار شد، سمینار هفتگی انجمن که چند سال نخست دایر بود و سمینارهای هفتگی دانشگاهها جای آن را گرفت، چندین سمینار تخصصی که با سمینار جبر اهواز در دیماه ۱۳۵۶ آغاز شد و هر از چند گاهی یک سمینار تخصصی نو ظاهر گشت. بدون شک با مشارکت در همایشها امکان گفت و شنود و تبادل نظر در زمینههای مختلف فراهم و در همین اثنا بذرهایی کاشته می شوند که امید می رود در آینده درختان تنومندی در زمینهٔ اندیشهٔ ریاضی به بار آورند. از این نظر می توان صرف نیروی انسانی و منابع مالی در راه برگزاری همایشها را توجیه کرد. اما اگر این افکار مدوّن و منتشر نشوند، دامنهٔ برد آنها از حیث زمان و مکان بسیار محدود خواهند ماند. متأسفانه راهی برای تدوین و نشر نشوند، دامنهٔ برد آنها از حیث زمان و مکان بسیار محدود خواهند ماند. متأسفانه راهی برای تدوین و نشر بازتاب یابند. اما میاحث اصلی کنفرانس متصور نیست مگر قسمتهای بسیار اندکی که ممکن است در خاطرات بازتاب یابند. اما میاحث اصلی کنفرانس یعنی مقالات ارائه شده را نه تنها می توان، بلکه باید به هر قیمت و رحمتی بخش کرد.

به طور کلی انتشارات ادواری و غیر ادواری انجمنهای علمی نیز جزء ارکان مهم فعالیتهای آنها هستند. انجمن ریاضی ایران خوشبختانه علاوه بر گزارش کنفرانسهای سالانه و سمینارهای تخصصی، چندین نشریهٔ ادواری دارد که هر یک دارای منشور و رسالت ویژهٔ خود است: بولتن انجمن ریاضی ایران ویژهٔ مقالات پژوهشی در سطح بینالمللی به زبان انگلیسی، فرهنگ و اندیشهٔ ریاضی ویژهٔ مقالات مروری به زبان فارسی، خبرنامه وگزارش ویژهٔ اطلاع رسانی و اظهار نظرها.

از سال ۱۳۷۸ همایشهای ماهانهٔ انجمن ریاضی ایران تشکیل گردید که هدف آن آشنا ساختن پژوهشگران جوان با رشتههای پژوهشی گوناگون است. شیوهٔ کار این همایش آن است که در کمیتهٔ همایش نسبت به دعوت از استادان مقیم داخل و خارج که در زمینهای پیشرفت قابل ملاحظهای داشته و کارهای پژوهشی را هدایت نموده باشند دعوت به عمل آید که مقالهای مروری ارائه دهند و آن را به زبان فارسی با انگلیسی در اختیار انجمن قرار دهند تا به شکل شایسته منتشر شود. کمیتهٔ اول در سال ۱۳۷۸ عمدتاً در شهر کتاب هفت همایش جالب برگزار کرد. کمیتهٔ دوم در سال ۱۳۷۹ همایش را در دانشگاههای مختلف،

در سال ۱۳۸۰ آن را در تالار فرهنگسرای دانشجو و دانشگاه تهران، و در سال ۱۳۸۱ همه را در دانشگاه تهران برگزار نمود.

از آنجا که ویرایش مقالات رسیدهٔ سال اول بیش از حد طول کشید و آخر سر هم امر ویرایش ناقص ماند، کمیته تصمیم گرفت در مسألهٔ ویرایش تجدید نظر کند و به اتکای آن که مقالات را استادان زبردست می نویسند، از ویرایش علمی و ادبی آنها به کلی صرف نظر نماید. بدین ترتیب، نویسندگانِ مقالات نهایت جدیت را مبذول فرمودند که مقاله به شکل نهایی و آمادهٔ تکثیر باشد. به علاوه، کمال مطلوب آن است که مقاله چندین روز پیش از برگزاری همایش به دست کمیته برسد تا به تعداد اندکی تکثیر شود و در اختیار شرکت کنندگان قرار گیرد. اگر این کمال مطلوب حاصل شود، در پایان هر سال مجلدی مشتمل بر مقالاتِ همان سال برای چاپ و انتشار آماده می شود و در اختیار عموم قرار می گیرد.

جلد اول هشت مقاله را در برگرفت که مشتمل بر مقالات فارسی دریافت شده بود. جلد دوم، یازده مقالهٔ رسیده به زبان انگلیسی را در بر میگیرد که آن نیز حاوی مقالاتی از سال ۱۳۷۸ تا بهار ۱۳۸۱ است. برخی از مقالات ارائه شده در طول ۴ سال گذشته به دست کمیته نرسیده است.

از طرف خود و همکارانم در کمیتهٔ همایش، از استادان بزرگواری که دعوت ما را پذیرفتند و به نگارش و تحویل مقالهٔ خود پرداختند متواضعانه سپاسگزارم. همچنین لازم می دانم از حمایتهای جناب آقای دکتر مهدی بهزاد رئیس سابق و جناب آقای دکتر سیدعبادالله محمودیان رئیس فعلی انجمن ریاضی ایران و همچنین از اعضای محترم شورای اجرایی و همکارانم در کمیتهٔ همایش ماهانهٔ انجمن خانمها دکتر: شیوا زمانی و مرگان محمودی و آقایان دکتر: علی آبکار، غلامحسین اسلامزاده، علی ایرانمنش، حسین حاج ابوالحسن و سیامک یاسمی سپاسگزاری کنم.

اجرای مراحل مختلف همایش مدیون زحمات اعضای فقال دبیرخانهٔ انجمن است. بدین وسیله از همهٔ این عزیزان به ویژه از تلاشهای خانم افسانه بختیاری و آقای مرتضی عبدیزاده که ظرف چند سال گذشته در برگزاری همایشها کمیته را یاری نمودند و خانم فریده صمدیان به خاطر حروف چینی بعضی مقالات و پردازش نهایی متن و یکنواختی آنها تشکر میکنم. سازماندهی و نظم کارهای دبیرخانه از زمان تصدی آقای منصور شکوهی به ریاست دبیرخانه در تحقق این انتشار مفید و مؤثر بوده است. از حسن ادارهٔ ایشان قدرشناسی میکنم.

در خاتمه، پیروزی اعضای جدید کمیتهٔ همایش ماهانه آقای دکتر علی آبکار، خانم دکتر لیلا خاتمی، آقای دکتر مهدی دهقان (رئیس کمیته)، آقای دکتر عمید رسولیان، آقای دکتر بهروز طایعه رضایی و خانم دکتر مژگان محمودی را خواستارم.

ارسلان شادمان تهران، بهار ۱۳۸۳

## بسمالله الرحمن الرحيم



## گزارش همایش ماهانه انجمن ریاضی ایران

جلد دوم، بهار ۸۳

اعضای کمیتهٔ همایش ماهانه:

على أبكار (٨١\_١٣٨٠)

غلامحسین اسلامزاده (۸۱\_۱۳۷۹)

مسعود امینی (۱۳۷۸)

على ايرانمنش (٨١\_١٣٧٨)

حسين حاج ابوالحسن (١٣٧٨)

آرش رستگار (۱۳۷۸)

شیوا زمانی (۱۳۷۸)

ارسلان شادمان (۸۱\_۱۳۷۹)

مژگان محمودی (۱۳۷۹\_۱۳۷۹)

سیامک یاسمی (۱۳۷۸)

حروفچینی: IATEX \_ فریده صمدیان

چاپ و صحافى: چاپ بهمن ـ قم

ناشر: بهمن آرا

شابک: ۸\_۳\_۹۵۲۶۸ ۹۶۴

نشانی: تهران، صندوق یستی ۴۱۸\_۱۳۱۴

انجمن رياضي ايران

تلفن و نمابر: ۵۸۸۵،۸۸۵، ۸۹۷۷۰۸، ۵۷۷۷ م۸۸

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نشاني الكترونيك:

حق چاپ محفوظ و مخصوص انجمن ریاضی ایران است.

همایش ماهانهٔ انجمن ریاضی ایران بهعنوان یکی از فعالیتهای مستمر انجمن از ۱۳۷۸ با اهداف زیر شروع شده است:

نقل از آیین نامهٔ مصوب ۲۸/۱/۲۶

 آشنایی جامعه ریاضی ایران با شاخههای مختلف ریاضیات.

آشنایی دانشجویان تحصیلات تکمیلی علوم ریاضی
 با مسائل و زمینههای مختلف تحقیقاتی ریاضی در ایران.

 ایجاد ارتباط بین محققان و دانشجویان ریاضی کشور،
 بهویژه دانشجویان تحصیلات تکمیلی رشته های مختلف علوم ریاضی.

• كمك به شناخت توان تحقيقاتي كشور در رياضيات.

 کمک به ایجاد جو مساعد برای رشد ریاضیات در ایران و باروری استعدادهای ریاضی.

مقالههای همایش معمولاً مروری تحقیقی هستند و به معرفی یک رشتهٔ فعال پژوهشی میپردازند. این مقالات توسط یکی از استادان مدعو داخل یا مقیم خارج ارائه میشوند. جهت تسریع در انتشار مقالات، کمیتهٔ همایش از ویرایش آنها معذور است. از این رو، درخواست می شود نویسنده مقاله را به شکل نهایی تحویل دهد. چاپ مقالهها در مجلات علمی ترویجی به همان شکل یا با اندکی تغییر بلامانع است. در این صورت توصیه می شود نویسندهٔ محترم به جلد و صفحات گزارش همایش ماهانه که مقاله در آن مندرج است اشاره فرماید.

از هر مقاله ۲۵ نسخه به مؤلف تقدیم می شود. زمان برگزاری همایش معمولاً آخرین چهارشنبه هر ماه جز مرداد و اسفند و با رعایت ماه مبارک رمضان است.

# Proceedings of the Monthly Colloquium of the Iranian Mathematical Society

Volume 2 (2004)

Edited by: Arsalan Chademan



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