MATH 578 Numerical Analysis, Fall 2020
Student Notes

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Abstract

This is just my notes going through the highlights from the lecture notes for MATH 578 Numerical Analysis, Fall 2020 (Tsogtgerel, 2020). As a computational statistician muggle taking this course for optimization and machine learning, this notes might not always be a good summary, I must say... As an overview, the first chapter covers up some basics of analysis functions – specifically, Lagrange interpolation (Taylor series expansion is a special case) and minimax polynomials; the second chapter covers up linear system part.

I Function Evaluation

1 Basic Computer Arithmetic $\forall a \in \mathbb{Z}$, a *base−β representation* exists for some $\beta \in \mathbb{N} \setminus \{1\}$:

$$a = \pm \sum_{k=0}^{\infty} a_k \beta^k$$

where $0 \leq a_k \leq \beta - 1$ is defined as the k−th digit of a in base β. And grade-school column sum/difference first carries out *Cauchy sum* or *difference*, which takes sum/difference for each digits; then it recursively perform carrying for addition for borrowing for subtraction. Let

$$n := \max\{k | a_k \neq 0\}, \ m := \max\{k | b_k \neq 0\}$$

So a, b will be $n+1$ and $m+1$ digit number. The bit complexity for addition/subtraction will then be $O(n+m+1)$. Column multiplication carries out similarly. However, multi-

plication can also be done row-wisely: the *Cauchy product*

$$ab = \left(\sum_{i=0}^{\infty} a_i \beta^i\right) \cdot b = \sum_{i=0}^{n} a_i \cdot \beta^i b$$

where $\beta^j b$ is simply shifting digits, and multiplication by a_i can be carried out as column addition. The bit complexity for column multiplication would then be $O(nm+1)$.

As for division algorithm, assume that the quotient is expressed as:

$$q = q_0 + q_{-1}\beta^{-1} + q_{-2}\beta^{-2} + \cdots$$

And let a, b here be positive and normalized. The *partial reminder* refers to the *normalized* reminder obtained in the division process. Two division algorithms for a/b were introduced here: i). *restoring division*: keeping performing subtraction see if the partial reminder goes below 0, and if it goes below 0, "restore" by adding the divisor back to it to prevent negative digits; ii). *non-restoring division*: the idea of non-restoring division is to use generalized digit, e.g. $\{-1, 1\}$ for binary computing, to allow negative sign in a digit, and a conversion back to standard digit will be indeed required *in the end*. To generalize non-restoring division to any radix β, note that the partial reminders are given by:

$$r_{j+1} = \beta r_j - q_{-j} b$$

the above two division processes determine q_{-j} both by subtracting b from βr_j, the difference is for restoring division, $0 \le q_{-j} < \beta$ gives partial reminder $0 \le r_{j+1} < b$; for non-restoring division, $-\beta < q_j < \beta$ gives partial reminder $-b \le r_{j+1} < b$.

However, both of above division algorithms are not efficient – especially not for bignums. WLOG, let a, b be integers here, the idea of *long division* is to determine the quotient by observing the first digit of the divisor and perform restoring division. In comparison, *SRT division* is non-restoring division with normalized divisor and reminder. *Error propagation* describes the idea of computation will alternate (mostly increase) the error of approximation numbers, such as floating point numbers. Usually error propagation is captured upper-boundedly by *conditional number*, e.g. conditional number of summation is

$$\kappa_+(x) = \frac{|x_1| + |x_2| + \cdots + |x_n|}{|x_1 + x_2 + \cdots + x_n|}$$

Furthermore, the following axiom is used for a wide-range of numerical error analysis for floating point numbers: For each $\star \in \{+, -, \times, /\}$, there exists a binary operation $\circledast : \tilde{\mathbb{R}} \times \tilde{\mathbb{R}} \mapsto \tilde{\mathbb{R}}$ s.t.

$$\left|x \star y - x \circledast y\right| \le \varepsilon \left|x \star y\right|, \ x, y \in \tilde{\mathbb{R}}$$

dividing by zero is excluded. Normally, ε is referred as "machine precision."

2 Evaluation of Power Series A function $f : (a,b) \mapsto \mathbb{R}$ is called *analytic* at $c \in (a,b)$ if it can be developable into a power series around c; and called analytic at (a,b) if analytic at c, $\forall c \in (a,b)$. For such class of analytic functions, a way to evaluate them is through Taylor series, backed by a generalized version of mean value theorem proposed by Lagrange: Let f be a $n+1$ times differentiable function in (c,x), with $f^{(n)}$ continuous in $[c,x)$. Then $\exists \xi \in (c,x)$ s.t.

$$f(x) = \sum_{k=0}^{n} \frac{f^{(k)}(c)}{k!}(x-c)^k + \frac{f^{(n+1)}(\xi)}{(n+1)!}(x-c)^{n+1}$$

See Lagrange interpolation section coming later for proof. This theorem gives an expression of the error as a result of approximating using n–th order Taylor series. Moreover, the following series are listed with their relative condition numbers:

$$\frac{1}{1-x} = \sum_{k=0}^{\infty} x^k, \ \forall |x| < 1$$

$$\kappa(x) = \left| \frac{(1-x)^{-2}}{(1-x)^{-1}/x} \right| = \left| \frac{x}{1-x} \right|$$

$$e^x = 1 + x + \frac{x^2}{2} + \cdots + \frac{x^n}{n!} + \cdots, \ \forall x \in \mathbb{R}$$

$$\kappa(x) = \left| \frac{(e^x)'}{e^x/x} \right| = |x|$$

$$\log(1+x) = \sum_{k=1}^{\infty} \frac{(-1)^{n-1} x^n}{k}, \ -1 < x \leq 1$$

$$\kappa(x) = \left| \frac{(1+x)^{-1}}{\log(1+x)/x} \right| = \frac{x}{(1+x)} \cdot \left| \frac{1}{\log(1+x)} \right|$$

$$\sin x = \sum_{k=0}^{\infty} (-1)^k \frac{x^{2k+1}}{(2k+1)!}, \ \forall x \in \mathbb{R}$$

$$\kappa(x) = \left| \frac{\cos x}{\sin x/x} \right| = |x \cot x|$$

$$\cos x = \sum_{k=0}^{\infty} (-1)^k \frac{x^{2k}}{(2k)!}, \ \forall x \in \mathbb{R}$$

$$\kappa(x) = \left| \frac{-\sin x}{\cos x/x} \right| = |x \tan x|$$

$$\arctan x = \sum_{n=0}^{\infty} \frac{(-1)^n x^{2n+1}}{2n+1}, \ |x| \leq 1$$

$$\arcsin x = x + \frac{1}{2}\cdot\frac{x^3}{3} + \frac{1\cdot 3}{2\cdot 4}\cdot\frac{x^5}{5} + \frac{1\cdot 3\cdot 5}{2\cdot 4\cdot 6}\cdot\frac{x^7}{7} + \cdots,\ -1 \le x < 1$$

And recall that the relative condition numbers is defined by:

$$\kappa := \lim_{\varepsilon\downarrow 0}\ \sup_{\|\delta x\|\le\varepsilon} \frac{\|\delta f\|/\|f(x)\|}{\|\delta x\|/\|x\|}$$

3 Acceleration of Convergence Two methods of acceleration of convergence are discussed here:

i). *Euler transform*: *Hausdorff moment characterization* says that

$$m_k = \int_0^1 x^k d\mu \text{ for some } \sigma - \text{additive Borel probability measure } \mu$$
$$\Leftrightarrow m_0 = 1,\ m \text{ is completely monotone; i.e. } (-1)^n \Delta^n m_k \ge 0,\ \forall n, k$$

The formula

$$\frac{\pi}{4} = \sum_{k=0}^{n} (-1)^k \frac{1}{2k+1}$$

can be accelerated by repeatedly taking average of two consecutive terms, called *Euler transform*. Applying Hausdorff moment characterization, error analysis for this can be done by noticing that

$$a_k = \frac{1}{k} = \int_0^1 t^k d\mu$$

and the rest follows from power series.

ii). *Aitken's Δ^2–process*: used to evaluate a noisy geometric series. For a series defined by

$$a_k = Cq^k + O\left(\delta^k\right), \text{ for some } 0 < \delta < q < 1$$
$$S_n = \sum_{k=1}^{n} a_k$$

Observe that

$$\begin{aligned} S &= S_n + \sum_{k=n+1}^{\infty} a_k \\ &= S_n + \sum_{k=n+1}^{\infty} Cq^k + O(\delta^n) \end{aligned}$$

$$
\begin{aligned}
&= S_n + \frac{Cq^{n+1}}{1-q} + O(\delta^n) \\
&= S_n + \frac{a_n^2}{a_{n-1} - a_n} + O(\delta^n)
\end{aligned}
$$

The last inequality above used the fact that

$$
\begin{aligned}
q &= \frac{a_n}{a_{n-1}} + O\left(\left(\frac{\delta}{q}\right)^n\right) \\
a_n &= Cq^n + O(\delta^n) \\
\Rightarrow Cq^{n+1} &= \left(\frac{a_n}{a_{n-1}} + O\left(\left(\frac{\delta}{q}\right)^n\right)\right)(a_n - O(\delta^n)) = \frac{a_n^2}{a_{n-1}} + O(\delta^n), \text{ and} \\
\frac{1}{1-q} &= \frac{a_{n-1}}{a_{n-1} - a_n} + O\left(\left(\frac{\delta}{q}\right)^n\right)
\end{aligned}
$$

Let

$$
\begin{aligned}
\Delta S_{n-1} &:= S_n - S_{n-1} = a_n \\
\Delta^2 S_{n-2} &:= a_n - a_{n-1} = \Delta a_{n-1}
\end{aligned}
$$

We then have

$$
S_n + \frac{a_n^2}{a_{n-1} - a_n} = S_n - \frac{(\Delta S_{n-1})^2}{\Delta^2 S_{n-2}}
$$

which gives the name "Δ^2"

4 Root Finding Fixed point iterations are based on a theorem: Let $\phi : (a,b) \mapsto (a,b)$ be continuous. Further, let $x_{k+1} = \phi(x_k)$, $x_0 \in (a,b)$, and

$$
\forall x, y \in (a,b),\ \exists \rho < 1 \text{ s.t. } \left|\phi(x) - \phi(y)\right| \leq \rho \left|x - y\right|
$$

moreover, assume that $\exists \alpha \in (a,b)$ s.t. $\phi(\alpha) = \alpha$. Then $\forall x_0 \in (a,b)$, $x_n \to \alpha$ as $n \to \infty$ (linear convergence). Note that possible underlying connection to Lipschitz continuity here. And recall that optimization can be more or less considered as a root finding procedure of the first-order optimality condition. The examples given here are chord method (corresponding to gradient descent), and Newton-Raphson method (local quadratic convergence).

5 Lagrange Interpolation The problem *Lagrange Interpolation* aims to solve is, given $(x_0, y_0), (x_1, y_1), \ldots, (x_n, y_n)$, find coefficients $a_0, \ldots, a_n$ for $p \in \mathbb{P}_n$ s.t.

$$p(x) := \begin{bmatrix} 1 & x_0 & x_0^2 & \cdots & x_0^n \\ 1 & x_1 & x_1^2 & \cdots & x_1^n \\ \vdots & & & \ddots & \vdots \\ 1 & x_n & x_n^2 & \cdots & x_n^n \end{bmatrix} \begin{bmatrix} a_0 \\ a_1 \\ \vdots \\ a_n \end{bmatrix} = \begin{bmatrix} y_0 \\ y_1 \\ \vdots \\ y_n \end{bmatrix}$$

One way to get the coefficients for the polynomial is to use *Lagrange coefficients*:

$$\phi_k(x) := \prod_{i=0,\ i\neq k}^{n} \frac{x - x_i}{x_k - x_i}$$

and

$$p(x) = \sum_{k=0}^{n} y_k \phi_k(x)$$

as we can observe that

$$\phi_j(x_i) = \begin{cases} 1, & i = j \\ 0, & i \neq j \end{cases}$$

Now define *Lagrange interpolation* as a map $\mathcal{L}_n : \mathcal{C}(a,b) \mapsto \mathbb{P}_n$, where $\{x_0, \ldots, x_n\} \subset (a,b)$ are distinct and fixed; i.e., to take $n+1$ points on f and construct the Lagrange polynomial passing through these $n+1$ points. Note that $\mathcal{L}_n$ is a projection, i.e. $\mathcal{L}_n \mathcal{L}_n = \mathcal{L}_n$. Recall we have seen how Lagrange generalized mean value theorem to higher-orders for Taylor series before, and here is the origin of Lagrange Theorem:

Let f be $n+1$th order differentiable in (a,b), and $x \in (a,b)$. Then $\exists \xi = \xi(x)$ s.t.

$$\min\{x_0, \ldots, x_n, x\} < \xi < \max\{x_0, \ldots, x_n, x\}, \text{ and} \tag{1}$$

$$f(x) - (\mathcal{L}_n f)(x) = \frac{(x - x_0) \cdots (x - x_n)}{(n+1)!} f^{(n+1)}(\xi)$$

The idea of proof is to construct the *Lagrange reminder*:

$$R(x) := f(x) - (\mathcal{L}_n f)(x);\ A := \frac{R(x)}{\prod_{i=0}^{n}(x - x_i)}$$

then the function

$$F(z) := f(z) - (\mathcal{L}_n f)(x) - A \prod_{i=0}^{n} (z - x_i)$$

has $n+2$ distinct zeros $\{x_0,\dots,x_n,x\}$; $F'(z)$ has $n+1$ distinct zeros; ...; $F^{(n+1)}(\xi) = f^{(n+1)}(\xi) - A(n+1)! = 0$ for some ξ in the convex hull as described in (1). This implies

$$f(x) - (\mathcal{L}_n f)(x) = R(x) = A\prod_{i=0}^{n}(x-x_i) = \frac{(x-x_0)\cdots(x-x_n)}{(n+1)!}f^{(n+1)}(\xi)$$

Note here how Rolle's theorem can be used to bridge the gap between higher-order in the last of the proof. Another interesting thing is that, Taylor's series expansion can be considered as Lagrange interpolation with repeated x_j.

6 Runge's Phenomenon *Runge's phenomenon* refers to the phenomenon that for a typical analytic function, equispaced Lagrange interpolation tends to oscillate more towards the boundary – that is, it tends to interpolate better in the middle. A typical analytic function will have $f^{(n)}(x) \sim \frac{n!}{\delta^n}$, and will have error$\sim \frac{\pi(x)}{\delta^n}$ for $\pi(x) = (x-x_0)\cdots(x-x_n)$. This suggests that *high-order polynomials on equispaced grid is not a good idea*, rather, it's a better idea to pick more points around the edge. Alternatively, it might be a better idea to approximate a function not by interpolating at certain points, but rather to minimize the upper bound of the approximation error norm – which leads to the discussion of the following three sections.

7 Weierstrass Approximation Theorem The *Weierstrass Approximation Theorem* states that a polynomial is dense in the space of continuous function in uniform norm: Let $f \in \mathcal{C}[a,b]$ and $\varepsilon > 0$; then $\exists n \in \mathbb{N}$, $\exists q \in \mathbb{P}_n(x)$ s.t.

$$\max_{x\in[a,b]} |f(x) - q(x)| \le \varepsilon$$

Bernstein proposed a constructive proof back in 1904. WLOG, $[a,b] = [0,1]$. Define *Bernstein polynomials* to have coefficients

$$\beta_{n,j}(x) = \binom{n}{j} x^j (1-x)^{n-j},\ j = 0,\dots,n$$

i.e., binomial polynomial if you study stats... It has a few simple properties:

1. $\beta_{n,j}(x) > 0,\ \forall x \in (0,1)$
2. $\sum_{j=0}^{n} \beta_{n,j}(x) = 1$
3. $\sum_{j=0}^{n} \frac{j}{n}\beta_{n,j}(x) = x$
4. $\sum_{j=0}^{n} \frac{j^2}{n^2}\beta_{n,j}(x) = \left(1-\frac{1}{n}\right)x^2 + \frac{1}{n}x$

And the interpolation proceeds as: let $x_j = \frac{j}{n}$, $j = 0, 1, \ldots, n$, and let $B_n f(x) = \sum_{j=0}^{n} f\left(x_j\right) \beta_{n,j}(x)$. Observe that

$$\begin{aligned} f(x) - B_n f(x) =& f(x) \sum_{j=0}^{n} \beta_{n,j}(x) - \sum_{j=0}^{n} f\left(x_j\right) \beta_{n,j}(x) \\ =& \sum_{j=0}^{n} \left[f(x) - f\left(x_j\right) \right] \beta_{n,j}(x) \end{aligned}$$

Now split the function into two components:

$$\begin{aligned} R_\delta(x) :=& \left| \sum_{|x - x_j| \le \delta} \left[f(x) - f\left(x_j\right) \right] \beta_{n,j}(x) \right| \\ \le& \underbrace{\left| \sum_{j=0}^{n} \beta_{n,j}(x) \right|}_{=1} \cdot \underbrace{\max_{y \in [0,1],\ |x-y| \le \delta} \left| f(x) - f(y) \right|}_{=:\omega(\delta)} \\ S_\delta(x) :=& \left| \sum_{|x - x_j| > \delta} \left[f(x) - f\left(x_j\right) \right] \beta_{n,j}(x) \right| \end{aligned}$$

and construct interpolation sequence of $\xi_1, \xi_2, \ldots, \xi_p$ between x and x_j s.t. the distance (in Euclidean norm) between two neighbor points $\le \delta$, then

$$\begin{aligned} \left| f(x) - f\left(x_j\right) \right| \le& |f(x) - f(\xi_1)| + |f(\xi_1) - f(\xi_2)| + \cdots + \left| f\left(\xi_p\right) - f\left(x_j\right) \right| \\ \le& (p+1)\, \omega(\delta) \\ \le& \left(1 + \frac{\left| x - x_j \right|}{\delta} \right) \omega(\delta) \end{aligned}$$

This further implies that

$$|S_\delta(x)| \le \underbrace{\sum_{|x - x_j| > \delta} \omega(\delta)\, \beta_{n,j}(x)}_{\le \omega(\delta)} + \frac{\omega(\delta)}{\delta} \underbrace{\sum_{|x - x_j| > \delta} \left| x - x_j \right| \beta_{n,j}(x)}_{=:A} \le \left(1 + \frac{1}{4\delta^2 n} \right) \omega(\delta)$$

where above inequality uses the fact that

$$\begin{aligned}
\delta A \leq & \sum_{|x-x_j|>\delta} \left(x-x_j\right)^2 \beta_{n,j}(x) \\
\leq & \sum_{j=0}^{n} \left(x-x_j\right)^2 \beta_{n,j}(x) \\
= & x^2 \sum_{j=0}^{n} \beta_{n,j}(x) - 2x \sum_{j=0}^{n} \frac{j}{n} \beta_{n,j}(x) + \sum_{j=0}^{n} \frac{j^2}{n^2} \beta_{n,j}(x) \\
= & x^2 - 2x^2 + \left(1-\frac{1}{n}\right) x^2 + \frac{1}{n} x \\
= & \frac{x(1-x)}{n} \\
\leq & \frac{1}{4n}
\end{aligned}$$

Hence,

$$|f(x) - B_n f(x)| \leq \left(2 + \frac{1}{4n\delta^2}\right) \omega(\delta), \ \forall \delta > 0 \text{ and } x \in [0,1]$$

Pick $\delta = \frac{1}{\sqrt{n}}$ completes the proof.

8 Minimax polynomials The *minimax polynomial* refers to the polynomial of a given degree that minimizes the uniform norm of the error for a continuous function on a closed interval, and its existence is ensured by the following theorem: Let $f \in \mathcal{C}[0,1]$ and $n \in \mathbb{N}_0$. Then $\exists p \in \mathbb{P}_n$ s.t.

$$\|f-p\|_\infty = \inf_{q \in \mathbb{P}_n} \|f-q\|_\infty$$

such q is called a *minimax polynomial* of degree n for f (on $[0,1]$).

The proof follows from continuous function achieves minimizer over a compact set (Weierstrass Theorem): For the sake of simplicity, let $a \in \mathbb{R}^{n+1}$ denote the coefficient vector for a nth order polynomial q, and

$$E(a) \coloneqq \|f-q\|_\infty = \max_{x \in [0,1]} |f(x) - q(x)|$$

First we are to prove the continuity of E:

$$\begin{aligned}
|E(a+\delta a)| \leq & \left| \|f-q-\delta q\|_\infty - \|f-q\|_\infty \right| \\
\leq & \|\delta q\|_\infty \\
\leq & |\delta a_0| + \cdots + |\delta a_n|
\end{aligned}$$

Now let $K := \left\{a \in \mathbb{R}^{n+1} | E(a) \leq \|f\|_\infty + 1\right\}$. Then:

1. K is closed, because $K = E^{-1}([0, \|f\|_\infty + 1])$ (pre-image of a closed set under continuous mapping is closed)
2. K is bounded, because $\|q\|_\infty \leq \underbrace{\|f - q\|_\infty}_{=:E(a)} + \|f\|_\infty$ and

$$\|a\| \leq \text{constant} \cdot \|q\|_\infty \Rightarrow E(a) \to \infty \text{ as } \|a\| \to \infty$$

3. Nonempty, because $0 \in K$

Thus, by Weierstrass Theorem, $\exists a^* \in K$ s.t. $E(a^*) = \inf_{a \in K} E(a)$ – but we still have to prove that $E(a^*) = \inf_{a \in \mathbb{R}^{n+1}} E(a)$:

$$E(a^*) \leq E(0) = \|f\|_\infty \leq \|f\|_\infty + 1 < E(a), \ \forall a \in \mathbb{R}^{n+1} \setminus K$$

9 Equioscillation Theorems Two important theorems are given to characterize minimax polynomials.

The first one is *De la Vallee Poussin Theorem*: $\forall f \in \mathcal{C}[a,b]$, $n \in \mathbb{N}_0$, $p \in \mathbb{P}_n$, if

$$f\left(x_j\right) - p\left(x_j\right) = (-1)^j e_j, \ \forall j = 0, 1, \ldots, n+1$$

where $a_0 \leq x_0 < x_1 < \cdots < x_{n+1} \leq b$, and $\operatorname{sgn} e_j = \text{constant}$ for $j = 0, 1, \ldots n+1$; then[1]

$$E_n(f) := \min_{q \in \mathbb{P}_n} \|f - q\|_\infty \geq \min_j \left|e_j\right|$$

The proof is by contradiction: assume that the conclusion is false, then

$$p\left(x_j\right) - q\left(x_j\right) = (-1)^j e_j + \underbrace{f\left(x_j\right) - q\left(x_j\right)}_{< |e_j|, \ \forall j = 0,1,\ldots,n+1}$$

$$\Rightarrow p - q \text{ has } n+1 \text{ (distinct) zeros}$$

$$\Rightarrow p \equiv q$$

but it contradicts our assumption on p and q

The second one is *Chebyshev's Oscillation Theorem*, which characterizes the minimax polynomials: $p \in \mathbb{P}_n$ is a minimax polynomial for $f \in \mathcal{C}[0,1]$ iff $f - p$ takes the value $\pm \|f - p\|_\infty$,

[1] existence of minimax polynomial was proved in Section 8

with alternating changes of sign, at least $n+2$ times in $[0,1]$. Moreover, this minimax polynomial is unique.

For statement besides uniqueness: Proof for "$\Leftarrow$" is done by DLVP, $\|f-p\|_\infty \le E_n(f) \Rightarrow \|f-p\|_\infty = E_n(f)$ by minimality of $E_n(f)$; proof for "$\Rightarrow$" is done by contradiction: assume the conclusion is false, i.e., $f-p$ takes the value $\pm\|f-p\|_\infty$ of k times for some $2 \le k \le n+1$[2], and let $\delta := \pm\|f-p\|_\infty$; then $f(x_i)-p(x_i) = (-1)^j\delta$ for $j = 1,\dots,k$. And WLOG this allows us to (quasi-)partition $[0,1]$ into k intervals split by $\xi_1,\xi_2,\dots,\xi_{k-1}$ s.t. on

$$(0,\xi_1),(\xi_2,\xi_3),\cdots : -\delta \le f-p \le \delta-\varepsilon$$
$$(\xi_1,\xi_2),(\xi_3,\xi_4),\cdots : -\delta+\varepsilon \le f-p \le \delta$$

for some $\varepsilon > 0$. Now let $r(x) = \pm(x-\xi_1)\cdots(x-\xi_{k-1})$ – we'll discuss choice of sign shortly after, and let $q(x) := p(x)-\alpha\cdot r(x)$ for some small $\alpha > 0$ s.t. $\|\alpha r\|_\infty \le \frac{\varepsilon}{2}$, then $f-q = f-p+\alpha r$. Thus on

$$(0,\xi_1),(\xi_2,\xi_3),\cdots : -\delta < -\delta+\alpha r \le f-q \le \delta-\frac{\varepsilon}{2}$$
$$(\xi_1,\xi_2),(\xi_3,\xi_4),\cdots : -\delta+\frac{\varepsilon}{2} \le f-q \le \delta+\alpha r < \delta$$

and we choose the sign of $r(x)$ s.t. $r > 0$ on the first line above and $r < 0$ on the second line above. Then q actually takes strictly less error than p, which contradicts that p is the minimax polynomial.

For uniqueness statement: let p,q both be minimax polynomials, and let $r := \frac{p+q}{2}$. Then

$$\begin{aligned}
&|f-r| \le \frac{1}{2}|f-p| + \frac{1}{2}|f-q| \le E_n(f)\\
\Rightarrow &|f-r| = E_n(f) \text{ at } n+2 \text{ distinct points}\\
\Rightarrow &f-p = f-q = \pm E_n(f) \text{ at those points – because } f-p = -(f-q) \Rightarrow f-r = 0\\
\Rightarrow &p = q \text{ at } n+2 \text{ distinct points}\\
\Rightarrow &p \equiv q
\end{aligned}$$

10 Chebyshev Polynomials Recall that the Runge's phenomenon suggests that the equispaced interpolation of polynomials does not approximate the function well, then we aims to position the interpolation points over a non-equal grid to approximate the function better. For example, we are to find the minimax polynomial in $\mathbb{P}_n$ for $f(x) = x^{n+1}$. Recall that $\sin,\cos$ usually brings oscillations, but they are not polynomials, then

[2] $k \ge 2$ because it's a minimax polynomial

Chebyshev introduced a polynomial variant from it:

$$t_n(x) := \cos(n \operatorname{arc}\cos x)$$

which will gives us $t_n(x) = 1$, $t_1(x) = x$. Recall that

$$\cos((n+1)\theta) + \cos((n-1)\theta) = 2\cos(n\theta)\cos\theta$$

translate this into $t_n(x)$, it is

$$t_{n+1}(x) = 2t_n(x)x - t_{n-1}(x)$$

these are called *Chebyshev polynomials*, the zeros of $t_{n+1}(x)$ satisfy $(n+1)\operatorname{arc}\cos x = \frac{\pi}{2} + k\pi$ for $k = 0, 1, \ldots, n$.

II Equation Solving

11 Gaussian Elimination The idea of *Gaussian Elimination*, is based on use upper rows to eliminate front-end matrix terms – one term at a time; and the resulting matrix will be an upper-triangular matrix. e.g.:

$$A = \underbrace{\begin{bmatrix} 2 & 1 & 1 \\ 4 & 3 & 3 \\ 8 & 7 & 9 \end{bmatrix}}_{A_1} \to \underbrace{\begin{bmatrix} 2 & 1 & 1 \\ 0 & 1 & 1 \\ 0 & 3 & 5 \end{bmatrix}}_{A_2} \to \underbrace{\begin{bmatrix} 2 & 1 & 1 \\ 0 & 1 & 1 \\ 0 & 0 & 2 \end{bmatrix}}_{A_3}$$

written in matrix form of above example, it will be

$$A_2 = \underbrace{\begin{bmatrix} 1 & 0 & 0 \\ -2 & 1 & 0 \\ -4 & 0 & 1 \end{bmatrix}}_{\Lambda_1} A_1,\ A_3 = \underbrace{\begin{bmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & -3 & 1 \end{bmatrix}}_{\Lambda_2} A_2$$

and note such Λ_k are always lower-triangular – in fact it only has nonzero entries at the kth column and all diagonal entries being 1, as we use upper rows to eliminate elements from lower rows.

12 LU-decomposition As a summary, in the example, $A_3 = \underbrace{\Lambda_2\Lambda_1}_{\Lambda} A$; as a product of lower-triangular matrices, Λ is also lower-triangular, hence Λ^{-1} is also lower-triangular.

And the decomposition for full rank matrix $A = \Lambda^{-1}A_3$ is called *LU-decomposition*, in practice:

1. LU-decomposition has arithmetic complexity of roughly $\frac{1}{3}n^3$ multiplications;
2. LU-decomposition breaks down if $(A_k)_{k,k} = 0$ for some k
3. L and U can be stored in a single $n \times n$ array (because Λ^{-1} always has diagonal elements all being 1)

LU decomposition of A exists iff all principal minors of A are nonzero. If exists, LU decomposition is unique. Prove by noticing that Gaussian elimination always preserves principal minors. For uniqueness, let

$$LU = \hat{L}\hat{U} \Rightarrow \underbrace{\hat{L}^{-1}L}_{\text{lower-trig}} = \underbrace{\hat{U}U^{-1}}_{\text{upper-trig}} = I \Rightarrow \hat{U} = U,\ \hat{L} = L$$

Now the issue still remains if we encounter $(A_k)_{k,k} = 0$ for some k. To solve this issue, and also to make most prominent values (measured by Euclidean norm) up to the top to ensure numerical stability, *pivoting* is introduced. *Partial pivoting* means row interchanges (arithmetic complexity n^2); and *complete pivoting* refers to row and column interchanges (arithmetic complexity $\frac{1}{3}n^3$). An example for partial pivoting row interchange:

$$\underbrace{\begin{bmatrix} 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 1 \\ 1 & 0 & 0 & 0 \\ 0 & 0 & 1 & 0 \end{bmatrix}}_{P} \cdot \begin{bmatrix} a_1 \\ a_2 \\ a_3 \\ a_4 \end{bmatrix} = \begin{bmatrix} a_2 \\ a_4 \\ a_1 \\ a_3 \end{bmatrix}$$

and the *permutation matrix* P has properties that $PP^T = P^TP = I$, and the product of permutation matrix is still a permutation matrix (recall it just interchanges rows). Now partial pivoting LU decomposition performs pivoting after each elimination step, specifically,

$$\begin{aligned} U =& \Lambda_n P_n \Lambda_{n-1} P_{n-1} \cdots \Lambda_1 P_1 A \\ =& \underbrace{\Lambda_n \left(P_n \Lambda_{n-1} P_n^{-1}\right)}_{\Lambda'_{n-1}} \underbrace{\left(P_n P_{n-1} \Lambda_{n-2} P_{n-1}^{-1} P_n^{-1}\right)}_{\Lambda'_{n-2}} \cdots \underbrace{\left(P_n \cdots P_2 \Lambda_1 P_2^{-1} \cdots P_n^{-1}\right)}_{\Lambda'_1} P_n \cdots P_2 P_1 A \end{aligned}$$

where $\Lambda'_.$ are *unit lower triangular matrices* – note that they are not lower triangular. And let $\Lambda' = \Lambda_n \Lambda'_{n-1} \Lambda'_{n-2} \cdots \Lambda'_1$, then $U = \Lambda' PA$, this gives

$$PA = LU$$

which is called *PLU-decomposition*. From the above pivoting process, it can be concluded that every square matrix has a PLU-decomposition.

13 Orthogonalization and QR-decomposition A matrix $Q \in \mathbb{R}^{n\times n}$ is called *orthogonal* if $Q^T Q = I$, i.e., if its column vectors form a orthonormal basis of $\mathbb{R}^n$. The idea of QR-decomposition comes from

$$Ax = QRx = b \Rightarrow QRx = b \Rightarrow Rx = Q^T b$$

then $Rx = Q^T b$ can be solved by back-substitution. If A, B are orthogonal, then AB and BA are both orthogonal. This allows us to perform QR-decomposition by a series of steps and times an orthogonal matrix at each step.

Recall that the projection of a on b is defined to be

$$\text{proj}_b a := \frac{\langle a, b\rangle}{\|a\| \cdot \|b\|} \cdot \|a\| \cdot \frac{b}{\|b\|} = \frac{\langle a, b\rangle}{\langle b, b\rangle} b$$

First we'll have a look at Gram-Schmidt method: let $a_1, a_2, \ldots, a_m \in \mathbb{R}^n$ be column vectors of $A \in \mathbb{R}^{n\times m}$, we can then form a orthonormal basis $q_1, q_2, \ldots, q_m$ by letting

$$\begin{aligned}
q_1 =& \frac{a_1}{\|a_1\|};\\
q'_2 =& a_2 - \langle a_2, q_1\rangle q_1,\ q_2 = \frac{q'_2}{\left\|q'_2\right\|};\\
&\vdots\\
q'_m =& a_m - \sum_{k=1}^{m-1} \langle a_m, q_k\rangle q_k,\ q_m = \frac{q'_m}{\|q'_m\|}.
\end{aligned}$$

where $\|\cdot\|$ denote Euclidean norm. i.e., in each Gram-Schmidt step, first take off the projection of a_k onto the existing orthonormal basis s.t. the remaining vector will be orthogonal to the existing basis, then normalize a_k. Applying Gram-Schmidt to perform QR-decomposition, each Gram-Schmidt step can be considered as multiplication with a triangular matrix (i.e., step k will normalize $a_k^{(k)}$, and subtract the projections on q_k from

$a_{k+1}^{(k)}, a_{k+2}^{(k)}, \ldots, a_m^{(k)}$):

$$\begin{bmatrix} a_1 & a_2 & \ldots & a_m \end{bmatrix} \begin{bmatrix} \frac{1}{\langle q_1, a_1 \rangle} & -\frac{\langle q_1, a_2 \rangle}{\langle q_1, a_1 \rangle} & -\frac{\langle q_1, a_3 \rangle}{\langle q_1, a_1 \rangle} & \cdots & -\frac{\langle q_1, a_m \rangle}{\langle q_1, a_1 \rangle} \\ 0 & 1 & 0 & & 0 \\ 0 & 0 & 1 & & 0 \\ \vdots & & & \ddots & \vdots \\ 0 & 0 & 0 & \cdots & 1 \end{bmatrix} = \begin{bmatrix} q_1 & a_2^{(2)} & a_3^{(2)} & \ldots & a_m^{(2)} \end{bmatrix}$$

Or view a_k as the sum of its projections on $q_1, q_2, \ldots, q_k$, from which we formulate

$$A = \begin{bmatrix} q_1 & q_2 & \ldots & q_m \end{bmatrix} \begin{bmatrix} \langle q_1, a_1 \rangle & \langle q_1, a_2 \rangle & \cdots & \langle q_1, a_m \rangle \\ 0 & \langle q_2, a_2 \rangle & & \langle q_2, a_m \rangle \\ \vdots & & \ddots & \vdots \\ 0 & 0 & \cdots & \langle q_m, a_m \rangle \end{bmatrix} = QR$$

where q_k is obtained using Gram-Schmidt – note that this ensures all the 0s below diagonal.

14 QR-decomposition by Triangularization *Triangularization* refers to the idea of triangularizing a matrix by zeroing its below-diagonal entries. Here two methods are discussed.

The first one is *triangularization by givens rotation*: a *(clockwise) givens rotation matrix*[3] is defined by

$$G = \begin{bmatrix} \cos\theta & \sin\theta \\ -\sin\theta & \cos\theta \end{bmatrix}$$

And for $a = \begin{bmatrix} a_1 \\ a_2 \end{bmatrix}$, to zero the second entry, i.e., to ensure

$$Ga = \begin{bmatrix} a_1 \cos\theta + a_2 \sin\theta \\ -a_1 \sin\theta + a_2 \cos\theta \end{bmatrix} = \begin{bmatrix} \sqrt{a_1^2 + a_2^2} \\ 0 \end{bmatrix}$$

we only need to let

$$\sin\theta = \frac{a_2}{\sqrt{a_1^2 + a_2^2}}$$

[3] recall we have seen them in complex analysis

$$\cos\theta = \frac{a_1}{\sqrt{a_1^2 + a_2^2}}$$

Note that givens rotation matrix is orthogonal; and generalization to zeroing the $n-$dimensional vector entry a_{k+1} can be simply done by taking an identity matrix and mutate $\begin{bmatrix} I_{kk} & I_{k(k+1)} \\ I_{(k+1)k} & I_{(k+1)(k+1)} \end{bmatrix}$ to be the givens rotation matrix. Then for $A \in \mathbb{R}^{n\times m}$, we can zero the entries of a_i, $\forall i = 1, 2, \ldots, m$ in an order of $n, n-1, \ldots, i+1$ – we have to stop at $i+1$ for a_i as further zeroing will mutate the sparse patterns for zeroed $a_1, a_2, \ldots, a_{i-1}$. This way, we can obtain an upper-triangular matrix.

The second QR-decomposition method is *Householder's reflector*. Different from how givens rotations method rotates the vector to zeroing an entry, Householder's method will reflect the vector by a hyperplane H s.t. the reflection can point to the desired direction – one column vector at a time. Specifically, for $a_1 \in \mathbb{R}^n$, we try to multiply by an orthogonal matrix Q_1 s.t. $Q_1 a_1 = \|a_1\| e_1$ where e_1 having first entry being 1 and rest of entries being 0, the hyperplane H will be orthogonal to $v \coloneqq \|a_1\| e_1 - a_1$, therefore

$$Q_1 = I - 2\frac{vv^T}{v^T v}$$

where Q_1 is orthogonal. In general,

$$Q_k = \begin{bmatrix} I & 0 \\ 0 & F \end{bmatrix}$$

where $I \in \mathbb{R}^{(k-1)\times(k-1)}$, and $F \in \mathbb{R}^{(n-k+1)\times(n-k+1)}$ s.t. $F\tilde{a}_k^{(k)} = \left\|\tilde{a}_k^{(k)}\right\| e_1$, where $\tilde{a}_k^{(k)} \in \mathbb{R}^{n-k+1}$ is $\left(\left(a_k^{(k)}\right)_k, \left(a_k^{(k)}\right)_{k+1}, \ldots, \left(a_k^{(k)}\right)_n\right)$; i.e., the upper-left sub-matrix I together with the two zero sub-matrices are to preserve obtained $a_1^{(k)}, a_2^{(k)}, \ldots, a_{k-1}^{(k)}$ from the first $k-1$ steps, and F is reflecting $\tilde{a}_k^{(k)}$ to obtain $a_k^{(k+1)}$ – which is to be preserved later; i.e., $a_k^{(k+1)} = a_k^{(k+2)} = \cdots = a_k^{(\min(m,n-1))}$.

15 Conditioning of $Ax = b$ Consider a linear system with numerical error:

$$\begin{aligned} &(A + \delta A)(x + \delta x) = b + \delta b \\ \Leftrightarrow\ &(A + \delta A)\,\delta x = \delta b - \delta A x \end{aligned}$$

From where, we would expect for all invertible A, $A+\delta A$ will also be invertible if δA is small. To solve this issue, first we'll look at *induced*[4] *matrix norm*, defined for any $A \in \mathbb{R}^{n\times m}$ by

$$\|A\| \coloneqq \sup_{x\in\mathbb{R}^m} \frac{\|Ax\|}{\|x\|} = \sup_{\|x\|\le 1} \frac{\|Ax\|}{\|x\|} = \sup_{\|x\|=1} \|Ax\|$$

Immediately following from the definition, we have

- $\|\alpha A\| = |\alpha|\|A\|$ (*absolutely homogeneous*)
- $\|A+B\| \le \|A\| + \|B\|$ (*triangle-inequality*)
- $\|A\| \ge 0$ and $\|A\| = 0 \Leftrightarrow A = 0$ (*positive-definiteness*)

and the well-known *Frobenius norm*:

$$\|A\|_F \coloneqq \sqrt{\sum_{j=1}^{m}\sum_{i=1}^{n}\left|a_{ij}\right|^2} = \operatorname{tr}\left(A^T A\right) = \operatorname{tr}\left(AA^T\right)$$

which leads to the use fact that: $\exists \alpha, \beta > 0$ s.t. $\alpha\|A\|_F \le \|A\|_* \le \beta\|A\|_F$, where $\|\cdot\|_*$ denotes any induced norm.

And for matrix geometric series, we are thinking of

$$(I-K)^{-1} = I + K + K^2 + \cdots \tag{2}$$

and (2) converges iff the $\ell - 2$ norm of all eigenvalues of A are strictly less than 1 – recall that $I-K$ is invertible iff 1 is not an eigenvalue of K. And specific for convergence proof, let

$$B_l \coloneqq I + K + \cdots + K^l$$

then we have

$$\begin{aligned}\|B_{l+m} - B_l\| &= \left\|K^{l+1} + \cdots + K^{l+m}\right\| \\ &\le \|K\|^{l+1} + \cdots + \|K\|^{l+m} \\ &\le \frac{\|K\|^{l+1}}{1-\|K\|}\end{aligned}$$

if $\|K\| < 1$. Then $\{B_l\}$ is Cauchy, which implies that $\exists B \in \mathbb{R}^{n\times m}$ s.t. $B_l \to B$ as $l \to \infty$.

Now go back to our problem, we need $A+\delta A = A\left(I + A^{-1}\delta A\right)$ to be invertible, then we'll have $(A+\delta A)^{-1} = \left(I + A^{-1}\delta A\right)^{-1} A^{-1}$; and $\left(I + A^{-1}\delta A\right)^{-1}$ exists if $\left\|A^{-1}\delta A\right\| < 1$, note that

[4] *induced* means matrix norm induced by vector norms

$\left\|A^{-1}\delta A\right\| \leq \left\|A^{-1}\right\| \|\delta A\|$, then we only need $\|\delta A\| < \frac{1}{\left\|A^{-1}\right\|}$. The rest of error analysis follows from matrix norm properties and matrix geometric series properties[5]. Eventually, it can be derived that

$$\frac{\|\delta x\|}{\|x\|} \leq \frac{\|A\| \left\|A^{-1}\right\|}{1 - \left\|A^{-1}\delta A\right\|} \left(\frac{\|\delta A\|}{\|A\|} + \frac{\|\delta b\|}{\|b\|} \right)$$

where we define the *condition number* as

$$\kappa(A) = \|A\| \left\|A^{-1}\right\|$$

16 Backward Error Analysis For floating point addition, we can treat them *as if* input were perturbed; e.g.:

$$x_1 \oplus x_2 = (x_1 + x_2)(1 + \delta) = (1 + \delta) x_1 + (1 + \delta) x_2 =: \tilde{x}_1 + \tilde{x}_2$$

Applying this treatment to the entire algorithm, we get *backward error analysis (BEA)*. Let $\tilde{f} : \mathbb{R}^n \mapsto \mathbb{R}^m$ be some algorithmic realization of $f : \mathbb{R}^n \mapsto \mathbb{R}^m$, then BEA refers to the idea of model the errors committed within $\tilde{f}$ by error in the input data. The algorithm is called *stable* if $\exists \tilde{x} \approx x$ s.t. $f(\tilde{x}) \approx \tilde{f}(x)$. If it is possible to make $f(\tilde{x}) = f(x)$, then the algorithm is called *backward stable*.

[5] which is frequently used when we deal with matrix inverse

References

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