

CONDITION: THE GEOMETRY OF NUMERICAL ALGORITHMS

MATERIAL ON A COURSE GIVEN AT JOURNÉES NATIONALES DE CALCUL FORMEL 2017
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PETER BÜRGISSER

ABSTRACT. The performance of numerical algorithms, both regarding stability and complexity, can be understood in a unified way in terms of condition numbers. This requires to identify the appropriate geometric settings and to characterize condition in geometric ways. A probabilistic analysis of numerical algorithms can be reduced to a corresponding analysis of condition numbers, which leads to fascinating problems of geometric probability and integral geometry.

This is the theme of my recent monograph *Condition*, written with Felipe Cucker, that appeared in 2013 in Springer’s Grundlehren series. The monograph is divided into three parts. Its first part deals with the solution of linear systems of equations, where many of the concepts can be explained in an elementary way. The second part is devoted to linear programming, i.e., the solution of systems of linear inequalities (there exist natural extensions to convex programming). The third part is devoted to the solution of systems of polynomial equations, focusing on Smale’s 17th problem, which asks to find a solution of a given system of n complex homogeneous polynomial equations in $n + 1$ unknowns. This problem can be solved in average (and even smoothed) polynomial time. Recently, Pierre Lairez succeeded in providing a complete solution of Smale’s 17th problem (“A deterministic algorithm to compute approximate roots of polynomial systems in polynomial average time,” to appear in *J. FoCM*).

The enclosed course material in the form of slides follows the three part structure of the monograph and attempts to illustrate the main unifying concepts and key ideas. The framework seems quite generally applicable. For instance, a numerical algorithm for computing eigenpairs of matrices, that is numerically stable and provably runs in average polynomial time, was recently developed along these lines (Armentano, Beltrán, Bürgisser, Cucker, and Shub, “A stable, polynomial-time algorithm for the eigenpair problem,” accepted for *J. EMS*).

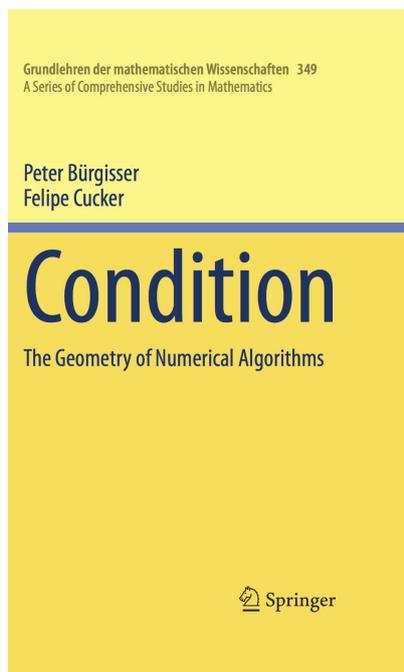
Condition: The Geometry of Numerical Algorithms

Peter Bürgisser
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Journées Nationales de Calcul Formel 2017

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For more details I refer to my new monograph (Springer 2013) with Felipe Cucker:



Outline

Overview

Part I: Linear Equalities

- Turing's Condition Number
- Average Probabilistic Analysis of $\kappa(A)$
- Smoothed Probabilistic Analysis of $\kappa(A)$
- Random Triangular Matrices

Part II: Linear Inequalities

- Interior-point methods
- Condition numbers of linear programming
- Average Analysis of GCC condition number
- Smoothed Analysis of GCC condition number
- Condition numbers of convex optimization

Part III: Polynomial Equations

- Smale's 17th Problem
- Approximate zeros, condition, and homotopy continuation
- Probabilistic analyses
- Solution to Smale's 17th problem
- Proof ideas

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- ▶ However, average analysis can rarely explain a good performance in practice. Its results strongly depend on the distribution of the inputs, which is unknown, and usually assumed to be Gaussian for rendering the mathematical analysis feasible.

Smoothed analysis

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Worst case analysis	Average case analysis	Smoothed analysis
$\sup_{a \in \mathbb{R}^p} T(a)$	$\mathbb{E}_{a \in \mathcal{D}} T(a)$	$\sup_{\bar{a} \in \mathbb{R}^p} \mathbb{E}_{a \in N(\bar{a}, \sigma^2)} T(a)$

\mathcal{D} distribution on \mathbb{R}^p , $N(\bar{a}, \sigma^2)$ Gaussian distribution centered at \bar{a} .

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- ▶ Recent new idea: **weak average-case analysis** (Amelunxen & Lotz, 2016). Take average after excluding a set of outliers of exponentially small probability.

Part I: Linear Equalities

Turing's condition number of a matrix

A. Turing, 1948

J. von Neumann and H. Goldstine, 1947

General definition of condition number

- ▶ Suppose we have a numerical computation problem

$$f: \mathbb{R}^p \rightarrow \mathbb{R}^q, \quad x \mapsto y = f(x).$$

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- ▶ Formal definition for differentiable f :

$$\kappa(f, x) := \|Df(x)\| \frac{\|x\|}{\|f(x)\|}$$

where $\|Df(x)\|$ denotes the operator norm of the Jacobian of f at x .

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- ▶ $\kappa(A)$ was introduced by **A. Turing** in 1948.

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- ▶ $\sqrt{\lambda_1}$ and $\sqrt{\lambda_n}$ are called largest and smallest singular value of A .

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where dist either refers to operator norm or to **Frobenius norm** (Euclidean norm on $\mathbb{R}^{n \times n}$) defined as

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- ▶ One calls $\log_{10} \left(\frac{\delta}{\epsilon_{\text{mach}}} \right)$ the **loss of precision** in decimal digits.

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- ▶ Let $A \in \mathbb{R}^{n \times n}$ be invertible and $b \in \mathbb{R}^n$. If the system $Ax = b$ is solved using the Householder QR factorization, the computed solution \tilde{x} has a **loss of precision bounded by**

$$\log_{10} \left(\frac{\|\tilde{x} - x\|}{\epsilon_{\text{mach}} \|x\|} \right) \leq 2 \log_{10} n + \log_{10} \kappa(A) + c,$$

where c denotes a universal constant c .

Method of conjugate gradients

- ▶ Consider a full-rank rectangular matrix $A \in \mathbb{R}^{m \times n}$ with $m > n$, a vector $c \in \mathbb{R}^m$, and the least squares problem

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- ▶ In order to achieve a relative error ε , it suffices to execute

$$\frac{1}{2} \sqrt{\kappa(S)} \ln \left(\frac{1}{\varepsilon} \right)$$

iterations (Hestenes and Stiefel, 1952).

Probabilistic Analysis of Turing's Condition Number

Average-case Analysis

H. Goldstine and J. von Neumann

Numerical inverting matrices of high order, II, 1951

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- ▶ It plays an important role in physics (cf. Wigner, 1967).

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- ▶ Application: the QR factorization has an **average loss of precision** $3 \log_{10} n + \mathcal{O}(1)$. Satisfactory result!
- ▶ There is an intuitive geometric way of deriving such results, that also has the virtue of generalizing to a wide variety of situations.

Link to geometry: reduction to sphere

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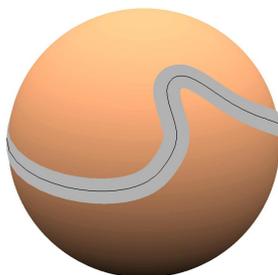
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$$\text{Prob}_A\{\kappa_F(A) \geq \varepsilon^{-1}\} = \text{Prob}_B\{\kappa_F(B) \geq \varepsilon^{-1}\} = \text{Prob}_B\{\text{dist}(B, \Sigma) \leq \varepsilon\}.$$

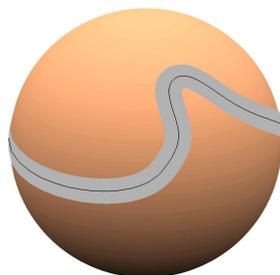
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Let $T(\Sigma_{\mathbb{S}}, \varepsilon)$ denote the neighborhood (or **tube**) of $\Sigma_{\mathbb{S}} := \Sigma \cap \mathbb{S}$ of radius $\arcsin \varepsilon$ in the sphere \mathbb{S} .



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$$\begin{aligned} \text{Prob}_B\{\kappa_F(B) \geq \varepsilon^{-1}\} &= \text{Prob}_B\{\text{dist}(B, \Sigma) \leq \varepsilon\} \\ &= \frac{\text{vol}(T(\Sigma_{\mathbb{S}}, \varepsilon))}{\text{vol}(\mathbb{S})} = \frac{\text{vol}(\Sigma_{\mathbb{S}}) \cdot 2\varepsilon}{\text{vol}(\mathbb{S})} + o(\varepsilon). \end{aligned}$$

Volume of determinant hypersurface $\Sigma_{\mathbb{S}}$ (1)

- ▶ Σ is the zero set of the determinant, a homogeneous polynomial of degree $d = n$.

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- ▶ Therefore,

$$\frac{\text{vol}(\Sigma_{\mathbb{S}})}{\text{vol}(\mathbb{S}')} = d \cdot \text{Prob}_P \{P \cap \Sigma_{\mathbb{S}} \neq \emptyset\} \leq d.$$

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we obtain the asymptotic tail bound

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- These ideas have been developed in detail by Demmel (1988).

Application to method of conjugate gradients

- ▶ The method of conjugate gradients, on input $S = A^T A$, takes

$$\frac{1}{2} \sqrt{\kappa(S)} \ln\left(\frac{1}{\varepsilon}\right) = \frac{1}{2} \kappa(A) \ln\left(\frac{1}{\varepsilon}\right).$$

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- ▶ However,

$$\text{Prob}\{\kappa(A) \geq t\} = \mathcal{O}\left(\frac{n}{t}\right)$$

implies

$$\mathbb{E}(\kappa(A)) = \infty.$$

This is inconsistent with the success of CGM in practice!

Explanation?

Condition of rectangular matrices

- ▶ CGM is usually applied to $S = R^T R$, where $R \in \mathbb{R}^{m \times n}$ is **rectangular** with $m \geq n$. (E.g., overdetermined least square problem with m linear constraints in n variables.)

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- ▶ E.g., for $4n \times n$ matrices R and large n , $\kappa(A) \simeq 3$.

Probabilistic Analysis of Turing's Condition Number

Smoothed Analysis

Smoothed analysis of $\kappa(A)$

- ▶ Take now any $\bar{A} \in \mathbb{R}^{n \times n}$, $0 < \sigma \leq 1$ and consider the isotropic Gaussian density

$$\rho(A) = \frac{1}{(\sigma\sqrt{2\pi})^{n^2}} \exp\left(-\frac{\|A - \bar{A}\|_F^2}{2\sigma^2}\right)$$

with mean \bar{A} and covariance matrix $\sigma^2 I$. Notation: $A \sim N(\bar{A}, \sigma^2 I)$.

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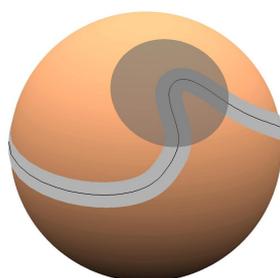
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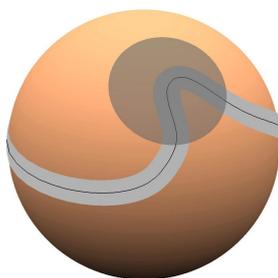
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$$\begin{aligned}
 \text{Prob}_{A \in B(\bar{A}, \sigma)} \{ \kappa_F(A) \geq \varepsilon^{-1} \} &= \text{Prob}_{A \in B(\bar{A}, \sigma)} \{ \text{dist}(A, \Sigma) \leq \varepsilon \} \\
 &= \text{Prob}_{A \in B(\bar{A}, \sigma)} \{ A \in T(\Sigma_S, \varepsilon) \} = \frac{\text{vol}(T(\Sigma_S, \varepsilon) \cap B(\bar{A}, \sigma))}{\text{vol}(B(\bar{A}, \sigma))}
 \end{aligned}$$

Uniform smoothed analysis means to provide relative bounds on the volume of tubes intersected with small spherical caps!

Heuristic estimation (1)

- ▶ Write $B := B(\bar{A}, \sigma)$. Then

$$\frac{\text{vol}(T(\Sigma_{\mathbb{S}}, \varepsilon) \cap B)}{\text{vol}(B)} \approx \frac{\text{vol}(\Sigma_{\mathbb{S}} \cap B) \cdot 2\varepsilon}{\text{vol}(B)}.$$

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- ▶ Therefore, writing $p := \dim \mathbb{S}$,

$$\begin{aligned} \frac{\text{vol}(\Sigma_{\mathbb{S}} \cap B)}{\text{vol}(B)} &= \frac{\text{vol}(\Sigma_{\mathbb{S}} \cap B)}{\text{vol}(\mathbb{S}')} \cdot \frac{\text{vol}(\mathbb{S}')}{\text{vol}(\mathbb{S})} \cdot \frac{\text{vol}(\mathbb{S})}{\text{vol} B} \\ &\lesssim d \cdot \text{Prob}_P \{P \cap B \neq \emptyset\} \cdot \sqrt{p} \cdot \frac{1}{\sigma^p}. \end{aligned}$$

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- The bound is worse by a factor n compared to Wschebor's result. But it has the advantage to be true in a much more general situation.

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B, Cucker, Lotz (2008)

For all $\sigma \in (0, 1]$ and all $t \geq (2d + 1)\frac{p}{\sigma}$,

$$\sup_{\bar{a} \in S^p} \text{Prob}_{a \in B(\bar{a}, \sigma)} \{ \mathcal{C}(a) \geq t \} \leq 26 dp \frac{1}{\sigma t}.$$

$$\sup_{\bar{a} \in S^p} \mathbb{E}_{a \in B(\bar{a}, \sigma)} (\ln \mathcal{C}(a)) \leq 2 \ln \left(\frac{dp}{\sigma} \right) + 4.7.$$



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- ▶ **Corollary:** For all $\bar{A} \in \mathbb{R}^{n \times n}$ of Frobenius norm one and $0 < \sigma \leq 1$

$$\mathbb{E}_{A \in B(\bar{A}, \sigma)}(\ln \kappa_{\text{eigen}}(A)) \leq 2 \ln \frac{n^4}{\sigma} + 5.$$

Random Triangular Matrices:

The classical condition number is not always appropriate!

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- ▶ We give a simple proof of a related result later on.
- ▶ Would the loss of precision in the solution of triangular systems conform to this bound, we would not be able to accurately find these solutions!

Explanation?

Componentwise relative errors

The classical condition number is the condition number of matrix inversion $A \mapsto A^{-1}$:

$$\kappa(A) = \lim_{\delta \rightarrow 0} \sup_{\text{RelError}(A) \leq \delta} \frac{\text{RelError}(A^{-1})}{\text{RelError}(A)}.$$

Here, we use the [normwise relative error](#)

$$\text{RelError}(A) := \frac{\|\tilde{A} - A\|}{\|A\|},$$

with the spectral norm $\|\cdot\|$.

Componentwise condition number

- ▶ Instead of RelError we may use the possibly much larger **componentwise relative error**

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$$\text{CwRelError}(A) := \max_{i,j} \frac{\|\tilde{a}_{ij} - a_{ij}\|}{\|a_{ij}\|}.$$

- ▶ We define the **componentwise condition number** of matrix inversion correspondingly as

$$\text{Cw}^\dagger(A) := \lim_{\delta \rightarrow 0} \sup_{\text{CwRelError}(A) \leq \delta} \frac{\text{CwRelError}(A^{-1})}{\text{CwRelError}(A)}.$$

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- ▶ This explains why linear triangular systems can be solved by backward substitution with high accuracy.

Why random triang. matrices are ill-conditioned (1)

Let $L = (\ell_{ij})$ denote a random **unit lower-triangular** matrix with $\ell_{ii} = 1$ and with independent standard Gaussian random entries ℓ_{ij} for $i > j$.

Then we have

$$\mathbb{E}(\|L^{-1}\|_F^2) = 2^n - 1.$$

In particular, $\mathbb{E}(\|L\|_F^2 \|L^{-1}\|_F^2) \geq n(2^n - 1)$, hence $\mathbb{E}(\kappa(L)^2)$ grows exponentially in n .

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- ▶ Hence s_i is a function of the first i rows of L and thus independent of the entries of L in the rows with index larger than i .

Why random triang. matrices are ill-conditioned (2)

- ▶ By squaring we obtain for $i \geq 2$

$$s_i^2 = \sum_{\substack{j \neq k \\ j, k < i}} l_{ij} l_{ik} s_j s_k + \sum_{j < i} l_{ij}^2 s_j^2.$$

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- ▶ Solving this recursion with $\mathbb{E}(s_1^2) = 1$ yields

$$\mathbb{E}(s_i^2) = 2^{i-2} \quad \text{for } i \geq 2.$$

Why random triang. matrices are ill-conditioned (3)

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- ▶ By an analogous argument one shows that

$$\mathbb{E}(\|v_k\|^2) = 2^{n-k}$$

for the k th column v_k of L^{-1} . Altogether, we obtain

$$\mathbb{E}(\|L^{-1}\|_F^2) = \mathbb{E}\left(\sum_{k=1}^n \|v_k\|^2\right) = \sum_{k=1}^n \mathbb{E}(\|v_k\|^2) = 2^n - 1.$$

□

Part II: Linear Inequalities

Interior-point methods for linear programming

Linear programming (1)

- ▶ **Standard primal form of linear programs:** given $A \in \mathbb{R}^{m \times n}$, $b \in \mathbb{R}^m$, $c \in \mathbb{R}^n$; look for optimal $x \in \mathbb{R}^n$

$$\min c^T x \quad \text{subject to } Ax = b, x \geq 0 \quad (\text{P})$$

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- ▶ It is known that $\max b^T y = \min c^T x$ if (P) and (D) are both feasible (duality).

Linear programming (1)

- ▶ **Standard primal form of linear programs:** given $A \in \mathbb{R}^{m \times n}$, $b \in \mathbb{R}^m$, $c \in \mathbb{R}^n$; look for optimal $x \in \mathbb{R}^n$

$$\min c^T x \quad \text{subject to } Ax = b, x \geq 0 \quad (\text{P})$$

- ▶ **Standard dual form of linear programs:** Given A, b, c , look for optimal $y \in \mathbb{R}^m$.

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- ▶ Suppose that (P) and (D) are both feasible, The vector $s := c - A^T y$ of **slack variables** satisfies

$$A^T y + s = c, \quad s \geq 0,$$

hence, using $Ax = b$,

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- ▶ Optimality is equivalent to $s^T x = 0$, which is equivalent to the **complementary slackness condition**

$$x_i s_i = 0, \quad i = 1, 2, \dots, n. \quad (1)$$

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- ▶ It can be shown that, if $\text{rank } A = m$, there is exactly one solution ζ_μ of this system, for all $\mu > 0$.

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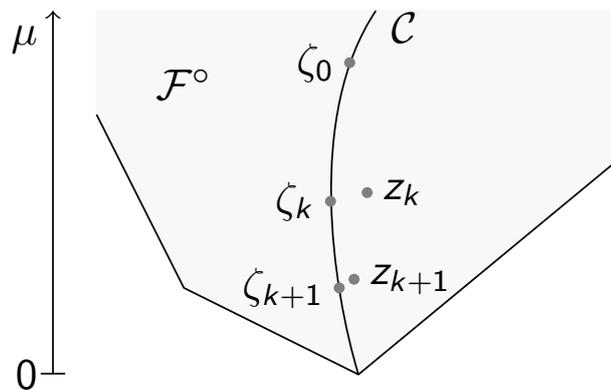
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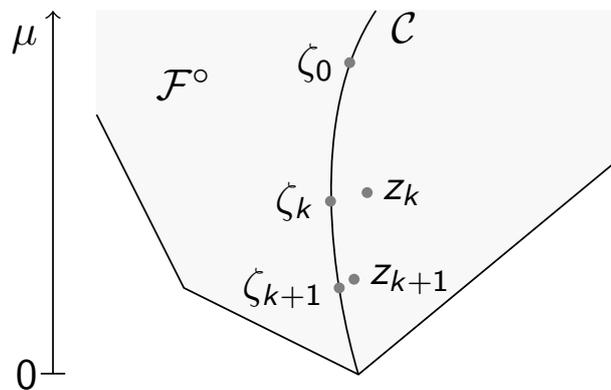
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- ▶ The **duality measure** of $z = (x, y, s) \in \mathcal{F}^\circ$ is defined as

$$\mu(z) := \frac{1}{n} \sum_{i=1}^n x_i s_i.$$

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$$z = (x, y, s) \mapsto F(z) = (A^T y + s - c, Ax - b, x_1 s_1, \dots, x_n s_n)$$

satisfying $\{\zeta_\mu\} = F^{-1}(0, 0, \mu e_n)$, where $e_n := (1, \dots, 1) \in \mathbb{R}^n$. The Jacobian matrix of F at z equals

$$DF(z) = \begin{bmatrix} 0 & A^T & I \\ A & 0 & 0 \\ S & 0 & X \end{bmatrix},$$

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- ▶ Fact: $DF(z)$ is invertible if $\text{rank } A = m$ and $s_i x_i \neq 0$ for all i .

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- ▶ Set $\zeta_k = \zeta_{\mu_k}$. Then $F(\zeta_k) = (0, 0, \mu_k e_n)$ for all $k \in \mathbb{N}$. A first order approximation gives

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- ▶ One easily checks for $z_{k+1} = z_k + (\Delta x, \Delta y, \Delta s)$

$$A^T(y + \Delta y) + (s + \Delta s) = c, \quad A(x + \Delta x) = b.$$

Primal-dual IPM

We choose $\sigma := 1 - \frac{1}{4\sqrt{n}}$.

Algorithm: Primal-Dual IPM

Input: $A \in \mathbb{R}^{m \times n}$, $b \in \mathbb{R}^m$, $c \in \mathbb{R}^n$ s.t. $\text{rank } A = m \leq n$.

Choose starting point $z_0 = (x^0, y^0, s^0) \in \mathcal{F}^\circ$ with duality measure μ_0 .

for $k = 0, 1, 2, \dots$

Solve

$$\begin{bmatrix} 0 & A^T & I \\ A & 0 & 0 \\ S^k & 0 & X^k \end{bmatrix} \cdot \begin{bmatrix} \Delta x^k \\ \Delta y^k \\ \Delta s^k \end{bmatrix} = \begin{bmatrix} 0 \\ 0 \\ \sigma^{k+1} \mu_0 e_n - X^k S^k e_n \end{bmatrix},$$

where $X^k = \text{diag}(x_1^k, \dots, x_n^k)$, $S^k = \text{diag}(s_1^k, \dots, s_n^k)$.

Set

$$(x^{k+1}, y^{k+1}, s^{k+1}) = (x^k, y^k, s^k) + (\Delta x^k, \Delta y^k, \Delta s^k).$$

until some stopping criterion is matched

Analysis of IPM

For the following, see Wright (1997).

Theorem.

Primal-Dual IPM produces, on a strictly feasible starting point z_0 on the central path (or close to it), a sequence of iterates $z_k \in \mathcal{F}^\circ$ such that $\mu(z_k) = \sigma^k \mu(z_0)$. After

$$k \geq 4\sqrt{n} \ln \frac{\mu_0}{\varepsilon}.$$

iterations we have $\mu(z^k) \leq \varepsilon$.

Condition numbers of linear programming

Jim Renegar, 1995

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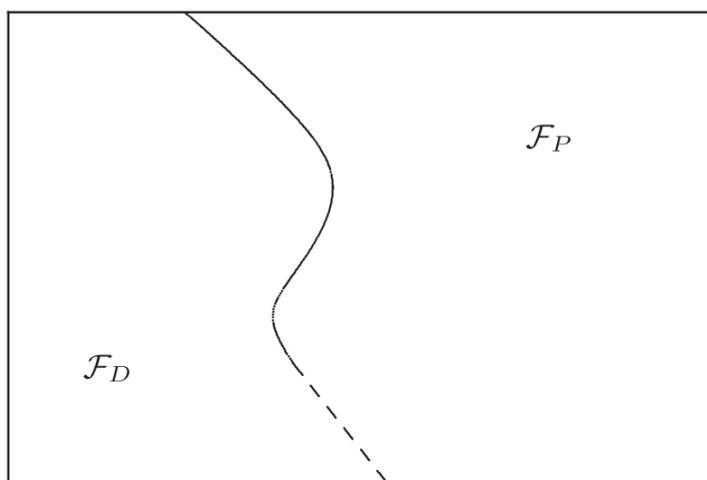
- ▶ Let \mathcal{F}_P° and \mathcal{F}_D° denote the set of instances where P and D are solvable, respectively.
- ▶ We have a disjoint union

$$\mathbb{R}^{n \times m} = \mathcal{F}_P^\circ \cup \mathcal{F}_D^\circ \cup \Sigma,$$

where the set of **ill-posed instances** Σ is the common boundary of \mathcal{F}_P° and \mathcal{F}_D° .

Linear Programming Feasibility Problem (2)

$$\mathbb{R}^{n \times m} = \mathcal{F}_P^\circ \cup \mathcal{F}_D^\circ \cup \Sigma,$$



The **Homogeneous Linear Programming Feasibility problem (HLPF)** is to decide for given A , whether $A \in \mathcal{F}_P^\circ$ or $A \in \mathcal{F}_D^\circ$.

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Primal-Dual IPM can be solved with a number of iterations bounded by

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- ▶ Consequence: HLPF can be solved in polynomial time for an integer matrix A , counting bit operations.

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- ▶ Recall $\Sigma = \overline{\mathcal{F}_P^\circ} \cap \overline{\mathcal{F}_D^\circ}$.

- ▶ Hence A is ill-posed, $A \in \Sigma$, iff Δ is contained in a closed halfspace and $0 \in \Delta$.

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$$\Delta(A) := \sup \left\{ \delta > 0 \mid \forall A' \in \mathbb{R}^{m \times n} \left(\max_{i \leq n} \frac{\|a'_i - a_i\|}{\|a_i\|} < \delta \Rightarrow A' \in \mathcal{F}_S^\circ \right) \right\},$$

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- ▶ The **GCC-condition number** of A is defined as (Goffin (1980), Cheung, Cucker (2001)):

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- ▶ Note that we measure the **relative size** of the perturbation **for each column a_i** with respect to the norm of a_i .
- ▶ Also, $\Delta(A)$ is scale invariant. We may therefore assume, without loss of generality, that $\|a_i\| = 1$ for all i .

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- ▶ We are going to define a variant of Renegar's condition number, that is better suited for probabilistic analysis.
- ▶ Suppose $A \in \mathcal{F}_S^\circ$ for $S \in \{P, D\}$. We define

$$\Delta(A) := \sup \left\{ \delta > 0 \mid \forall A' \in \mathbb{R}^{m \times n} \left(\max_{i \leq n} \frac{\|a'_i - a_i\|}{\|a_i\|} < \delta \Rightarrow A' \in \mathcal{F}_S^\circ \right) \right\},$$

where a'_i stands for the i th column of A' .

- ▶ The **GCC-condition number** of A is defined as (Goffin (1980), Cheung, Cucker (2001)):

$$\mathcal{C}(A) := 1/\Delta(A).$$

- ▶ Note that we measure the **relative size** of the perturbation **for each column a_i** with respect to the norm of a_i .
- ▶ Also, $\Delta(A)$ is scale invariant. We may therefore assume, without loss of generality, that $\|a_i\| = 1$ for all i .
- ▶ Hence we can interpret A with columns a_1, \dots, a_n as an element in the **product $\mathbb{S}^n = \mathbb{S} \times \dots \times \mathbb{S}$ of spheres $\mathbb{S} := \mathbb{S}^{m-1}$** .

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- ▶ We note that HLPF can be solved by a primal-dual interior-point method with a number of iterations

$$\mathcal{O}\left(\sqrt{n} \log(n \mathcal{C}(A))\right) .$$

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Cheung & Cucker (2001)

$$d_{\mathbb{S}}(A, \Sigma) = \begin{cases} \frac{\pi}{2} - \rho(A) & \text{if } A \in \mathcal{F}_D^\circ \\ \rho(A) - \frac{\pi}{2} & \text{if } A \in \mathbb{S}^n \setminus \mathcal{F}_D^\circ \end{cases} .$$

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Average Analysis of GCC condition number

GCC condition number and coverage processes (1)

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- ▶ Consequence: the expected number of iterations of interior point methods for HLPF is $\mathcal{O}(\sqrt{n} \log n)$.

Closed formula for $p(n, m, \alpha)$

- ▶ For $\alpha \geq \pi/2$, setting $\varepsilon := |\cos(\alpha)|$,

$$p(n, m+1, \alpha) = \sum_{k=1}^m \binom{n}{k+1} C(m, k) \int_{\varepsilon}^1 t^{m-k} (1-t^2)^{\frac{1}{2}km-1} \lambda_m(t)^{n-k-1} dt.$$

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- ▶ Let \mathcal{O}_m denote the m -dimensional volume of the sphere S^m .

$$\begin{aligned} \frac{\text{vol}(\Sigma) \cdot \varepsilon}{\text{vol}(S)^n} + o(\varepsilon^2) &= \\ &= \text{Prob} \{A \in \mathcal{F}_D^\circ, \mathcal{C}(A)^{-1} \leq \varepsilon\} = p(n, m, \pi/2) - p(n, m, \alpha) \\ &= \binom{n}{m+1} (m+1) \frac{\mathcal{O}_{m-1}}{\mathcal{O}_m} \frac{1}{2^{n-2}} \varepsilon + o(\varepsilon^2). \end{aligned}$$

Smoothed Analysis of GCC condition number

Gaussian smoothed analysis

- ▶ Model for local perturbations: $\bar{A} \in \mathbb{R}^{m \times n}$, Gaussians $A \in \mathbb{R}^{m \times n}$.

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- ▶
- ▶ This implies the bound $\mathcal{O}(\sqrt{n} \log \frac{n}{\sigma})$ on the smoothed expected number of iterations of the IPM considered for HLPF. **Excellent result!**

Uniform smoothed analysis of \mathcal{C}

- ▶ Model for smoothed analysis on product of spheres: $\bar{a}_1, \dots, \bar{a}_n \in \mathbb{S}$, independently choose a_i uniformly at random in spherical cap $B(\bar{a}_i, \sigma)$ of \mathbb{S} centered at \bar{a}_i with angular radius $\arcsin \sigma$. That is, choose $A \in B(\bar{A}, \sigma) := \prod_i B(\bar{a}_i, \sigma)$ uniformly.

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$$\sup_{\bar{A} \in \mathbb{S}^n} \text{Prob}_{A \in B(\bar{A}, \sigma)} \{A \in \mathcal{F}_D^\circ, \mathcal{C}(A) \geq \varepsilon^{-1}\} \leq 6.5 nm^2 \frac{\varepsilon}{\sigma}.$$

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- ▶ We even obtain robustness results.

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- ▶ The proof idea is similar to the previously mentioned (volume of tubes, integral geometry, counting argument).
- ▶ In particular, Poincaré's formula implies

$$\frac{\text{vol}(\partial K)}{\text{vol}(\mathbb{S}')} \leq 1.$$

Indeed, by convexity, the intersection of ∂K with a hyperequator \mathbb{S}' of \mathbb{S} in general position consists of at most two points.

Sketch of proof (2)

- ▶ **Crucial Lemma.** Let $A = (a_1, \dots, a_n) \in \mathcal{F}_D^\circ$ and $\mathcal{C}(A) \geq m\varepsilon^{-1}$. Then there exists $i \in \{1, \dots, n\}$ such that $a_i \in T_o(\partial K_i, \varepsilon)$, where $-K_i$ is the spherical convex hull of $a_1, \dots, a_{i-1}, a_{i+1}, \dots, a_n$.

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- ▶ The Lemma yields with $t = m/\varepsilon$

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- ▶ We bound the probability on the right-hand side for $i = n$ by an integral of probabilities conditioned on $A' := (a_1, \dots, a_{n-1})$:

$$\begin{aligned} & \text{Prob}\{A' \in \mathcal{F}_D^\circ \text{ and } a_n \in T_o(\partial K_n, \varepsilon)\} \\ &= \frac{1}{\text{vol}B(\bar{A}', \sigma)} \int_{A' \in \mathcal{F}_D^\circ \cap B(\bar{A}', \sigma)} \text{Prob}\{a_n \in T_o(\partial K_n, \varepsilon) \mid A'\} dA'. \end{aligned}$$

Sketch of proof (3)

- Fix now $A' \in \mathcal{F}_{n-1,m}$ and consider the convex set K_n in \mathbb{S} . The volume bound (*) yields

$$\text{Prob}\{a_n \in T_o(\partial K_n, \varepsilon) \mid A'\} = \frac{\text{vol}(T_o(\partial K_n, \varepsilon) \cap B(\bar{a}_n, \sigma))}{\text{vol}B(\bar{a}_n, \sigma)} \leq 6.5 m \frac{\varepsilon}{\sigma}.$$

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We conclude that

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- The same upper bound holds for any K_i . Altogether, we obtain

$$\text{Prob}\{A \in \mathcal{F}_D^\circ \text{ and } \mathcal{C}(A) \geq t\} \leq 6.5 nm^2 \frac{1}{\sigma t},$$

□

Condition Numbers of Convex Optimization

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- ▶ **Homogeneous convex feasibility problem (HCFP)**

Input $A \in \mathbb{R}^{m \times n}$ ($n > m$)

Decide the alternative

$$\exists x \in \mathbb{R}^n \setminus \{0\} : Ax = 0, x \in \check{C} \quad (\text{P})$$

$$\exists y \in \mathbb{R}^m \setminus \{0\} : A^T y \in C \quad (\text{D})$$

Convex homogeneous feasibility problem

- ▶ Most important cases:

Linear Programming : $C = \mathbb{R}_+^n = \mathbb{R}_+ \times \dots \times \mathbb{R}_+$

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$$\mathcal{F}_P \quad := \quad \{A \mid (P) \text{ is feasible}\},$$

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- ▶ Renegar's condition number is defined as:

$$\mathcal{C}_R(A) := \frac{\|A\|}{\text{dist}(A, \Sigma)}.$$

Convex homogeneous feasibility problem

- ▶ The probabilistic analyses for LP-condition numbers rely on the product structure of the cone $C = \mathbb{R}_+^n = \mathbb{R}_+ \times \dots \times \mathbb{R}_+$.

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- ▶ For general cones (like SDP), we look for a different, more coordinate-free approach.
- ▶ Suppose $A \in \mathbb{R}^{m \times n}$ has rank m . Consider the m -dimensional linear subspace $W := \text{im}A^T$ of \mathbb{R}^n .

$$\begin{array}{ccc}
 \exists x \in \mathbb{R}^n \setminus \{0\} : Ax = 0 & \text{(P)} & \exists y \in \mathbb{R}^m \setminus \{0\} : A^T y \in C & \text{(D)} \\
 x \in \check{C} & & & \\
 \Leftrightarrow & & \Leftrightarrow & \\
 \underbrace{\ker A}_{=: W^\perp} \cap \check{C} \neq \{0\} & & \underbrace{\text{im}A^T}_{=: W} \cap C \neq \{0\} &
 \end{array}$$

Grassmann condition number (1)

- ▶ Consider the inputs as an element of the Grassmann manifold

$$W \in \mathbb{G}_{n,m} := \{W \subseteq \mathbb{R}^n \mid W \text{ lin. subspace, } \dim W = m\}.$$

We have to decide the alternative

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- ▶ $\mathbb{G}_{n,m}$ is a compact Riemannian manifold. We have thus well-defined notions of (geodesic) distance (“angle”) and volume.

Grassmann condition number (2)

- ▶ We define the **Grassmann condition number** for $W \in \mathbb{G}_{n,m}$ as

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- ▶ The following result (Amelunxen (2011), Belloni, Freund (2009)) separates Renegar's condition number into an "intrinsic" and "extrinsic" part.

For $A \in \mathbb{R}^{m \times n}$ of rank m and $W := \text{im}A^T$ we have

$$\mathcal{C}_{\mathbb{G}}(A) \leq \mathcal{C}_R(A) \leq \kappa(A) \cdot \mathcal{C}_{\mathbb{G}}(A).$$

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$$\text{Prob} \left(\mathcal{L}_{\mathbb{G}}(A) \geq \frac{1}{\varepsilon} \right) \leq 6 \cdot n \varepsilon \quad \text{if } \varepsilon < n^{-\frac{3}{2}}.$$

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- ▶ It is a challenge to extend this result to a uniform smoothed analysis.

Part III: Polynomial Equations

Complexity of Bezout's Theorem

(Shub and Smale 1993–1996)

Smale's 17th problem

The 17th of S. Smale's problems for the 21st century asks:

*Can a zero of n complex polynomial equations in n unknowns be found **approximately, on the average**, in polynomial time with a uniform algorithm?*

Notations

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$$\mathcal{H}_d := \{f = (f_1, \dots, f_n) \mid f_i \in \mathbb{C}[X_0, \dots, X_n] \text{ homogeneous of degree } d_i\}.$$

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- ▶ The input size is $N := \dim_{\mathbb{C}} \mathcal{H}_d$.
- ▶ We look for zeros ζ of f in complex projective space \mathbb{P}^n : $f(\zeta) = 0$.

Bombieri-Weyl inner product

- ▶ For homogeneous polynomials of degree d_i ,

$$f_i(x) = \sum_{|\alpha|=d_i} a_\alpha^i X^\alpha, \quad g_i(x) = \sum_{|\alpha|=d_i} b_\alpha^i X^\alpha,$$

we define the **Bombieri-Weyl hermitian inner product** as

$$\langle f_i, g_i \rangle := \sum_{|\alpha|=d_i} a_\alpha^i \overline{b_\alpha^i} \binom{d_i}{\alpha}^{-1}.$$

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- ▶ For $f, g \in \mathcal{H}_d$ we define $\langle f, g \rangle := \sum_{i=1}^n \langle f_i, g_i \rangle$ and $\|f\| := \langle f, f \rangle^{1/2}$.

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- ▶ We have a **standard Gaussian** distribution on \mathcal{H}_d with density

$$\rho(f) = \frac{1}{\sqrt{2\pi}^{2N}} \exp\left(-\frac{1}{2}\|f\|^2\right).$$

Approximate zeros

- ▶ For $f \in \mathcal{H}_d$ we define the **Newton operator** $N_f: \mathbb{P}^n \rightarrow \mathbb{P}^n$ by

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- ▶ Here the distance d refers to the geodesic distance (angle) on the Riemannian manifold \mathbb{P}^n (Fubini-Study metric).

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- ▶ We define the **condition number** of f at (f, ζ) by

$$\mu(f, \zeta) := \|f\| \cdot \|DG(f)\|.$$

Radius of quadratic convergence

Put $D := \max_i d_i$.

Combining Smale's Gamma Theorem (1986) with the developments in Shub and Smale (1993–1996), one obtains:

Version of Smale's Gamma Theorem

If $d(x, \zeta) \leq \frac{0.3}{D^{3/2} \mu(f, \zeta)}$, then x is an approximate zero of f associated with ζ .

Adaptive linear homotopy continuation (1)

- ▶ Given a start system $(g, \zeta) \in V$ and an input $f \in \mathcal{H}_d$.

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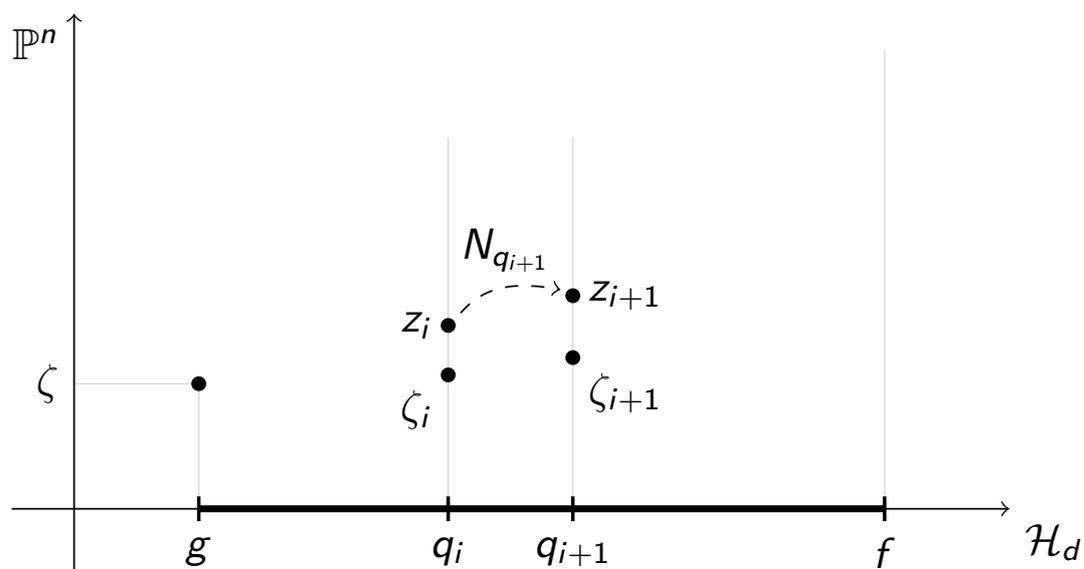
$$q_t := (1 - t)g + tf \quad \text{for } t \in [0, 1].$$

- ▶ If $[g, f]$ does not meet the discriminant variety (i.e., none of the q_t has a multiple zero), then there exists a unique lifting to V ,

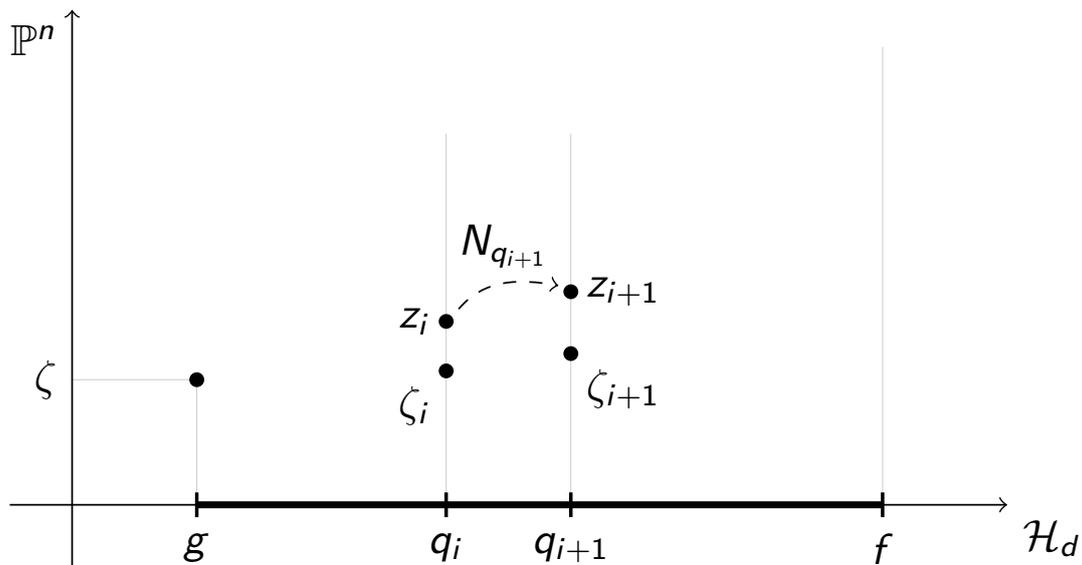
$$\gamma: [0, 1] \rightarrow V, t \mapsto (f_t, \zeta_t),$$

such that $f_0 = g$.

Adaptive linear homotopy continuation (2)



Adaptive linear homotopy continuation (2)



- The idea is to **follow the path γ numerically**: partition $[0, 1]$ into $t_0 = 0, \dots, t_k = 1$. Writing $q_i := q_{t_i}$, successively compute approximations z_i of ζ_{t_i} by Newton's method starting with $z_0 := \zeta$. More specifically, compute

$$z_{i+1} := N_{q_{i+1}}(z_i).$$

Complexity of adaptive linear homotopy continuation

- ▶ We compute t_{i+1} **adaptively** from t_i such that

$$d(q_{i+1}, q_i) = \frac{c}{D^{3/2} \mu^2(q_i, z_i)}.$$

This defines the **Adaptive Linear Homotopy ALH** algorithm.

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Shub & Smale (1994), Shub (2007)

z_i is an approximate zero of ζ_i for all i . Moreover,

$$K(f, g, \zeta) \leq 217 D^{3/2} \int_0^1 \mu_{\text{norm}}(\gamma(t))^2 \|\dot{\gamma}(t)\| dt.$$

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- ▶ **Las Vegas Algorithm LV**
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- ▶ LV has the **expected "running time"**

$$K(f) := \mathbb{E}_{g, \zeta} K(f, g, \zeta).$$

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- ▶ Note that randomness enters here in two ways: as an algorithmic tool and as a way to measure the performance of algorithms.

Smoothed expected polynomial time

Smoothed analysis: let $\bar{f} \in \mathcal{H}_d$ and suppose that f is isotropic Gaussian with mean \bar{f} and variance σ^2 .

Smoothed analysis (B, Cucker (2011))

$$\sup_{\|f\| \leq 1} \mathbb{E}_{f \sim N(\bar{f}, \sigma^2 I)} K(f) = \mathcal{O}\left(\frac{D^{3/2} N n}{\sigma}\right).$$

A near solution to Smale's 17th problem

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There is a **deterministic algorithm** for Smale's 17th problem taking on standard Gaussian input $f \in \mathcal{H}_d$ an expected number of arithmetic operations $T(f)$ bounded by

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- ▶ If $D \leq n$, the algorithm runs ALH with the start system (g, ζ) , where

$$g_i = X_i^{d_i} - X_0^{d_i}, \quad \zeta = (1, \dots, 1)$$

(the zeros of g_i consist of roots of unity).

A near solution to Smale's 17th problem

B, Cucker (2011)

There is a **deterministic algorithm** for Smale's 17th problem taking on standard Gaussian input $f \in \mathcal{H}_d$ an expected number of arithmetic operations $T(f)$ bounded by

$$\mathbb{E}_f T(f) = N^{\mathcal{O}(\log \log N)}.$$

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- ▶ If $D \geq n$, the algorithm runs ALH with the start system (g, ζ) , where

$$g_i := X_0^{d_i-i} X_i, \quad \zeta = (1, 0, \dots, 0);$$

cf. Armentano, B, Béltran, Cucker, Shub, (2016).

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Lairez (2016)

There is a **deterministic algorithm** for Smale's 17th problem taking on standard Gaussian input $f \in \mathcal{H}_d$ an expected number of arithmetic operations $T(f)$ bounded by polynomial in the input size N .

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- ▶ $\mathbb{E}(\|f\|^2) = 2N$ (chi-square). Replace $\|f\|$ by \sqrt{N} (cheating a bit).

Proof idea for smoothed analysis of ALH (2)



$$\mathbb{E}_{\zeta \in V(\mathbf{g})} K(\mathbf{f}, \mathbf{g}, \zeta) \leq c D^{3/2} N \int_0^1 \frac{\mu_2(\mathbf{q}_t)^2}{\|\mathbf{q}_t\|^2} dt.$$

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▶ By Fubini,

$$\mathbb{E}_{f \sim N(\bar{f}, \sigma^2 I)} \mathbb{E}_{g \sim N(0, I)} \mathbb{E}_{\zeta \in V(g)} K(f, g, \zeta) \leq c D^{3/2} N \int_0^1 \mathbb{E} \left(\frac{\mu_2(q_t)^2}{\|q_t\|^2} \right) dt.$$

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► For fixed t , $q_t = (1-t)g + tf$ is **again Gaussian**, $q_t \sim N(\bar{q}_t, \sigma_t^2 I)$, with

$$\bar{q}_t = t\bar{f}, \quad \sigma_t^2 = (1-t)^2 + \sigma^2 t^2.$$

Proof idea for smoothed analysis of ALH (3)

Main technical contribution of proof

$$\mathbb{E}_{q \sim N(\bar{q}, \sigma^2 I)} \left(\frac{\mu_2(q)^2}{\|q\|^2} \right) = \mathcal{O}\left(\frac{n}{\sigma^2}\right).$$

Using this,

$$\mathbb{E}_{f \sim N(\bar{f}, \sigma^2 I)} K(f) \leq c D^{3/2} N \int_0^1 \frac{n}{(1-t)^2 + \sigma^2 t^2} dt = c D^{3/2} N \frac{n}{\sigma}. \quad \square$$

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A slightly better estimate, with a simpler proof, was obtained by Armentano, B, Béltran, Cucker, Shub (2016).

On proving the main technical contribution (1)

- ▶ Put $\mathcal{M} := \mathbb{C}^{n \times (n+1)}$ and consider the map (slightly cheating ...)

$$\Psi: V \rightarrow \mathcal{M}, (q, \zeta) \mapsto M := \text{diag}(\sqrt{d_1}, \dots, \sqrt{d_n})^{-1} Df(\zeta).$$

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Recall $\mu(q, \zeta) / \|q\| = \|M^\dagger\|$.

- ▶ The noncentered Gaussian on \mathcal{H}_d defines a distribution on V (choose q and then one of its \mathcal{D} zeros uniformly at random). Then

$$\mathbb{E}_{\mathcal{H}_d} \left(\frac{\mu_2(q)^2}{\|q\|^2} \right) = \mathbb{E}_V \left(\frac{\mu(q, \zeta)^2}{\|q\|^2} \right) = \mathbb{E}_{\mathcal{M}} (\|M^\dagger\|^2)$$

where the last expectation is w.r.t. the distribution on \mathcal{M} induced by Ψ .

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- ▶ L_ζ is isometrically isomorphic to $\mathcal{M}_\zeta := \{M \in \mathcal{M} : M\zeta = 0\}$ inducing a Gaussian $N(\bar{M}_\zeta, \sigma^2 I)$ on the fiber \mathcal{M}_ζ .

On proving the main technical contribution (3)

- ▶ For $M \in \mathcal{M}$ of full rank with zero ζ one can show that

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$$\begin{aligned} \mathbb{E}_{\mathcal{M}}(\|M^{\dagger}\|^2) &= \int_{\mathcal{M}} \|M^{\dagger}\|^2 \rho_{\mathcal{M}}(M) dM \\ &= \mathbb{E}_{\zeta \in \mathbb{P}^n} \left(\mathbb{E}_{\tilde{\rho}_{\mathcal{M}_{\zeta}}} (\|M^{\dagger}\|^2) \right) \end{aligned}$$

Right expectation is over induced distribution of the zeros ζ of M , second expectation is w.r.t. the following **conditional density on \mathcal{M}_{ζ}** :

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- ▶ Hence $\mathbb{E}_{\mathcal{M}}(\|M^{\dagger}\|^2) = \mathcal{O}\left(\frac{n}{\sigma^2}\right)$. □

Thank you for your attention!

REFERENCES

- [1] E.L. Allgower and K. Georg. *Numerical Continuation Methods*. Springer-Verlag, 1990.
- [2] D. Amelunxen. *Geometric analysis of the condition of the convex feasibility problem*. PhD thesis, University of Paderborn, 2011.
- [3] D. Amelunxen and P. Bürgisser. Probabilistic analysis of the Grassmann condition number. *Found. Comput. Math.*, 15(1): 3–51, 2015.
- [4] D. Amelunxen and P. Bürgisser. A coordinate-free condition number for convex programming. *SIAM J. Optim.*, 22(3):1029–1041, 2012.
- [5] D. Amelunxen and P. Bürgisser. Robust smoothed analysis of a condition number for linear programming. *Math. Programming*, 131(1, Ser. A):221–251, 2012.
- [6] D. Amelunxen and M. Lotz. Average-case analysis without the black swans. Accepted for *J. Compl.*, arXiv:1512.09290.
- [7] D. Armentano. Stochastic perturbations and smooth condition numbers. *J. Complexity*, 26(2):161–171, 2010.
- [8] D. Armentano. Complexity of path-following methods for the eigenvalue problem. *Found. Comput. Math.*, 14(2): 185–236, 2014.
- [9] D. Armentano, C. Beltrán, P. Bürgisser, F. Cucker, and M. Shub. Condition length and complexity for the solution of polynomial systems. *Found. Comput. Math.*, 16(6): 1401–1422, 2016.
- [10] E. Barbier. Note sur le problème de l’aiguille et le jeu du joint couvert. *J. Math. Pures et Appl.*, 5(2):273–286, 1860.
- [11] V. Bargmann, D. Montgomery, and J. von Neumann. Solution of linear systems of high order (Princeton, 1946). In A.H. Taub, editor, *John von Neumann Collected Works*, volume 5. Pergamon, Elmsford, NY, 1963.
- [12] D.J. Bates, J.D. Hauenstein, A.J. Sommese, and C.W. Wampler. Software for numerical algebraic geometry: a paradigm and progress towards its implementation. In *Software for algebraic geometry*, volume 148 of *IMA Vol. Math. Appl.*, pages 1–14. Springer, New York, 2008.
- [13] A. Belloni and R. M. Freund. A geometric analysis of Renegar’s condition number, and its interplay with conic curvature. *Math. Program.*, 119(1, Ser. A):95–107, 2009.
- [14] A. Belloni, R.M. Freund, and S. Vempala. An efficient rescaled perceptron algorithm for conic systems. *Math. Oper. Res.*, 34:621–641, 2009.
- [15] C. Beltrán. A continuation method to solve polynomial systems and its complexity. *Numer. Math.*, 117(1):89–113, 2011.
- [16] C. Beltrán, J.-P. Dedieu, G. Malajovich, and M. Shub. Convexity properties of the condition number II. *SIAM J. Matrix Anal. Appl.*, 33(3):905–939, 2012.
- [17] C. Beltrán, J.-P. Dedieu, G. Malajovich, and M. Shub. Convexity properties of the condition number. *SIAM J. Matrix Anal. Appl.*, 31(3):1491–1506, 2009.
- [18] C. Beltrán and A. Leykin. Certified numerical homotopy tracking. *Experimental Mathematics*, 21(1):69–83, 2012.
- [19] C. Beltrán and L. M. Pardo. Estimates on the distribution of the condition number of singular matrices. *Found. Comput. Math.*, 7(1):87–134, 2007.
- [20] C. Beltrán and L.M. Pardo. On the complexity of non universal polynomial equation solving: old and new results. In *Foundations of computational mathematics, Santander 2005*, volume 331 of *London Math. Soc. Lecture Note Ser.*, pages 1–35. Cambridge Univ. Press, Cambridge, 2006.
- [21] C. Beltrán and L.M. Pardo. On Smale’s 17 problem: a probabilistic positive solution. *Found. Comput. Math.*, 8:1–43, 2008.
- [22] C. Beltrán and L.M. Pardo. Smale’s 17th problem: average polynomial time to compute affine and projective solutions. *J. Amer. Math. Soc.*, 22(2):363–385, 2009.
- [23] C. Beltrán and L.M. Pardo. Fast linear homotopy to find approximate zeros of polynomial systems. *Found. Comput. Math.*, 11(1):95–129, 2011.
- [24] C. Beltrán and M. Shub. Complexity of Bézout’s Theorem VII: distance estimates in the condition metric. *Found. Comput. Math.*, 9:179–195, 2009.

- [25] C. Beltrán and M. Shub. On the geometry and topology of the solution variety for polynomial system solving. *Found. Comput. Math.*, 12(6): 719–763, 2012.
- [26] Å. Björck. Component-wise perturbation analysis and error bounds for linear least squares solutions. *BIT*, 31(2):238–244, 1991.
- [27] L. Blum. Lectures on a theory of computation and complexity over the reals (or an arbitrary ring). In E. Jen, editor, *Lectures in the Sciences of Complexity II*, pages 1–47. Addison-Wesley, 1990.
- [28] L. Blum, F. Cucker, M. Shub, and S. Smale. *Complexity and real computation*. Springer-Verlag, New York, 1998. With a foreword by R.M. Karp.
- [29] L. Blum and M. Shub. Evaluating rational functions: infinite precision is finite cost and tractable on average. *SIAM J. Comput.*, 15(2):384–398, 1986.
- [30] K.-H. Borgwardt. The average number of pivot steps required by the simplex-method is polynomial. *Z. Oper. Res. Ser.* 26(5):157–177, 1982.
- [31] Comte de Buffon, G.-L. Leclerc. Essai d’arithmétique morale. In *Supplément à l’Histoire naturelle*, volume 4, pages 46–148. Imprimerie Royale, Paris, 1777.
- [32] P. Bürgisser. Condition of intersecting a projective variety with a varying linear subspace. *SIAM Journal on Applied Algebra and Geometry*, 2017. To appear, arXiv:1510.04142v2.
- [33] P. Bürgisser and F. Cucker. Smoothed analysis of Moore-Penrose inversion. *SIAM J. Matrix Anal. Appl.*, 31(5):2769–2783, 2010.
- [34] P. Bürgisser and F. Cucker. On a problem posed by Steve Smale. *Annals of Mathematics*, 174:1785–1836, 2011.
- [35] P. Bürgisser and F. Cucker. *Condition: The geometry of numerical algorithms*, volume 349 of *Grundlehren der Mathematischen Wissenschaften*. Springer, Heidelberg, 2013.
- [36] P. Bürgisser, F. Cucker, and E.R. Cardozo. On the condition of the zeros of characteristic polynomials. arXiv:1510.04419.
- [37] P. Bürgisser, F. Cucker, and M. Lotz. Smoothed analysis of complex conic condition numbers. *J. Math. Pures Appl. (9)*, 86(4):293–309, 2006.
- [38] P. Bürgisser, F. Cucker, and M. Lotz. The probability that a slightly perturbed numerical analysis problem is difficult. *Mathematics of Computation*, 77:1559–1583, 2008.
- [39] P. Bürgisser, F. Cucker, and M. Lotz. Coverage processes on spheres and condition numbers for linear programming. *Annals of Probability*, 38:570–604, 2010.
- [40] Z.-Z. Cheng and J.J. Dongarra. Condition numbers of Gaussian random matrices. *SIAM J. Matrix Anal. Appl.*, 27:603–620, 2005.
- [41] S. Chern. On the kinematic formula in integral geometry. *J. Math. Mech.*, 16:101–118, 1966.
- [42] D. Cheung and F. Cucker. Smoothed analysis of componentwise condition numbers for sparse matrices. *IMA J. Numer. Anal.*, 35(1):74–88, 2015.
- [43] D. Cheung and F. Cucker. A new condition number for linear programming. *Math. Program.*, 91(1, Ser. A):163–174, 2001.
- [44] D. Cheung and F. Cucker. Probabilistic analysis of condition numbers for linear programming. *Journal of Optimization Theory and Applications*, 114:55–67, 2002.
- [45] D. Cheung and F. Cucker. Solving linear programs with finite precision: I. Condition numbers and random programs. *Math. Program.*, 99:175–196, 2004.
- [46] D. Cheung and F. Cucker. Solving linear programs with finite precision: II. Algorithms. *Journal of Complexity*, 22:305–335, 2006.
- [47] D. Cheung and F. Cucker. Componentwise condition numbers of random sparse matrices. *SIAM J. Matrix Anal. Appl.*, 31:721–731, 2009.
- [48] D. Cheung and F. Cucker. On the average condition of random linear programs. *SIAM J. Optim.*, 23(2): 799–810, 2013.
- [49] D. Cheung, F. Cucker, and R. Hauser. Tail decay and moment estimates of a condition number for random linear conic systems. *SIAM J. Optim.*, 15(4):1237–1261 (electronic), 2005.
- [50] D. Cheung, F. Cucker, and J. Peña. Unifying condition numbers for linear programming. *Math. Oper. Res.*, 28(4):609–624, 2003.

- [51] D. Cheung, F. Cucker, and Ye. Y. Linear programming and condition numbers under the real number computation model. In Ph. Ciarlet and F. Cucker, editors, *Handbook of Numerical Analysis*, volume XI, pages 141–207. North-Holland, 2003.
- [52] F. Cucker. Approximate zeros and condition numbers. *J. of Complexity*, 15:214–226, 1999.
- [53] F. Cucker, H. Diao, and Y. Wei. On mixed and componentwise condition numbers for Moore-Penrose inverse and linear least squares problems. *Mathematics of Computation*, 76:947–963, 2007.
- [54] F. Cucker and J. Peña. A primal-dual algorithm for solving polyhedral conic systems with a finite-precision machine. *SIAM J. Optim.*, 12(2):522–554 (electronic), 2001/02.
- [55] F. Cucker and S. Smale. Complexity estimates depending on condition and round-off error. *Journal of the ACM*, 46:113–184, 1999.
- [56] F. Cucker and M. Wschebor. On the expected condition number of linear programming problems. *Numer. Math.*, 94:419–478, 2003.
- [57] J.-P. Dedieu. *Points fixes, zéros et la méthode de Newton*, volume 54 of *Mathématiques & Applications (Berlin) [Mathematics & Applications]*. Springer, Berlin, 2006. With a preface by Steve Smale.
- [58] J.-P. Dedieu, G. Malajovich, and M. Shub. On the curvature of the central path of linear programming theory. *Found. Comput. Math.*, 5(2):145–171, 2005.
- [59] J.-P. Dedieu, P. Priouret, and G. Malajovich. Newton’s method on Riemannian manifolds: convariant alpha theory. *IMA J. Numer. Anal.*, 23(3):395–419, 2003.
- [60] J.P. Dedieu. Condition number analysis for sparse polynomial systems. In *Foundations of computational mathematics (Rio de Janeiro, 1997)*, pages 75–101. Springer, Berlin, 1997.
- [61] J.W. Demmel. On condition numbers and the distance to the nearest ill-posed problem. *Numer. Math.*, 51:251–289, 1987.
- [62] J.W. Demmel. The probability that a numerical analysis problem is difficult. *Math. Comp.*, 50:449–480, 1988.
- [63] J.W. Demmel. *Applied Numerical Linear Algebra*. SIAM, 1997.
- [64] J. Dunagan, D.A. Spielman, and S.-H. Teng. Smoothed analysis of condition numbers and complexity implications for linear programming. *Math. Program.*, 126(2, Ser. A):315–350, 2011.
- [65] C. Eckart and G. Young. The approximation of one matrix by another of lower rank. *Psychometrika*, 1(3):211–218, 1936.
- [66] A. Edelman. Eigenvalues and condition numbers of random matrices. *SIAM J. Matrix Anal. Appl.*, 9(4):543–560, 1988.
- [67] A. Edelman. On the distribution of a scaled condition number. *Math. Comp.*, 58(197):185–190, 1992.
- [68] M. Epelman and R.M. Freund. A new condition measure, preconditioners, and relations between different measures of conditioning for conic linear systems. *SIAM J. Optim.*, 12(3):627–655 (electronic), 2002.
- [69] H. Federer. Curvature measures. *Trans. Amer. Math. Soc.*, 93:418–491, 1959.
- [70] R. Fletcher. Expected conditioning. *IMA J. Numer. Anal.*, 5(3):247–273, 1985.
- [71] R.A. Fisher. The sampling distribution of some statistics obtained from nonlinear equations. *Ann. Eugenics*, 9:238–249, 1939.
- [72] R.M. Freund and J.R. Vera. Condition-based complexity of convex optimization in conic linear form via the ellipsoid algorithm. *SIAM J. Optim.*, 10(1):155–176 (electronic), 1999.
- [73] R.M. Freund and J.R. Vera. Some characterizations and properties of the “distance to ill-posedness” and the condition measure of a conic linear system. *Math. Program.*, 86:225–260, 1999.
- [74] F. Gao, D. Hug, and R. Schneider. Intrinsic volumes and polar sets in spherical space. *Math. Notae*, 41:159–176 (2003), 2001/02. Homage to Luis Santaló. Vol. 1 (Spanish).
- [75] S. Geman. A limit theorem for the norm of random matrices. *Ann. Probab.*, 8(2):252–261, 1980.
- [76] A.J. Geurts. A contribution to the theory of condition. *Numer. Math.*, 39:85–96, 1982.
- [77] E.N. Gilbert. The probability of covering a sphere with N circular caps. *Biometrika*, 52:323–330, 1965.
- [78] S. Glasauer. Integral geometry of spherically convex bodies. *Diss. Summ. Math.*, 1(1-2):219–226, 1996.

- [79] J.-L. Goffin. The relaxation method for solving systems of linear inequalities. *Math. Oper. Res.*, 5(3):388–414, 1980.
- [80] I. Gohberg and I. Koltracht. Mixed, componentwise, and structured condition numbers. *SIAM J. Matrix Anal. Appl.*, 14:688–704, 1993.
- [81] H.H. Goldstine and J. von Neumann. Numerical inverting matrices of high order, II. *Proc. Amer. Math. Soc.*, 2:188–202, 1951.
- [82] G.H. Golub and C.F. Van Loan. *Matrix computations*. Johns Hopkins Studies in the Mathematical Sciences. Johns Hopkins University Press, Baltimore, MD, fourth edition, 2013.
- [83] M. Grötschel, L. Lovász, and A. Schrijver. *Geometric algorithms and combinatorial optimization*, volume 2 of *Algorithms and Combinatorics: Study and Research Texts*. Springer-Verlag, Berlin, 1988.
- [84] M.R. Hestenes and E. Stiefel. Methods of conjugate gradients for solving linear systems. *J. Research Nat. Bur. Standards*, 49:409–436 (1953), 1952.
- [85] N.J. Higham. *Accuracy and stability of numerical algorithms*. Society for Industrial and Applied Mathematics (SIAM), Philadelphia, PA, second edition, 2002.
- [86] H. Hotelling. Some new methods in matrix calculation. *Ann. Math. Statistics*, 14:1–34, 1943.
- [87] R. Howard. The kinematic formula in Riemannian homogeneous spaces. *Mem. Amer. Math. Soc.*, 106(509):vi+69, 1993.
- [88] P.L. Hsu. On the distribution of roots of certain determinantal equations. *Ann. Eugenics*, 9:250–258, 1939.
- [89] S. Janson. Random coverings in several dimensions. *Acta Math.*, 156:83–118, 1986.
- [90] W. Kahan. Numerical linear algebra. *Canad. Math. Bull.*, 9:757–801, 1966.
- [91] L.V. Kantorovich. *On Newton's method*, volume 28 of *Trudy Mat. Inst. Steklov.*, pages 104–144. Acad. Sci. USSR, Moscow–Leningrad, 1949. In Russian.
- [92] N. Karmarkar. A new polynomial time algorithm for linear programming. *Combinatorica*, 4:373–395, 1984.
- [93] M. Karow, D. Kressner, and F. Tisseur. Structured eigenvalue condition numbers. *SIAM J. Matrix Anal. Appl.*, 28(4):1052–1068 (electronic), 2006.
- [94] L.G. Khachiyan. A polynomial algorithm in linear programming. *Dokl. Akad. Nauk SSSR*, 244:1093–1096, 1979. (In Russian, English translation in *Soviet Math. Dokl.*, 20:191–194, 1979.)
- [95] D.A. Klain and G.-C. Rota. *Introduction to geometric probability*. Lezioni Lincee. [Lincei Lectures]. Cambridge University Press, Cambridge, 1997.
- [96] E. Kostlan. On the distribution of the roots of random polynomials. In M. Hirsch, J.E. Marsden, and M. Shub, editors, *From Topology to Computation: Proceedings of the Smalefest*, pages 419–431. Springer-Verlag, 1993.
- [97] P. Lairez. A Deterministic Algorithm to Compute Approximate Roots of Polynomial Systems in Polynomial Average Time. To appear in *Found. Comput. Math.*
- [98] M. Ledoux and M. Talagrand. *Probability in Banach spaces*, volume 23 of *Ergebnisse der Mathematik und ihrer Grenzgebiete (3) [Results in Mathematics and Related Areas (3)]*. Springer-Verlag, Berlin, 1991. Isoperimetry and processes.
- [99] T.Y. Li. Numerical solution of polynomial systems by homotopy continuation methods. In Ph. Ciarlet and F. Cucker, editors, *Handbook of Numerical Analysis*, volume XI, pages 209–304. North-Holland, Amsterdam, 2003.
- [100] J. De Loera, B. Sturmfels, and C. Vinzant. The central curve of linear programming. *Found. Comput. Math.*, 12(4): 509–540, 2012.
- [101] M. Lotz. On the volume of tubular neighborhoods of real algebraic varieties. *Proc. Amer. Math. Soc.*, 143(5): 1875–1889, 2015.
- [102] G. Malajovich and J.M. Rojas. High probability analysis of the condition number of sparse polynomial systems. *Theoret. Comput. Sci.*, 315(2-3):524–555, 2004.
- [103] R.E. Miles. The asymptotic values of certain coverage probabilities. *Biometrika*, 56:661–680, 1969.
- [104] T. Motzkin and I.Y. Schönberg. The relaxation method for linear inequalities. *Canadian Journal of Mathematics*, 6:393–404, 1954.

- [105] R.J. Muirhead. *Aspects of multivariate statistical theory*. John Wiley & Sons Inc., New York, 1982. Wiley Series in Probability and Mathematical Statistics.
- [106] Y. Nesterov and A. Nemirovsky. *Interior-Point Polynomial Algorithms in Convex Programming*. SIAM, 1994.
- [107] J. von Neumann and H.H. Goldstine. Numerical inverting matrices of high order. *Bull. Amer. Math. Soc.*, 53:1021–1099, 1947.
- [108] M. Nunez and R.M. Freund. Condition measures and properties of the central trajectory of a linear program. *Math. Program.*, 83:1–28, 1998.
- [109] V.Y. Pan. Optimal and nearly optimal algorithms for approximating polynomial zeros. *Comput. Math. Appl.*, 31(12):97–138, 1996.
- [110] V.Y. Pan. Solving a polynomial equation: some history and recent progress. *SIAM Rev.*, 39(2):187–220, 1997.
- [111] J. Peña and J. Renegar. Computing approximate solutions for conic systems of constraints. *Math. Program.*, 87:351–383, 2000.
- [112] J. Peña. Understanding the geometry on infeasible perturbations of a conic linear system. *SIAM J. Optim.*, 10:534–550, 2000.
- [113] J. Peña. A characterization of the distance to infeasibility under block-structured perturbations. *Linear Algebra Appl.*, 370:193–216, 2003.
- [114] J. Renegar. On the efficiency of Newton’s method in approximating all zeros of a system of complex polynomials. *Math. Oper. Res.*, 12(1):121–148, 1987.
- [115] J. Renegar. On the worst-case arithmetic complexity of approximating zeros of systems of polynomials. *SIAM J. Comput.*, 18:350–370, 1989.
- [116] J. Renegar. Is it possible to know a problem instance is ill-posed? *Journal of Complexity*, 10:1–56, 1994.
- [117] J. Renegar. Some perturbation theory for linear programming. *Math. Program.*, 65:73–91, 1994.
- [118] J. Renegar. Incorporating condition measures into the complexity theory of linear programming. *SIAM J. Optim.*, 5:506–524, 1995.
- [119] J. Renegar. Linear programming, complexity theory and elementary functional analysis. *Math. Program.*, 70:279–351, 1995.
- [120] J. Renegar. *A Mathematical View of Interior-Point Methods in Convex Optimization*. SIAM, 2000.
- [121] J.R. Rice. A theory of condition. *SIAM J. Numer. Anal.*, 3:217–232, 1966.
- [122] S.N. Roy. p-statistics or some generalizations in the analysis of variance appropriate to multivariate problems. *Sankhya*, 4:381–396, 1939.
- [123] M. Rudelson and R. Vershynin. Smallest singular value of a random rectangular matrix. *Comm. Pure Appl. Math.*, 62(12):1707–1739, 2009.
- [124] S.M. Rump. Structured perturbations part I: normwise distances. *SIAM J. Matrix Anal. Appl.*, 25:1–30, 2003.
- [125] S.M. Rump. Structured perturbations part II: componentwise distances. *SIAM J. Matrix Anal. Appl.*, 25:31–56, 2003.
- [126] A. Sankar, D.A. Spielman, and S.-H. Teng. Smoothed analysis of the condition numbers and growth factors of matrices. *SIAM J. Matrix Anal. Appl.*, 28(2):446–476 (electronic), 2006.
- [127] L.A. Santaló. *Integral geometry and geometric probability*. Addison-Wesley Publishing Co., Reading, Mass.-London-Amsterdam, 1976. With a foreword by Mark Kac, Encyclopedia of Mathematics and its Applications, Vol. 1.
- [128] R. Schneider and W. Weil. *Stochastic and integral geometry*. Probability and its Applications (New York). Springer-Verlag, Berlin, 2008.
- [129] M. Shub. Some remarks on Bezout’s theorem and complexity theory. In *From Topology to Computation: Proceedings of the Smalefest (Berkeley, CA, 1990)*, pages 443–455. Springer, New York, 1993.
- [130] M. Shub. Complexity of Bézout’s Theorem VI: geodesics in the condition (number) metric. *Found. Comput. Math.*, 9:171–178, 2009.
- [131] M. Shub and S. Smale. Computational complexity. On the geometry of polynomials and a theory of cost. I. *Ann. Sci. École Norm. Sup. (4)*, 18(1):107–142, 1985.

- [132] M. Shub and S. Smale. Computational complexity: on the geometry of polynomials and a theory of cost. II. *SIAM J. Comput.*, 15(1):145–161, 1986.
- [133] M. Shub and S. Smale. Complexity of Bézout’s Theorem I: geometric aspects. *J. Amer. Math. Soc.*, 6:459–501, 1993.
- [134] M. Shub and S. Smale. Complexity of Bézout’s Theorem II: volumes and probabilities. In F. Eyssette and A. Galligo, editors, *Computational Algebraic Geometry*, volume 109 of *Progress in Mathematics*, pages 267–285. Birkhäuser, 1993.
- [135] M. Shub and S. Smale. Complexity of Bézout’s Theorem III: condition number and packing. *Journal of Complexity*, 9:4–14, 1993.
- [136] M. Shub and S. Smale. Complexity of Bézout’s Theorem V: polynomial time. *Theoretical Computer Science*, 133:141–164, 1994.
- [137] M. Shub and S. Smale. Complexity of Bézout’s Theorem IV: probability of success; extensions. *SIAM J. of Numer. Anal.*, 33:128–148, 1996.
- [138] J.W. Silverstein. The smallest eigenvalue of a large-dimensional Wishart matrix. *Ann. Probab.*, 13(4):1364–1368, 1985.
- [139] S. Smale. The fundamental theorem of algebra and complexity theory. *Bull. Amer. Math. Soc.*, 4:1–36, 1981.
- [140] S. Smale. On the average number of steps of the simplex method of linear programming. *Math. Programming*, 27(3):241–262, 1983.
- [141] S. Smale. Newton’s method estimates from data at one point. In R. Ewing, K. Gross, and C. Martin, editors, *The Merging of Disciplines: New Directions in Pure, Applied, and Computational Mathematics*. Springer-Verlag, 1986.
- [142] S. Smale. Complexity theory and numerical analysis. In A. Iserles, editor, *Acta Numerica*, pages 523–551. Cambridge University Press, 1997.
- [143] S. Smale. Mathematical problems for the next century. *Math. Intelligencer*, 20(2):7–15, 1998.
- [144] A.J. Sommese and C.W. Wampler, II. *The numerical solution of systems of polynomials*. World Scientific Publishing Co. Pte. Ltd., Hackensack, NJ, 2005. Arising in engineering and science.
- [145] D.A. Spielman and S.-H. Teng. Smoothed analysis of algorithms: why the simplex algorithm usually takes polynomial time. In *Proceedings of the Thirty-Third Annual ACM Symposium on Theory of Computing*, pages 296–305 (electronic), New York, 2001. ACM.
- [146] D.A. Spielman and S.-H. Teng. Smoothed analysis of algorithms. In *Proceedings of the International Congress of Mathematicians*, volume I, pages 597–606, 2002.
- [147] D.A. Spielman and S.-H. Teng. Smoothed analysis: Why the simplex algorithm usually takes polynomial time. *Journal of the ACM*, 51(3):385–463, 2004.
- [148] J. Steiner. Über parallele Flächen. *Monatsber. preuss. Akad. Wiss.*, pages 114–118, 1840.
- [149] G.W. Stewart. On the perturbation of pseudo-inverses, projections and linear least squares problems. *SIAM Rev.*, 19(4):634–662, 1977.
- [150] G.W. Stewart. Stochastic perturbation theory. *SIAM Rev.*, 32(4):579–610, 1990.
- [151] G.W. Stewart and J.-G. Sun. *Matrix perturbation theory*. Computer Science and Scientific Computing. Academic Press Inc., Boston, MA, 1990.
- [152] T. Tao and V. Vu. Smooth analysis of the condition number and the least singular value. *Math. Comp.*, 79(272):2333–2352, 2010.
- [153] L.N. Trefethen and R.S. Schreiber. Average-case stability of Gaussian elimination. *SIAM J. Matrix Anal. Appl.*, 11:335–360, 1990.
- [154] A.M. Turing. Rounding-off errors in matrix processes. *Quart. J. Mech. Appl. Math.*, 1:287–308, 1948.
- [155] S.A. Vavasis and Y. Ye. Condition numbers for polyhedra with real number data. *Oper. Res. Lett.*, 17:209–214, 1995.
- [156] S.A. Vavasis and Y. Ye. A primal-dual interior point method whose running time depends only on the constraint matrix. *Math. Program.*, 74:79–120, 1996.
- [157] D. Viswanath and L.N. Trefethen. Condition numbers of random triangular matrices. *SIAM J. Matrix Anal. Appl.*, 19:564–581, 1998.

- [158] P.-Å. Wedin. Perturbation theory for pseudo-inverses. *Nordisk Tidskr. Informationsbehandling (BIT)*, 13:217–232, 1973.
- [159] N. Weiss, G.W. Wasilkowski, H. Woźniakowski, and M. Shub. Average condition number for solving linear equations. *Linear Algebra Appl.*, 83:79–102, 1986.
- [160] J.G. Wendel. A problem in geometric probability. *Math. Scand.*, 11:109–111, 1962.
- [161] H. Weyl. On the volume of tubes. *Amer. J. Math.*, 61(2):461–472, 1939.
- [162] E. Wigner. Random matrices in physics. *SIAM Rev.*, 9:1–23, 1967.
- [163] J.H. Wilkinson. Error analysis of direct methods of matrix inversion. *J. Assoc. Comput. Mach.*, 8:281–330, 1961.
- [164] J.H. Wilkinson. *Rounding Errors in Algebraic Processes*. Prentice Hall, 1963.
- [165] J.H. Wilkinson. *The Algebraic Eigenvalue Problem*. Clarendon Press, 1965.
- [166] J.H. Wilkinson. Modern error analysis. *SIAM Review*, 13:548–568, 1971.
- [167] J.H. Wilkinson. Note on matrices with a very ill-conditioned eigenproblem. *Numer. Math.*, 19:176–178, 1972.
- [168] J. Wishart. The generalized product moment distribution in samples from a normal multivariate population. *Biometrika*, 20A(272):32–43, 1928.
- [169] R. Wongkew. Volumes of tubular neighbourhoods of real algebraic varieties. *Pacific J. Math.*, 159(1):177–184, 1993.
- [170] H. Woźniakowski. Numerical stability for solving nonlinear equations. *Numer. Math.*, 27(4):373–390, 1976/77.
- [171] M.H. Wright. The interior-point revolution in optimization: history, recent developments, and lasting consequences. *Bull. Amer. Math. Soc. (N.S.)*, 42(1):39–56, 2005.
- [172] S. Wright. *Primal-Dual Interior-Point Methods*. SIAM, 1997.
- [173] M. Wschebor. Smoothed analysis of $\kappa(A)$. *Journal of Complexity*, 20(1):97–107, 2004.
- [174] Y. Ye. Toward probabilistic analysis of interior-point algorithms for linear programming. *Math. of Oper. Res.*, 19:38–52, 1994.
- [175] T.J. Ypma. Historical development of the Newton-Raphson method. *SIAM Rev.*, 37(4):531–551, 1995.

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