

LINEAR PROGRAMMING AND CIRCUIT IMBALANCES

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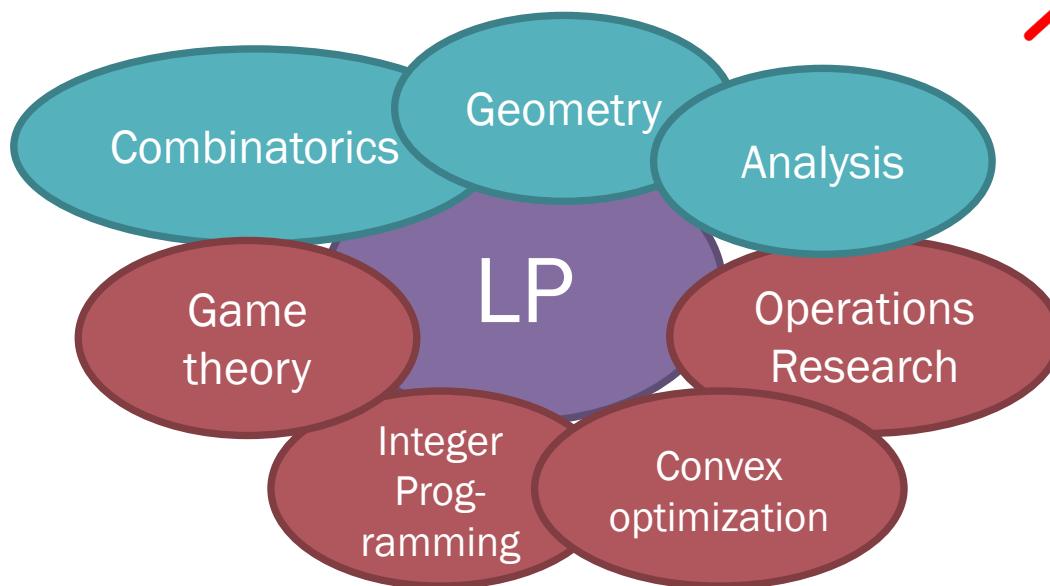
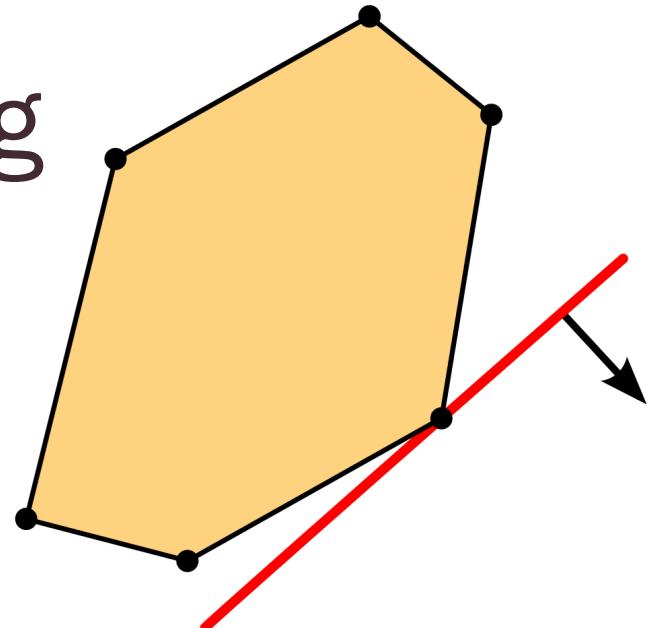
Slides available at
<https://personal.lse.ac.uk/veghi/ipco>

Linear programming

$$\min c^T x$$

$$Ax = b$$

$$x \geq 0$$



Facets of linear programming

Discrete

- Basic solutions
- Polyhedral combinatorics
- Exact solution



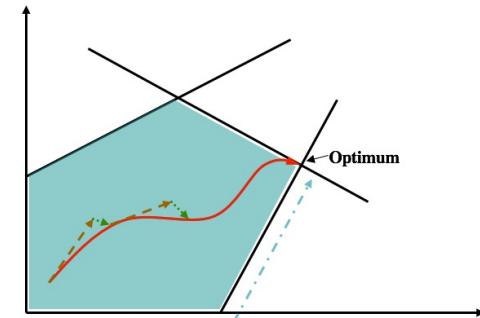
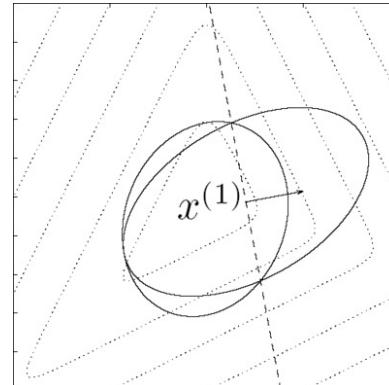
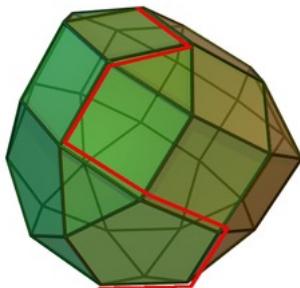
Continuous

- Continuous solutions
- Convex program
- Approximate solution

Linear programming algorithms

$$\begin{aligned} & \min c^T x \\ & Ax = b \\ & x \geq 0 \end{aligned}$$

- n variables, m constraints
- L : total bit-complexity of the rational input (A, b, c)
- Simplex method: Dantzig, 1947
- Weakly polynomial algorithms: $\text{poly}(L)$ running time
 - Ellipsoid method: Khachiyan, 1979
 - Interior point method: Karmarkar, 1984



Weakly vs strongly polynomial algorithms for LP

$$\begin{aligned} \min c^T x \\ Ax = b \\ x \geq 0 \end{aligned}$$

- n variables, m constraints, total encoding L .
- Strongly polynomial algorithm:
 - $\text{poly}(n, m)$ elementary arithmetic operations $(+, -, \times, \div, \geq)$, independent of L .
 - PSPACE: The bit-length of numbers during the algorithm remain polynomially bounded in the size of the input.
 - Can also be defined in the real model of computation

Is there a strongly polynomial
algorithm for Linear
Programming?



Smale's 9th question

Strongly polynomial algorithms for some classes of Linear Programs

- Systems of linear inequalities with at most two nonzero variables per inequality: Megiddo '83
- Network flow problems
 - Maximum flow: Edmonds-Karp-Dinitz '70-72, ...
 - Min-cost flow: Tardos '85, Fujishige '86, Goldberg-Tarjan '89, Orlin '93, ...
 - Generalized flow: V '17, Olver-V '20
- Discounted Markov Decision Processes: Ye '05, Ye '11, ...

Dependence on the constraint matrix only

$$\min_x c^T x, \quad Ax = b \quad x \geq 0$$

- Algorithms with running time dependent only on A , but not on b and c .

- Combinatorial LP's: integer matrix $A \in \mathbb{Z}^{m \times n}$.

$$\Delta_A = \max\{|\det(B)| : B \text{ submatrix of } A\}$$

Tardos '86: $\text{poly}(n, m, \log \Delta_A)$ black box LP algorithm

- Layered-least-squares (LLS) Interior Point Method

Vavasis-Ye '96: $\text{poly}(n, m, \log \bar{\chi}_A)$ LP algorithm in the real model of computation

$\bar{\chi}_A$: condition number

- Dadush-Huiberts-Natura-V '20: $\text{poly}(n, m, \log \bar{\chi}_A^*)$

$\bar{\chi}_A^*$: optimized version of $\bar{\chi}_A$

Outline of the lectures

1. Tardos's algorithm for min-cost flows
2. The circuit imbalance measure κ_A and the condition measure $\bar{\chi}_A$
3. Solving LPs: from approximate to exact
 4. Optimizing circuit imbalances
 5. Interior point methods: basic concepts
 6. Layered-least-squares interior point methods

- Dadush-Huiberts-Natura-V '20: A scaling-invariant algorithm for linear programming whose running time depends only on the constraint matrix
- Dadush-Natura-V '20: Revisiting Tardos's framework for linear programming: Faster exact solutions using approximate solvers



Part 1

Tardos's algorithm for min-cost flows *circuits, proximity, and variable fixing*



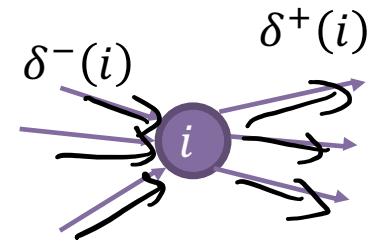
The minimum-cost flow problem

- Directed graph $G = (V, E)$, node demands $b: V \rightarrow \mathbb{R}$ with $b(V) = 0$, costs $c: E \rightarrow \mathbb{R}$.

$$\min \underbrace{c^\top x}_{\text{cost}}$$

$$\text{s. t. } \sum_{ji \in \delta^-(i)} x_{ji} - \sum_{ij \in \delta^+(i)} x_{ij} = b_i \quad \forall i \in V$$

$$x \geq 0$$

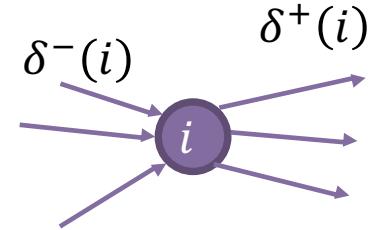


- Form with arc capacities can be reduced to this form.
- Constraint matrix is totally unimodular (TU)

		arcs
		ij
		nodes
	i	-1
	j	1

The minimum-cost flow problem: optimality

- Directed graph $G = (V, E)$, node demands $b: V \rightarrow \mathbb{R}$ with $b(V) = 0$, costs $c: E \rightarrow \mathbb{R}$.



$$\min c^\top x$$

$$\text{s. t. } \sum_{(j,i) \in \delta^-(i)} x_{ji} - \sum_{(i,j) \in \delta^+(i)} x_{ij} = b_i \quad \forall i \in V$$
$$x \geq 0$$

- Dual program:

$$\max b^\top \pi$$
$$\text{s.t. } \pi_j - \pi_i \leq c_{ij} \quad \forall ij \in E$$

- Optimality: $f_{ij} > 0 \implies \pi_j - \pi_i = c_{ij}$

Dual solutions: potentials

- Dual program: max cost feasible potential

$$\max b^T \pi$$

$$\text{s. t. } \pi_j - \pi_i \leq c_{ij} \quad \forall ij \in E$$

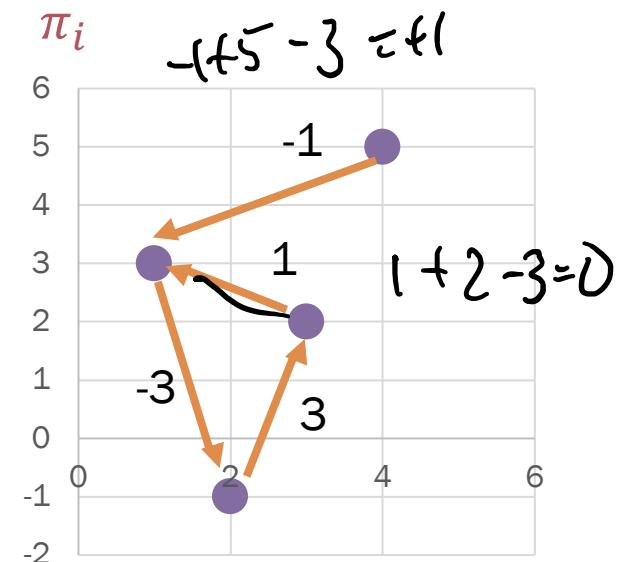
- Residual cost:

$$c_{ij}^\pi = c_{ij} + \pi_i - \pi_j \geq 0$$

- Residual graph:

$$E_f = E \cup \{(j, i) : f_{ij} > 0\}$$

$$c_{ji} = -c_{ij}$$



LEMMA: The primal feasible f is optimal \Leftrightarrow

$\exists \pi: c_{ij}^\pi \geq 0$ for all $(i, j) \in E$ and $c_{ij}^\pi = 0$ if $f_{ij} > 0$ \Leftrightarrow

$\exists \pi: c_{ij}^\pi \geq 0$ for all $(i, j) \in E_f$

Variable fixing by proximity

- If for some $(i, j) \in E$ we can show that $\underline{f}_{ij}^* = 0$ in every optimal solution, then we can remove (i, j) from the graph.
- Overall goal: in strongly polynomial number of steps, guarantee that we can infer this for at least one arc.

PROXIMITY THEOREM: Let $\tilde{\pi}$ be the optimal dual potential for costs \tilde{c} , and \underline{f}^* an optimal primal solution for the original costs c . Then,

$$c_{ij}^{\tilde{\pi}} > |V| \cdot \|c - \tilde{c}\|_{\infty} \Rightarrow \underline{f}_{ij}^* = 0$$

Circulations and cycle decompositions

- For the node-arc incidence matrix A , $\ker(A) \subseteq \mathbb{R}^E$ is the set of circulations:

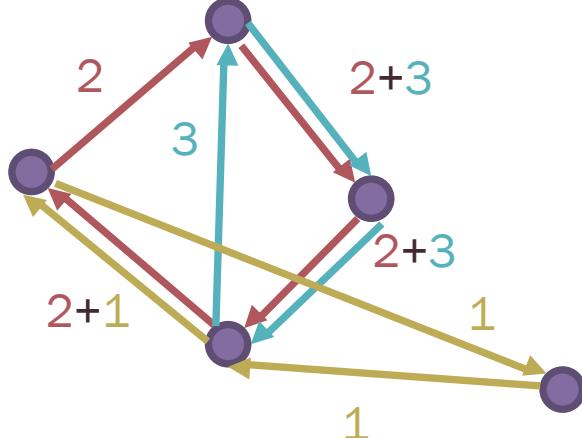
in-flow=out-flow

- LEMMA: every circulation $f \geq 0$ can be decomposed as

$$\underline{f} = \sum_i \lambda_i \chi_{C_i}, \quad \lambda_i \geq 0$$

for directed cycles C_i

*Can the 'ology
decomy.*



Circulations and cycle decompositions

- **LEMMA:** Let f and f' be two feasible flows for the same demand vector b . Then, we can write

$$f' = f + \sum_i \lambda_i \chi_{C_i}, \quad \lambda_i \geq 0$$

$$\underbrace{f' - f}_{\text{sign-consistent}}$$

for **sign-consistent** directed cycles C_i in \vec{E} :

- If $f'_{ij} > f_{ij}$ then cycles may only contain ij but not ji .
- If $f_{ij} > f'_{ij}$ then cycles may only contain ji but not ij .
- If $f_{ij} = f'_{ij}$ then no cycle contains ij or ji .

Every cycle is moving from f towards f' .

PROXIMITY THEOREM: Let $\tilde{\pi}$ be the optimal dual potential for costs \tilde{c} , and f^* an optimal primal solution for the original costs c . Then,

$$c_{ij}^{\tilde{\pi}} > |V| \cdot \underbrace{\|c - \tilde{c}\|_{\infty}}_{\epsilon} \Rightarrow f_{ij}^* = 0$$

PROOF:

$$\begin{aligned}
 & \text{Given } c \rightsquigarrow f^* \text{ opt} \\
 & \tilde{c} \rightsquigarrow (\tilde{f}, \tilde{\pi}) \\
 & \tilde{c}_{ij}^{\tilde{\pi}} > |V|\epsilon \\
 & \tilde{c}_{ij}^{\tilde{\pi}} > |V|\epsilon \quad (|V|-1)\epsilon > 0 \Rightarrow \tilde{f}_{ij}^* < 0 \\
 & \underline{f_{ij}^* > 0} \\
 & f^* = \tilde{f} + \sum_k \lambda_k X_{C_k} \\
 & \tilde{f} + \lambda_1 X_{C_1}, \tilde{f}^* - \lambda_1 X_{C_1} \text{ are feasible solns.} \\
 & c(C_1) \leq 0, \quad \tilde{c}(C_1) \leq \epsilon |C_1| \\
 & \tilde{c}(C_1) = \tilde{c}^{\tilde{\pi}}(C_1) = \sum_{e \in C_1} \tilde{c}_e^{\tilde{\pi}} \geq \tilde{c}_{ij}^{\tilde{\pi}} > (|V|-1)\epsilon \quad \square
 \end{aligned}$$

Rounding the costs

- Rescale c such that $\|c\|_\infty = |V|\sqrt{|E|}$
- Round costs as $\tilde{c}_{ij} = \lfloor c_{ij} \rfloor$
- For \tilde{c} we can find optimal primal and dual solutions in strongly polynomial time, e.g. the Out-of-Kilter method by Ford and Fulkerson 1962.
- For the optimal dual $\tilde{\pi}$, fix all arcs to 0 that have
$$c_{ij}^{\tilde{\pi}} > |V| > |V| \cdot \underbrace{\|c - \tilde{c}\|_\infty}_{\epsilon}$$
- QUESTION: Why would such an arc exist?

$$|c_{ij}^{\tilde{\pi}}| < 0.001\epsilon$$

Minimum-norm projections

- Residual cost:

$$c_{ij}^\pi = c_{ij} + \pi_i - \pi_j \geq 0$$

- The cost vectors

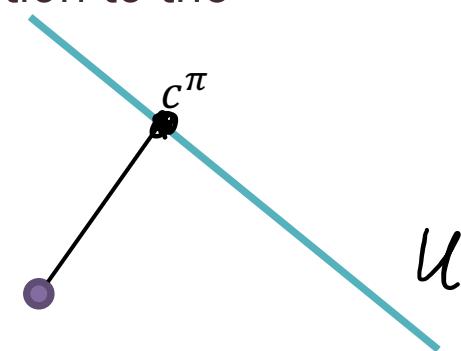
$$\underbrace{U = \{c^\pi : \pi \in \mathbb{R}^V\}}_{\text{form an affine subspace.}} \subset \mathbb{R}^E$$

- For any feasible flow f and any residual cost c^π ,

$$(c^\pi)^\top f = \underbrace{c^\top f}_{\text{ }} + \underbrace{b^\top \pi}_{\text{ }}$$

- Solving the problem for c and c^π is equivalent.
- If $0 \in U$, i.e. $\exists \pi : c^\pi \equiv 0$, then every feasible flow is optimal
- **IDEA:** Replace the input c by the min norm projection to the affine subspace U :

$$c^\pi = \arg \min_{\pi \in \mathbb{R}^V} \|c^\pi\|_2$$



Rounding the costs

- Assume c is chosen as a min norm projection:

$$\underbrace{\|c^\pi\|_2 \geq \|c\|_2}_{\forall \pi \in \mathbb{R}^V}$$

- Rescale c such that $\|c\|_\infty = |V| \sqrt{|E|}$
- Round costs as $\tilde{c}_{ij} = \lfloor c_{ij} \rfloor$
- For the optimal dual $\tilde{\pi}$, fix all arcs to 0 that have

$$c_{ij}^{\tilde{\pi}} > |V| > |V| \cdot \|c - \tilde{c}\|_\infty$$

- **LEMMA:** There exist at least one such arc.

PROOF:

$$\underbrace{\|c^{\tilde{\pi}}\|_\infty}_{\geq \frac{\|c^{\tilde{\pi}}\|_2}{\sqrt{|E|}}} \geq \frac{\|c\|_2}{\sqrt{|E|}} \geq \frac{\|c\|_\infty}{\sqrt{|E|}} = |V|$$

Also note that

$$\underbrace{c_{ij}^{\tilde{\pi}} \geq \tilde{c}_{ij}^{\tilde{\pi}} \geq 0}_{\text{arcs to 0}}$$

Summary of Tardos's algorithm

- Variable fixing based on proximity that can be shown by cycle decomposition.
- Replace the input cost by an equivalent min-cost projection
- Round to small integer costs \tilde{c}
- Find optimal dual $\tilde{\pi}$ for \tilde{c} with simple classical method
- Identify a variable $f_{ij}^* = 0$ as one where $c_{ij}^{\tilde{\pi}}$ is large and remove all such arcs.
- Iterate



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Part 2

The circuit imbalance measure κ_A
and the condition measure $\bar{\chi}_A$



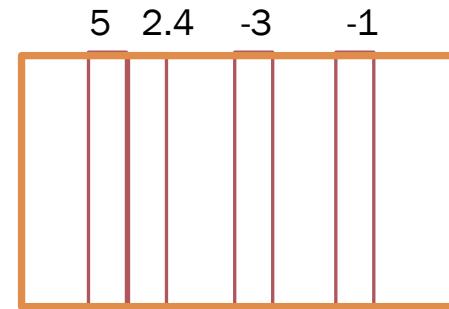
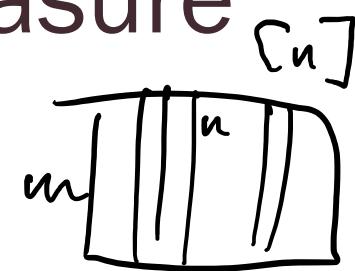
The circuit imbalance measure

- The matrix $A \in \mathbb{R}^{m \times n}$ defines a **linear matroid** on $[n] = \{1, 2, \dots, n\}$: a set $I \subseteq [n]$ is **independent** if the columns $\{a_i : i \in I\}$ are linearly independent.
- $C \subseteq [n]$ is a **circuit** if $\{a_i : i \in C\}$ is a linearly dependent set minimal for containment.
- For a circuit C , there exists a vector $g^C \in \mathbb{R}^C$ unique up to a scalar multiplier such that

$$\sum_{i \in C} g_i^C a_i = 0$$

- \mathcal{C}_A : set of all circuits.
- The **circuit imbalance measure** is defined as

$$\kappa_A = \max \left\{ \frac{|g_j^C|}{|g_i^C|} : C \in \mathcal{C}_A, i, j \in C \right\}$$

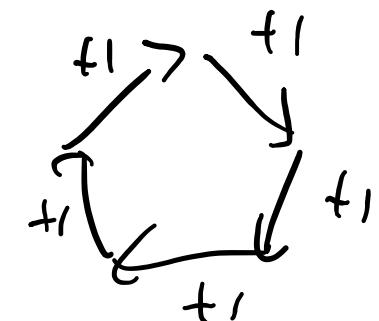


$$\frac{5}{|-1|} = 5$$

Properties of κ_A

$$\kappa_A = \max \left\{ \frac{|g_j^C|}{|g_i^C|} : C \in \mathcal{C}_A, i, j \in C \right\}$$

- This measure depends only on the linear subspace $W = \ker(A)$: if $\ker(A) = \ker(B)$ then $\kappa_A = \kappa_B$
- We will use $\kappa_W = \kappa_A$ for $W = \ker(A)$

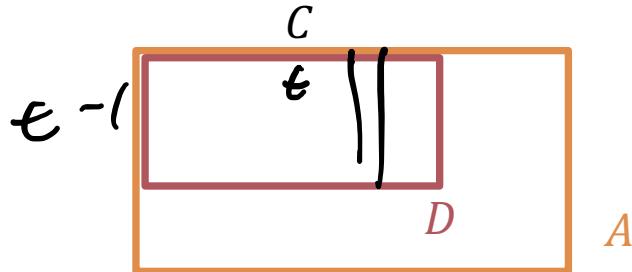


Connection to subdeterminants:

- For an integer matrix $A \in \mathbb{Z}^{m \times n}$,

$$\underline{\Delta_A} = \max \{ |\det(B)| : B \text{ submatrix of } A \}$$

- For a circuit $C \in \mathcal{C}_A$, with $|C| = t$ let $D = A_{J,C} \in \mathbb{R}^{(t-1) \times t}$ be a submatrix with linearly independent rows.



$D^{(i)} \in \mathbb{R}^{(t-1) \times (t-1)}$ remove the i -th column from D . By **Cramer's rule**

$$g^C = (\det(D^{(1)}), \det(D^{(2)}), \dots, \det(D^{(t)}))$$

Properties of κ_A

- **LEMMA:** For an integer matrix $A \in \mathbb{Z}^{m \times n}$,

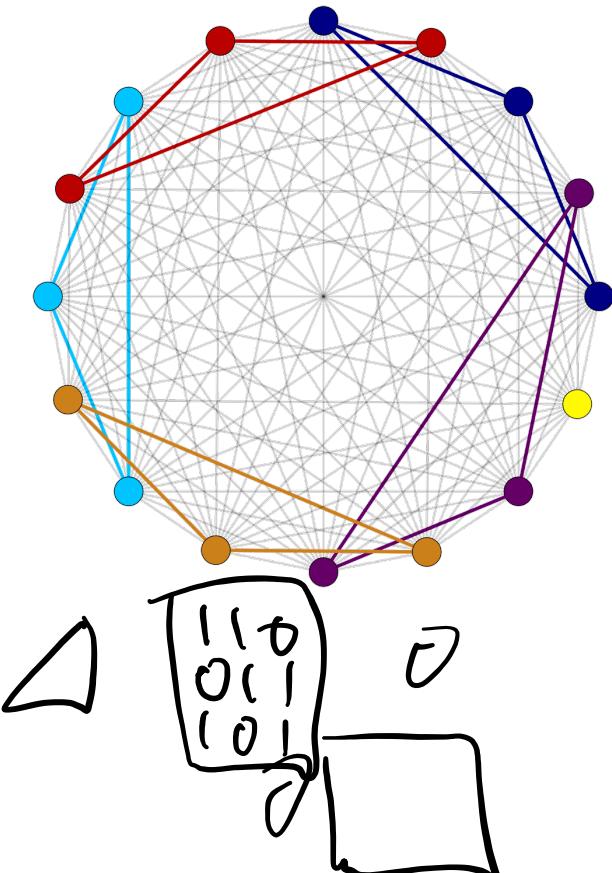
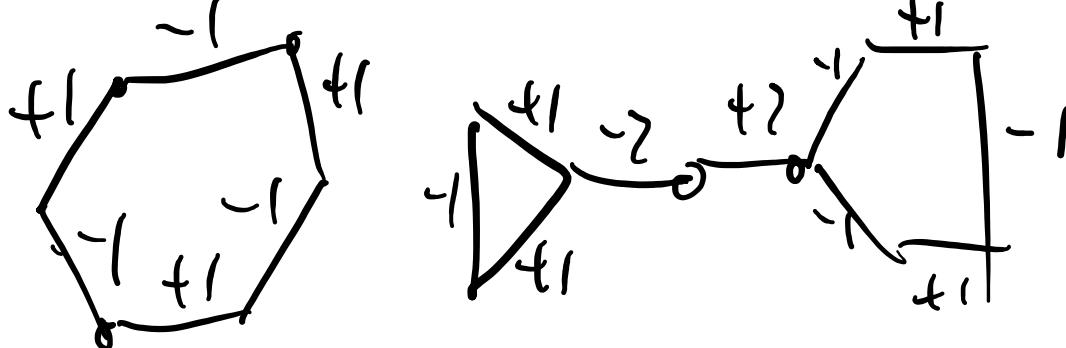
$$\kappa_A \leq \Delta_A$$

For a totally unimodular matrix A , $\kappa_A = 1$

- **EXERCISE:**

- If A is the node-edge incidence matrix of an undirected graph, then $\kappa_A \in \{1,2\}$
- For the incidence matrix of a complete undirected graph on n nodes,

$$\Delta_A \geq 2^{\left\lceil \frac{n}{3} \right\rceil}$$



Circuit imbalance and TU matrices

THEOREM (Cederbaum, 1958): If $A \in \mathbb{Z}^{m \times n}$ is a TU-matrix, then $\kappa_A = 1$. Conversely, if $\kappa_W = 1$ for a linear subspace $W \subset \mathbb{R}^n$ then there exists a TU-matrix A such that $W = \ker(A)$.

PROOF (Ekbatani & Natura):

$$W = \ker(A) \quad A' = A_B^{-1} A$$

basis B

$0, \pm 1$

$$\begin{array}{|c|c|c|} \hline 1 & 0 & M \\ \hline 0 & 1 & \\ \hline \end{array} \quad m \times n$$

$$M^{-1} A' \quad \begin{array}{|c|c|c|} \hline B & M^{-1} & 1 & 0 & 0 \\ \hline & & 1 & 0 & 1 \\ \hline \end{array}$$

$$1 = \det(M) \cdot \det(M^{-1}) \stackrel{M^{-1}}{\Rightarrow} \det(M) = \pm 1$$

Duality of circuit imbalances

THEOREM: For every linear subspace $W \subset \mathbb{R}^n$, we have

$$\kappa_W = \kappa_{W^\perp}$$

Circuits in optimization

- Appear in various LP algorithms directly or indirectly
- IPCO summer school 2020: Laura Sanità's lectures discussed *circuit augmentation* algorithms and *circuit diameter*
- Integer programming: κ has a natural integer variant that is related to Graver bases
- ...

The condition number $\bar{\chi}_A$

$$\bar{\chi}_A = \sup\{\|A^T(ADA^T)^{-1}AD\| : D \text{ is positive diagonal matrix}\}$$

- Measures the norm of *oblique* projections
- Introduced by Dikin 1967, Stewart 1989, Todd 1990
- **THEOREM (Vavasis-Ye 1996):** There exists a $\text{poly}(n, m, \log \bar{\chi}_A)$ LP algorithm for $\min c^T x, Ax = b, x \geq 0, A \in \mathbb{R}^{m \times n}$
- **LEMMA**
 - i. If A is an integer matrix with bit encoding length L , then $\bar{\chi}_A \leq 2^{O(L)}$
 - ii. $\bar{\chi}_A = \max\{\|B^{-1}A\| : B \text{ nonsingular } m \times m \text{ submatrix of } A\}$
 - iii. $\bar{\chi}_A$ only depends on the subspace $W = \ker(A)$
 - iv. $\bar{\chi}_W = \bar{\chi}_{W^\perp}$

The lifting operator

$$w = \ker(\lambda)$$

- For a linear subspace $W \subset \mathbb{R}^n$ and index set $I \subseteq [n]$, we let

$$\pi_I: \mathbb{R}^n \rightarrow \mathbb{R}^I$$

denote the coordinate projection, and

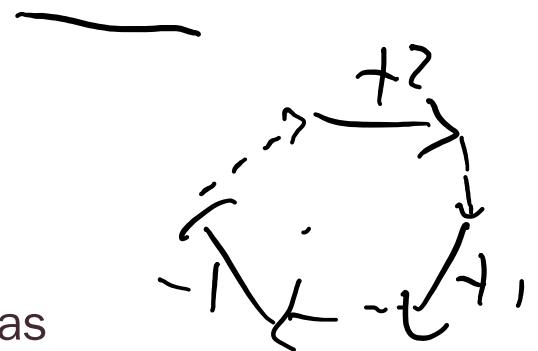
$$\pi_I(W) = \{x_I: x \in W\}$$

- The lifting operator $L_I^W: \mathbb{R}^I \rightarrow \mathbb{R}^n$ is defined as

$$L_I^W(z) = \arg \min_{\underline{\quad}} \{ \|x\|_2 : x \in W, x_I = z \}$$

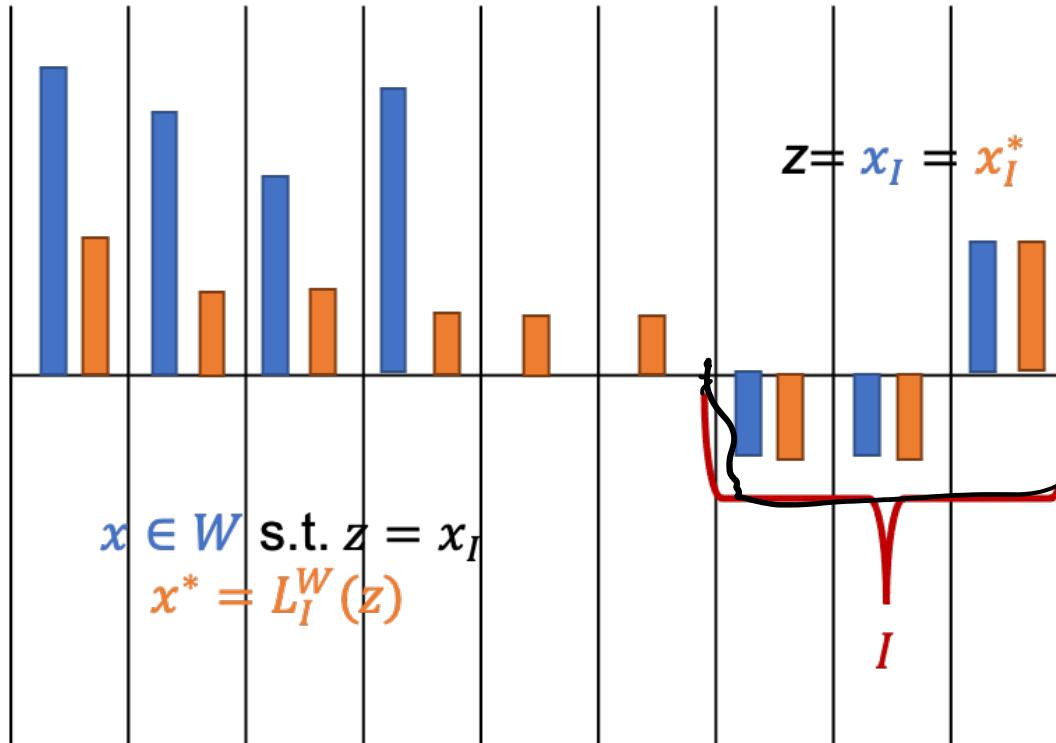
- This is a linear operator; we can efficiently compute a projection matrix $H \in \mathbb{R}^{n \times I}$ such that $L_I^W(z) = Hz$.
- LEMMA:**

$$\bar{\chi}_A = \max_{I \subseteq [n]} \|L_I^W\| = \max \left\{ \frac{\|L_I^W(z)\|_2}{\|z\|_2} : I \subseteq [n], z \in \pi_I(W) \setminus \{0\} \right\}$$



The lifting operator

$$L_I^W(z) = \arg \min\{\|x\|_2 : x \in W, x_I = z\}$$



The lifting operator

$$\ell_1 \rightarrow \ell_\infty$$

$$\forall j \notin I \quad \forall i \in C_i \cap \bar{I} \quad |g_j^{C_i}| \leq \sum_{\ell \in I} |g_\ell^{C_i}|$$

LEMMA:

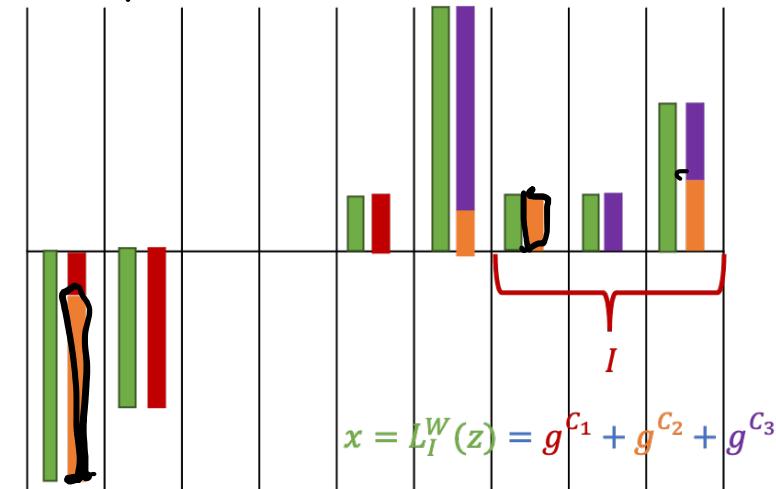
$$\kappa_A = \max \left\{ \frac{\|L_I^W(z)\|_\infty}{\|z\|_1} : I \subseteq [n], z \in \pi_I(W) \setminus \{0\} \right\}$$

PROOF: $\kappa_A \|L_I^W(z)\|_\infty \leq \kappa_A \|z\|_1$,

Carathéodory "x
 $x = \sum_i g^{C_i}$ sign-consistent comb.

$\forall i \quad C_i \cap \bar{I} \neq \emptyset$

$$x' = x - g \quad \forall i \in W \quad \|x'_i\|_2 \leq \|x\|_2 \quad \Rightarrow \quad \forall j \notin I \quad |x_j| = \sum_i |g_j^{C_i}| \leq K \sum_{i \in I} |x_i| = K \|z\|_1$$



$$x = L_I^W(z) = g^{C_1} + g^{C_2} + g^{C_3}$$

The condition numbers κ_A and $\bar{\chi}_A$

THEOREM: For every matrix $A \in \mathbb{R}^{m \times n}$, $n \geq 2$

$$\sqrt{1 + \kappa_A^2} \leq \bar{\chi}_A \leq n\kappa_A$$

Approximability of κ_A and $\bar{\chi}_A$:

LEMMA (Tunçel 1999): It is NP-hard to approximate $\bar{\chi}_A$ by a factor better than $2^{\text{poly}(\text{rank}(A))}$

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Part 3

Solving LPs:

from approximate to exact



Fast approximate LP algorithms

$$\begin{aligned} & \min c^T x \\ & Ax = b \\ & x \geq 0 \end{aligned}$$

- ε -approximate solution:
 - Approximately feasible: $\|Ax - b\| \leq \varepsilon(\|A\|_F R + \|b\|)$
 - Approximately optimal: $c^T x \leq \text{OPT} + \varepsilon \|c\| R$
- Finding an approximate solution with $\log\left(\frac{1}{\varepsilon}\right)$ running time dependence implies a weakly polynomial exact algorithm.

Fast approximate LP algorithms

$$\min_x c^T x \quad Ax = b \quad x \geq 0$$

- n variables, m equality constraints, Randomized vs. Deterministic
- Significant recent progress:
 - R $O\left((\text{nnz}(A) + m^2)\sqrt{m} \log^{O(1)}(n) \log\left(\frac{n}{\varepsilon}\right)\right)$ Lee-Sidford '13-'19
 - R $O\left(n^\omega \log^{O(1)}(n) \log\left(\frac{n}{\varepsilon}\right)\right)$ Cohen, Lee, Song '19
 - D $O\left(n^\omega \log^2(n) \log\left(\frac{n}{\varepsilon}\right)\right)$ van den Brand '20
 - R $O\left((mn + m^3) \log^{O(1)}(n) \log\left(\frac{n}{\varepsilon}\right)\right)$ van den Brand, Lee, Sidford, Song '20
 - R $O\left((mn + m^{2.5}) \log^{O(1)}(n) \log\left(\frac{n}{\varepsilon}\right)\right)$ van den Brand, Lee, Liu, Saranurak, Sidford, Song, Wang '21

Some important techniques:

- weighted and stochastic central paths
- fast approximate linear algebra
- efficient data structures

Fast exact LP algorithms with κ_A dependence

$$\begin{aligned} & \min c^\top x \\ & Ax = b \\ & x \geq 0 \end{aligned}$$

- n variables, m equality constraints

THEOREM (Dadush, Natura, V. '20) There exists a $\text{poly}(n, m, \log \kappa_A)$ algorithm for solving LP exactly.

- Feasibility: m calls to an approximate solver
- Optimization: mn calls to an approximate solver

with $\varepsilon = 1/(\text{poly}(n, \kappa_A))$. Using van den Brand '20, this gives a deterministic exact $O(mn^{\omega+1} \log^2(n) \log(\kappa_A+n))$ time LP optimization algorithm

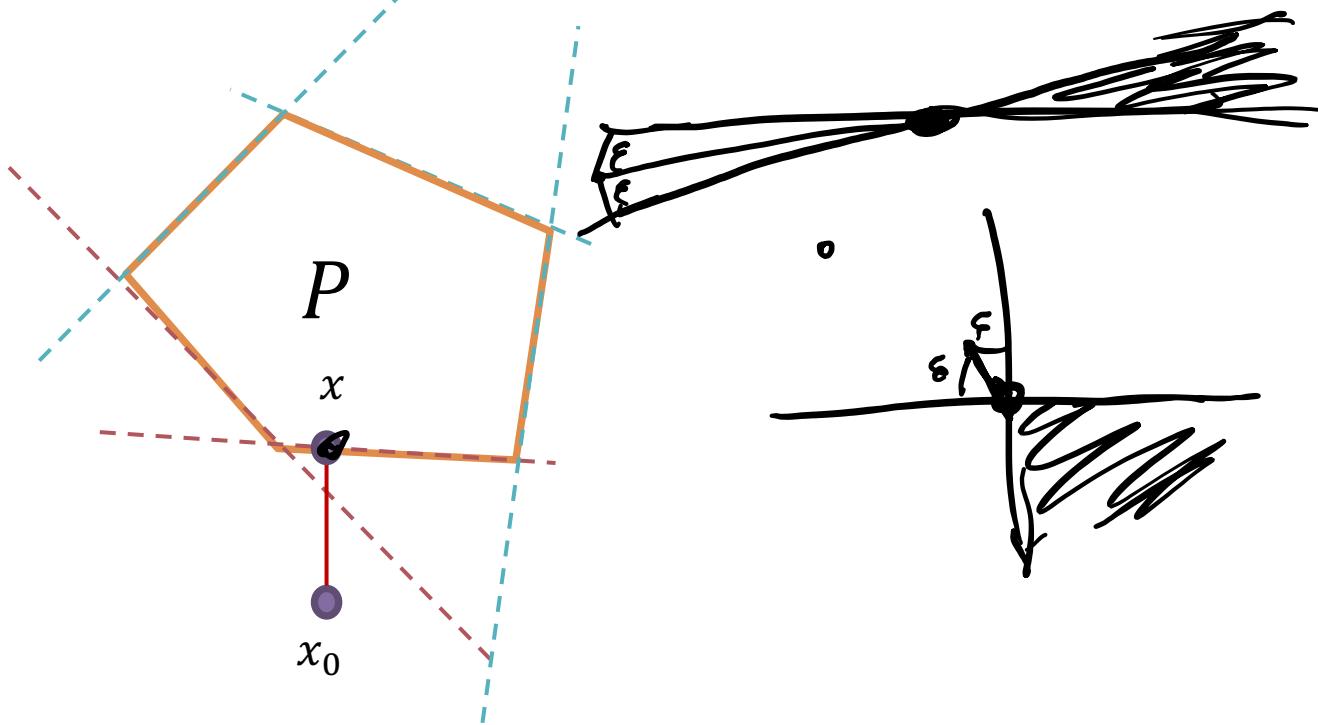
- Generalization of Tardos '86 for real constraint matrices and with directly working with approximate solvers.
- Main difference: arguments in Tardos '86 heavily rely on integrality assumptions

Hoffman's proximity theorem

Polyhedron $\underbrace{P = \{x \in \mathbb{R}^n : Ax \leq b\}}$, point $x_0 \notin P$, norms $\|\cdot\|_\alpha, \|\cdot\|_\beta$

THEOREM (Hoffman, 1952): There exists a constant $H_{\alpha,\beta}(A)$ such that

$$\exists x \in P: \|x - x_0\|_\alpha \leq H_{\alpha,\beta}(A) \underbrace{\|(Ax_0 - b)^+\|_\beta}_{\text{distance from } x_0 \text{ to the boundary}}$$



Alan J. Hoffman
1924-2021

LP in subspace form

- Matrix form: $A \in \mathbb{R}^{m \times n}, b \in \mathbb{R}^m, c \in \mathbb{R}^n$

$$\begin{array}{l} \min c^\top x \\ \underbrace{Ax = b}_{x \geq 0} \end{array}$$

$$\begin{array}{l} \max b^\top y \\ A^\top y + s = c \\ s \geq 0 \end{array}$$

- Subspace form: $W = \ker(A)$, $\underline{d} \in \mathbb{R}^n$ s.t. $\underline{Ad} = b$

$$\begin{array}{l} \min c^\top x \\ x \in W + \underline{d} \\ \tilde{x} \geq 0 \end{array}$$

$$\begin{array}{l} \max \underline{d}^\top (c - s) \\ s \in W^\perp + c \\ s \geq 0 \end{array}$$

$W = \ker(A)$

$W^\perp = \text{im}(A^\top)$

Proximity theorem with κ_A

THEOREM: For $A \in \mathbb{R}^{m \times n}$, $d \in \mathbb{R}^n$, consider the system

$$W = \text{ker}(A^\top)$$

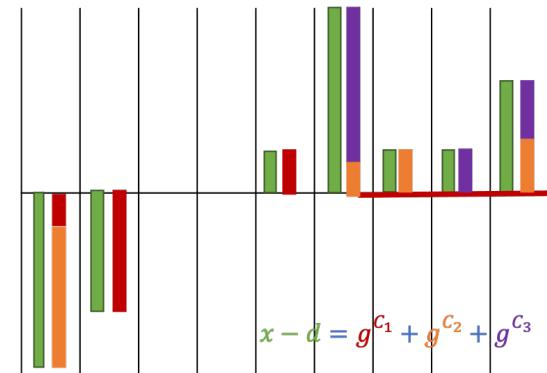
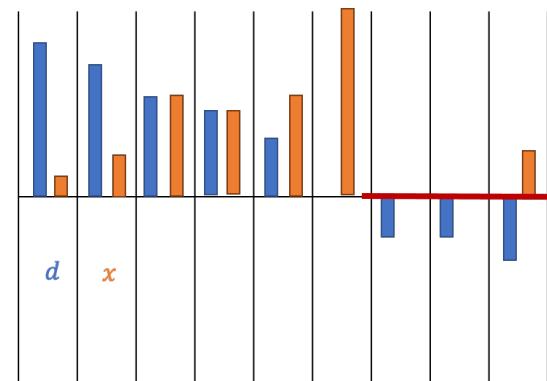
Assume it is feasible. $x \in W + d, x \geq 0$.

There exists a feasible solution x such that

$$\|x - d\|_\infty \leq \kappa_W \|d^\perp\|_1$$

PROOF:

$$\begin{aligned} Ad &= b \\ Ax &= Ad \\ x &\geq 0 \end{aligned}$$



Linear feasibility algorithm

Linear feasibility problem

$$x \in W + d, \quad x \geq 0.$$

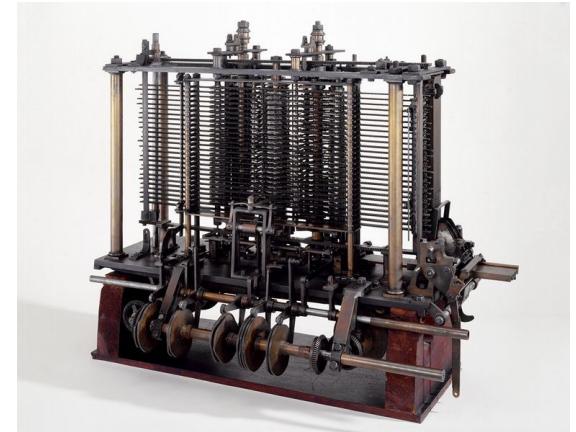
- Recursive algorithm using a **stronger** problem formulation:

$$x \in W + d, \quad x \geq 0.$$

$$\|x - d\|_\infty \leq C' \kappa_W^2 \|d^-\|_1$$

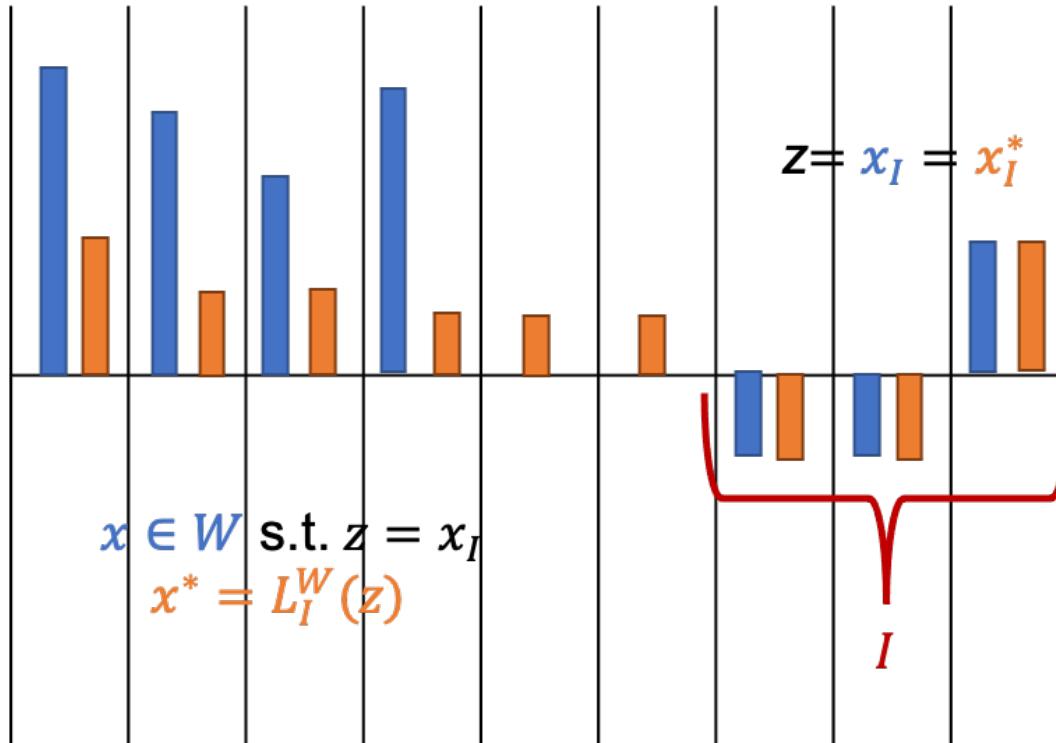
- Black box oracle for $\varepsilon = 1/(\text{poly}(n, \kappa_A))$

$x \in W + d$
proximity $\|x - d\|_\infty \leq C \kappa_W \|d^-\|_1$
error $\|x^-\|_\infty \leq \varepsilon \|d^-\|_1$



The lifting operator

$$L_I^W(z) = \arg \min\{\|x\|_2 : x \in W, x_I = z\}$$



The linear feasibility algorithm

1. Call the black box solver to find a solution z for $\varepsilon = 1/(\kappa_W n)^4$

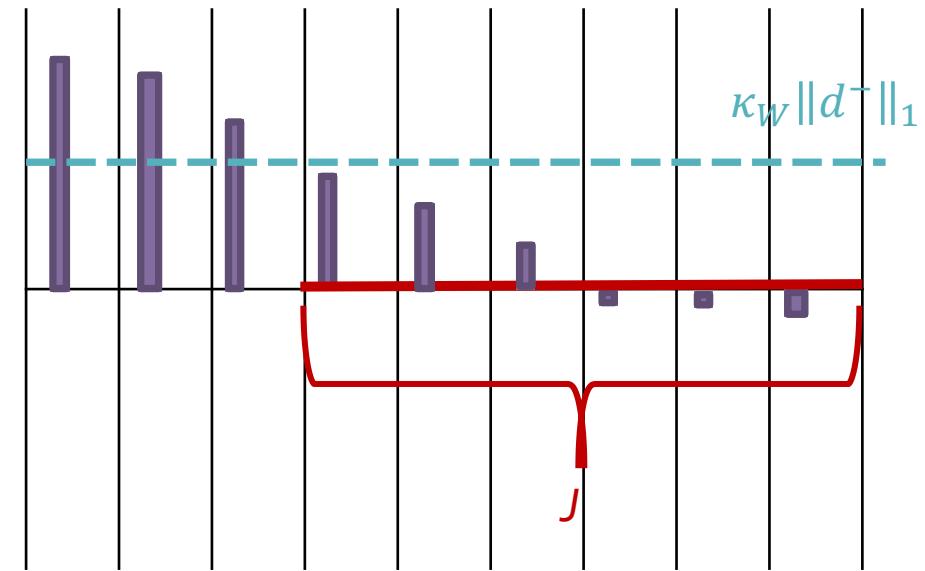


$$\begin{aligned} z &\in W + d \\ \|z - d\|_\infty &\leq C\kappa_W \|d^-\|_1 \\ \|z^-\|_\infty &\leq \varepsilon \|d^-\|_1 \end{aligned}$$

2. Set $J = \{i \in [n] : z_i < \kappa_W \|d^-\|_1\}$; assume $J \neq [n]$.
3. Recursively obtain $\tilde{x} \in \mathbb{R}_+^J$ from $\mathcal{F}(\pi_J(W), z_J)$
4. Return $x = z + L_J^W(\tilde{x} - z_J)$

Problem $\mathcal{F}(W, d)$

$$\begin{aligned} x &\in W + d \\ \|x - d\|_\infty &\leq C'\kappa_W^2 \|d^-\|_1 \\ x &\geq 0 \end{aligned}$$



1. Call the black box solver to find a solution z for $\varepsilon = 1/(\kappa_W n)^4$

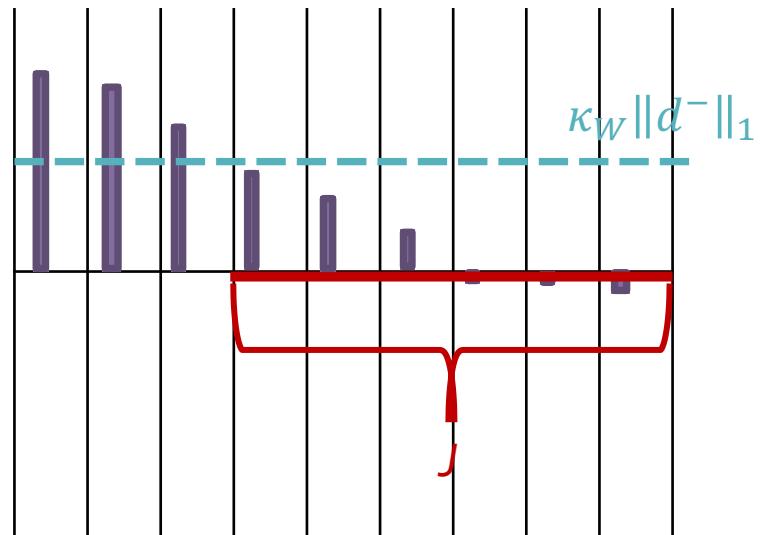


$$\begin{aligned} z &\in W + d \\ \|z - d\|_\infty &\leq C\kappa_W \|d^-\|_1 \\ \|z^-\|_\infty &\leq \varepsilon \|d^-\|_1 \end{aligned}$$

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Problem $\mathcal{F}(W, d)$

$$\begin{aligned} x &\in W + d \\ \|x - d\|_\infty &\leq C'\kappa_W^2 \|d^-\|_1 \\ x &\geq 0 \end{aligned}$$



The linear feasibility algorithm

$$J = \{i \in [n] : z_i < \kappa_W \|d^-\|_1\};$$

- If $J = [n]$, then we replace d by its projection to W^\perp
- Bound n on the number of recursive calls; can be decreased to m
- $O(mn^{\omega+o(1)} \log(\kappa_W + n))$ feasibility algorithm using van den Brand '20.

Certification

- In case of infeasibility we return an exact Farkas certificate
- κ_W is hard to approximate within $2^{O(n)}$ Tunçel 1999
- We use an estimate M in the algorithm
- The algorithm may fail if $\|L_J^W(\tilde{x} - z_J)\|_\infty > M\|\tilde{x} - z_J\|_1$
- In this case, we restart with

$$\max \left\{ M^2, \frac{\|L_J^W(\tilde{x} - z_J)\|_\infty}{\|\tilde{x} - z_J\|_1} \right\}$$

- Our estimate never overshoots κ_W by much, but can be significantly better.

Proximity for optimization

$$\begin{array}{ll} \min c^\top x & \max d^\top (c - s) \\ x \in W + d & s \in W^\perp + c \\ x \geq 0 & s \geq 0 \end{array}$$

THEOREM: Let $s \in W^\perp + c, s \geq 0$ be a feasible dual solution, and assume the primal is also feasible. Then there exists a primal optimal $x^* \in W + d, x^* \geq 0$ such that

$$\|x^* - d\|_\infty \leq \kappa_W \left(\|d^- \|_1 + \|d_{\text{supp}(s)}\|_1 \right).$$

Optimization algorithm

$$\begin{array}{ll}\min c^T x \\ x \in W + d \\ x \geq 0\end{array}$$

$$\begin{array}{ll}\max d^T(c - s) \\ s \in W^\perp + c \\ s \geq 0\end{array}$$

- $\textcolor{brown}{nm}$ calls to the black box solver
- $\leq n$ Outer Loops, each comprising $\leq m$ Inner Loops
- Each Outer Loop finds \tilde{d} with $\|d - \tilde{d}\|$ "small", and (x, s) primal and dual optimal solutions to
$$\min c^T x \text{ } s.t. x \in W + \tilde{d}, d \geq 0$$
- Using proximity, we can use this to conclude $x_I > 0$ for a certain variable set $I \subseteq n$ and recurse.

Outline of the lectures

1. Tardos's algorithm for min-cost flows
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Part 4

Optimizing circuit imbalances



Diagonal rescaling of LP

$$\begin{array}{ll} \min c^\top x & \max b^\top y \\ Ax = b & A^\top y + s = c \\ x \geq 0 & s \geq 0 \end{array}$$

Positive diagonal matrix $D \in \mathbb{R}^{n \times n}$

$$\begin{array}{ll} \min (\textcolor{brown}{D}c)^\top x' & \max b^\top y' \\ A\textcolor{brown}{D}x' = b & (AD)^\top y' + s' = \textcolor{brown}{D}c \\ x' \geq 0 & s' \geq 0 \end{array}$$

Mapping between solutions:

$$x' = \textcolor{brown}{D}^{-1}x, \quad y' = y, \quad s' = \textcolor{brown}{D}s$$

Diagonal rescaling of LP

Positive diagonal matrix $D \in \mathbb{R}^{n \times n}$

$$\begin{array}{ll}\min & (\mathbf{D}c)^\top x' \\ A\mathbf{D}x' = b & \\ x' \geq 0 & \end{array} \quad \begin{array}{ll}\max & b^\top y' \\ (AD)^\top y' + s' = \mathbf{D}c \\ s' \geq 0 & \end{array}$$

Mapping between solutions:

$$x' = \mathbf{D}^{-1}x, \quad y' = y, \quad s' = \mathbf{D}s$$

- Natural symmetry of LPs and many LP algorithms.
- The **Central Path** is invariant under diagonal scaling.
- Most “standard” interior point methods are invariant.

Dependence on the constraint matrix only

$$\min c^\top x, Ax = b \quad x \geq 0$$

- Algorithms with running time dependent only on A , but not on b and c .

- Combinatorial LP's: integer matrix $A \in \mathbb{Z}^{m \times n}$.

$$\Delta_A = \max\{|\det(B)| : B \text{ submatrix of } A\}$$

Tardos '86: $\text{poly}(n, m, \log \Delta_A)$ LP algorithm 

- Layered-least-squares (LLS) Interior Point Method
Vavasis-Ye '96: $\text{poly}(n, m, \log \bar{\chi}_A)$ LP algorithm in the real model of computation

$\bar{\chi}_A$: condition number 

- Dadush-Huiberts-Natura-V '20: $\text{poly}(n, m, \log \bar{\chi}_A^*)$
 $\bar{\chi}_A^*$: optimized version of $\bar{\chi}_A$ 

Optimizing κ_A and $\bar{\chi}_A$ by rescaling

\mathcal{D} = set of $n \times n$ positive diagonal matrices

$$\kappa_A^* = \inf\{\kappa_{AD} : D \in \mathcal{D}\}$$

$$\bar{\chi}_A^* = \inf\{\bar{\chi}_{AD} : D \in \mathcal{D}\}$$

- A scaling invariant algorithm with $\bar{\chi}_A$ dependence automatically yields $\bar{\chi}_A^*$ dependence.
- Recall $\sqrt{1 + \kappa_A^2} \leq \bar{\chi}_A \leq n\kappa_A$.

THEOREM (Dadush-Huiberts-Natura-V '20): Given $A \in \mathbb{R}^{m \times n}$, in $O(n^2m^2 + n^3)$ time, one can

- approximate the value κ_A within a factor $(\kappa_A^*)^2$, and
- compute a rescaling $D \in \mathcal{D}$ satisfying $\kappa_{AD} \leq (\kappa_A^*)^3$.

THEOREM (Tunçel 1999): It is NP-hard to approximate $\bar{\chi}_A$ (and thus κ_A) by a factor better than $2^{\text{poly}(\text{rank}(A))}$

Approximating κ_A^*

\mathcal{D} = set of $n \times n$ positive diagonal matrices

$$\kappa_A^* = \inf\{\kappa_{AD} : D \in \mathcal{D}\}$$

- **EXAMPLE:** Let $\ker(A) = \text{span}((0,1,1,M), (1,0,M,1))$

Pairwise circuit imbalances

- For a circuit C , there exists a vector $g^C \in \mathbb{R}^C$ unique up to a scalar multiplier such that

$$\sum_{i \in C} g_i^C a_i = 0$$

- \mathcal{C}_A : set of all circuits.

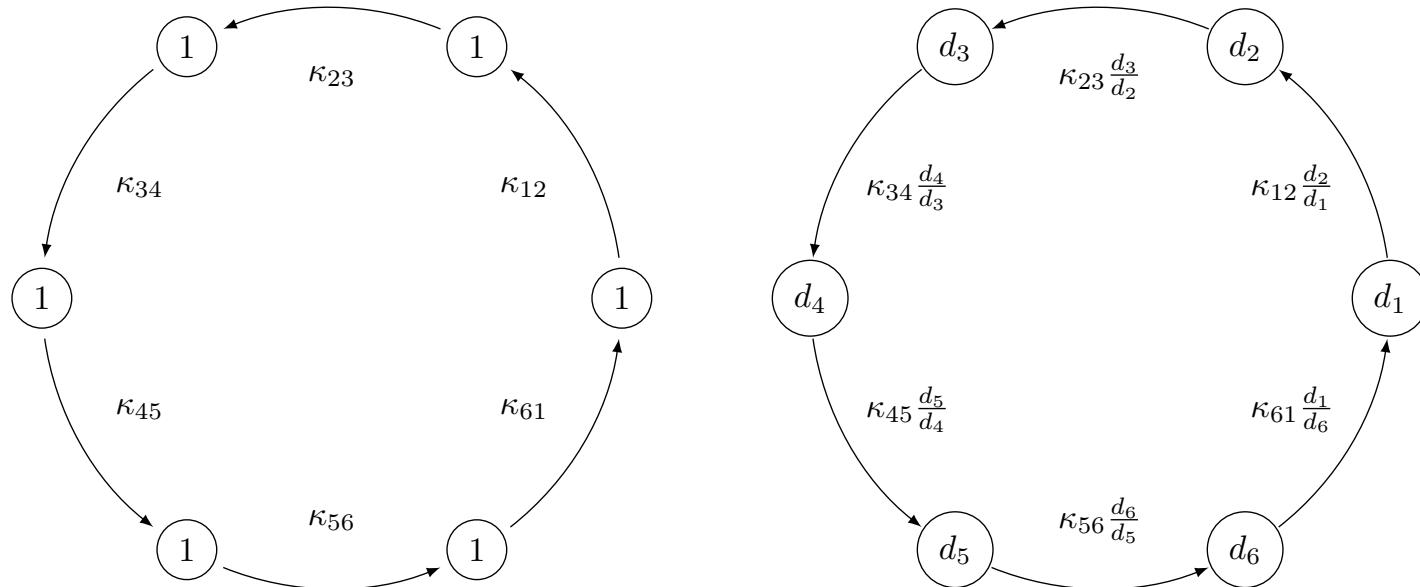
- For any $i, j \in [n]$,

$$\kappa_{ij} = \max \left\{ \frac{|g_j^C|}{|g_i^C|} : C \in \mathcal{C}_A, \text{s.t. } i, j \in C \right\}$$

- The circuit imbalance measure is

$$\kappa_A = \max_{i, j \in [n]} \kappa_{ij}$$

Cycles are invariant under scaling



LEMMA For any directed cycle H on $\{1, 2, \dots, n\}$

$$(\kappa_A^*)^{|H|} \geq \prod_{(i,j) \in H} \kappa_{ij}$$

Circuit imbalance min-max formula

THEOREM (Dadush-Huiberts-Natura-V '20):

$$\kappa_A^* = \max \left\{ \left(\prod_{(i,j) \in H} \kappa_{ij} \right)^{1/|H|} : H \text{ directed cycle on } \{1, 2, \dots, n\} \right\}$$

PROOF:

Circuit imbalance min-max formula

THEOREM (Dadush-Huiberts-Natura-V '20):

$$\kappa_A^* = \max \left\{ \left(\prod_{(i,j) \in H} \kappa_{ij} \right)^{1/|H|} : H \text{ directed cycle on } \{1, 2, \dots, n\} \right\}$$

- BUT: Computing the κ_{ij} values is NP-complete...
- **LEMMA:** For any circuit $C \in \mathcal{C}_A$ s.t. $i, j \in C$,

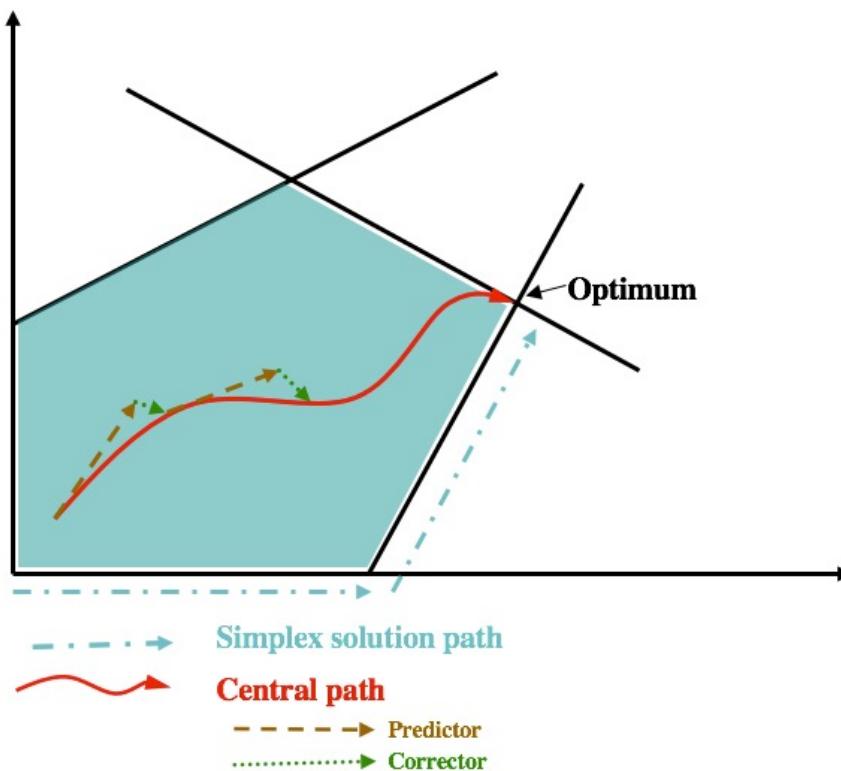
$$\frac{|g_j^C|}{|g_i^C|} \geq \frac{\kappa_{ij}}{(\kappa_W^*)^2}$$

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Part 5

Interior point methods: basic concepts



Primal and dual LP

- Matrix form: $A \in \mathbb{R}^{m \times n}, b \in \mathbb{R}^m, c \in \mathbb{R}^n$

$$\begin{array}{ll}\min c^\top x & \max b^\top y \\ Ax = b & A^\top y + s = c \\ x \geq 0 & s \geq 0\end{array}$$

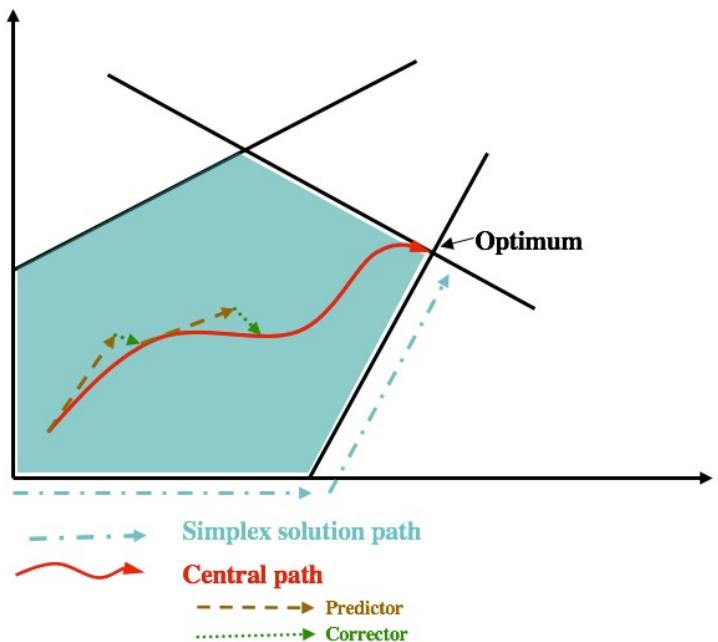
- Subspace form: $W = \ker(A)$, $d \in \mathbb{R}^n$ s.t. $Ad = b$

$$\begin{array}{ll}\min c^\top x & \max d^\top(c - s) \\ x \in W + d & s \in W^\top + c \\ x \geq 0 & s \geq 0\end{array}$$

- Complementary slackness: Primal and dual solutions (x, s) are optimal if $x^\top s = 0$: for each $i \in [n]$, either $x_i = 0$ or $s_i = 0$.
- Optimality gap:

$$c^\top x - d^\top(c - s) = x^\top s.$$

The central path

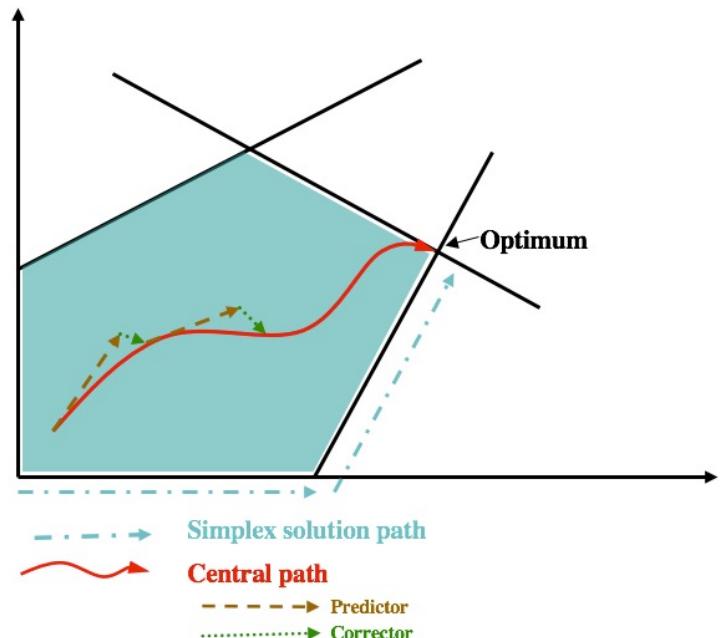


- For each $\mu > 0$, there exists a unique solution $w(\mu) = (x(\mu), y(\mu), s(\mu))$ such that
$$x(\mu)_i s(\mu)_i = \mu \quad \forall i \in [n]$$
the **central path element** for μ .
- The **central path** is the algebraic curve formed by $\{w(\mu) : \mu > 0\}$
- For $\mu \rightarrow 0$, the central path converges to an optimal solution $w^* = (x^*, y^*, s^*)$.
- The optimality gap is $s(\mu)^T x(\mu) = n\mu$.
- **Interior point algorithms:** walk down along the central path with μ decreasing geometrically.

The Mizuno-Todd-Ye Predictor-Corrector Algorithm

- Start from point $w_0 = (x_0, y_0, s_0)$ 'near' the central path at some $\mu_0 > 0$.
- Alternate between
 - **Predictor steps:** 'shoot down' the central path, decreasing μ by a factor at least $1 - \beta/n$.
May move slightly 'farther' from the central path.
 - **Corrector steps:** do not change parameter μ , but move back 'closer' to the central path.

Within $O(n)$ iterations, μ decreases by a factor 2.



The predictor step

- Step direction $\Delta w = (\Delta x, \Delta y, \Delta s)$

$$A\Delta x = 0$$

$$A^\top \Delta y + \Delta s = 0$$

$$s_i \Delta x_i + x_i \Delta s_i = -x_i s_i \quad \forall i \in [n]$$

- Pick the largest $\alpha \in [0,1]$ such that w' is still “close enough” to the central path
 $w' = w + \alpha \Delta w = (x + \alpha \Delta x, y + \alpha \Delta y, s + \alpha \Delta s)$

- Long step: $|\Delta x_i \Delta s_i|$ small for every $i \in [n]$
- New optimality gap is $(1 - \alpha)\mu$.

The predictor step – subspace view

$$\begin{aligned} A\Delta x &= 0 \\ A^T \Delta y + \Delta s &= 0 \\ s_i \Delta x_i + x_i \Delta s_i &= -x_i s_i \quad \forall i \in [n] \end{aligned}$$

- Assume the current point $w = (x, y, s)$ is on the central path. The steps can be found as minimum norm projections in the $(1/x)$ and $(1/s)$ rescaled norms

$$\Delta x = \arg \min \sum_{i=1}^n \left(\frac{x_i + \Delta x_i}{x_i} \right)^2 \text{ s.t. } x \in W = \ker(A)$$

$$\Delta s = \arg \min \sum_{i=1}^n \left(\frac{s_i + \Delta s_i}{s_i} \right)^2 \text{ s.t. } s \in W^\perp = \text{im}(A^T)$$

Some recent progress on interior point methods

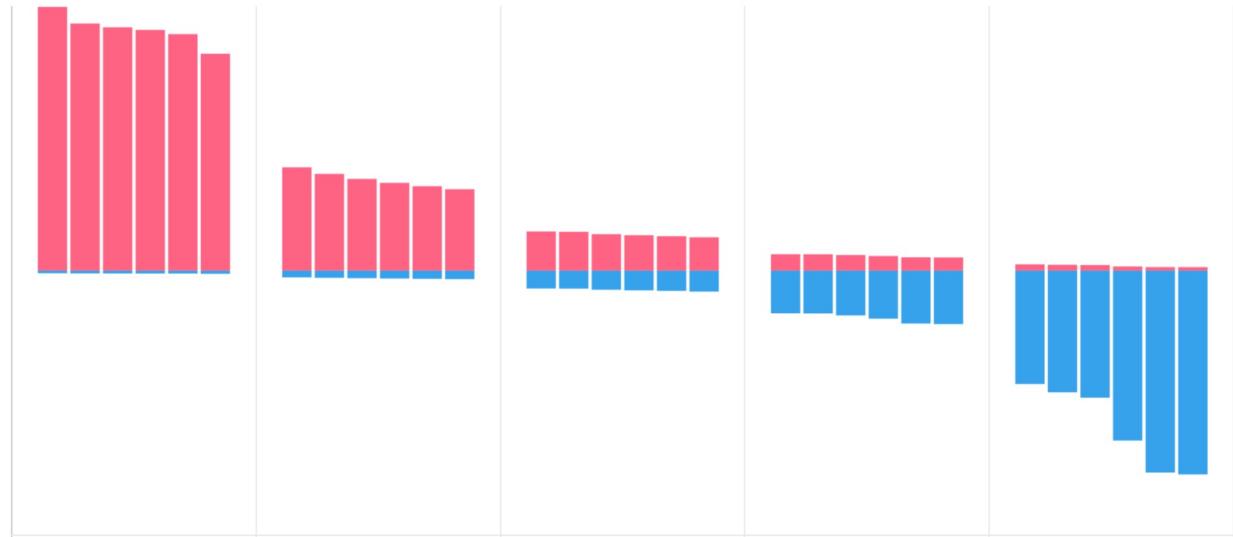
- Tremendous recent progress on fast approximate variants LS'14 – '19,
CLS'19, vdB'20, vdBLLSSW'20, vdBLLSSSW'21
- Fast approximate algorithms for combinatorial problems flows, matching and MDPs:
DS'08, M'13, M'16, CMSV'17, AMV'20,
vdBLNPTSSW'20, vdBLLSSSW'21

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Part 6

Layered-least-squares interior point methods



Layered-least-squares (LLS) Interior Point Methods:

Dependence on the constraint matrix only

$$\bar{\chi}_A^* = \inf\{\bar{\chi}_{AD} : D \in \mathcal{D}\}$$

- Vavasis-Ye '96: $O(n^{3.5} \log(\bar{\chi}_A + n))$ iterations
- Monteiro-Tsuchiya '03 $O(n^{3.5} \log(\bar{\chi}_A^* + n) + n^2 \log \log 1/\varepsilon)$ iterations
- Lan-Monteiro-Tsuchiya '09 $O(n^{3.5} \log(\bar{\chi}_A^* + n))$ iterations, but the running time of the iterations depends on b and c
- Dadush-Huiberts-Natura-V '20: scaling invariant LLS method with $O(n^{2.5} \log(n) \log(\bar{\chi}_A^* + n))$ iterations

Near monotonicity of the central path

LEMMA For $w = (x, y, s)$ on the central path, and for any solution $w' = (x', y', s')$ s.t. $(x')^\top s' \leq x^\top s$, we have

$$\sum_{i=1}^n \frac{x'_i}{x_i} + \frac{s'_i}{s_i} \leq 2n$$

PROOF:

IPM learns gradually improved upper bounds on the optimal solution.

Variable fixing...—or not?

LEMMA After every iteration, there exists variables x_i and s_j such that

$$\frac{1}{O(n)} \leq \frac{x_i}{x_i^*}, \frac{s_j}{s_j^*} \leq O(n)$$

For the optimal (x^*, y^*, s^*) . Thus, x_i and s_j have “*converged*” to their final values.

- **PROOF:** Can be shown using the form of the predictor step:

$$\Delta x = \arg \min \sum_{i=1}^n \left(\frac{x_i + \Delta x_i}{x_i} \right)^2 \text{ s. t. } x \in W$$

$$\Delta s = \arg \min \sum_{i=1}^n \left(\frac{s_i + \Delta s_i}{s_i} \right)^2 \text{ s. t. } s \in W^\perp$$

and bounds on the stepsize.

Variable fixing...—or not?

LEMMA After every iteration, there exists variables x_i and s_j such that

$$\frac{1}{O(n)} \leq \frac{x_i}{x_i^*}, \frac{s_j}{s_j^*} \leq O(n)$$

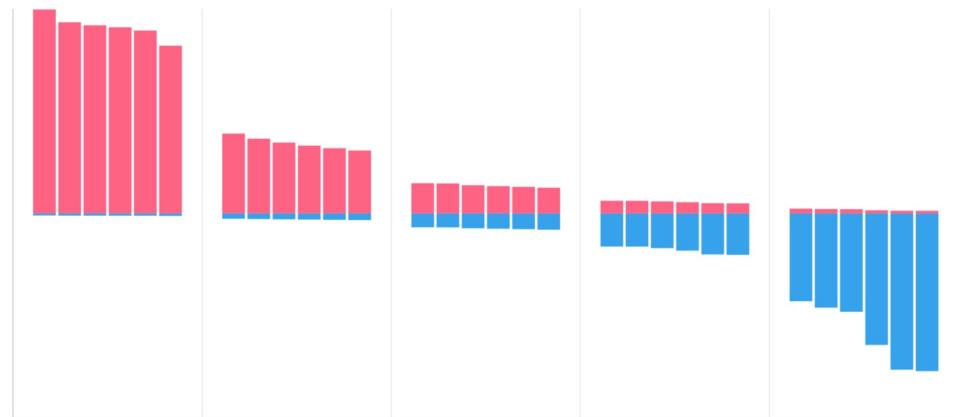
Thus, x_i and s_j have “converged” to their final values.

We cannot identify these indices,
just show their existence



Layered least squares methods

- Instead of the standard predictor step, split the variables into layers.
- Variables on different layers “behave almost like separate LPs”
- Force new primal and dual variables that must have converged.



Recap: the lifting operator and κ_A

- For a linear subspace $W \subset \mathbb{R}^n$ and index set $I \subseteq [n]$, we let

$$\pi_I: \mathbb{R}^n \rightarrow \mathbb{R}^I$$

denote the coordinate projection, and

$$\pi_I(W) = \{x_I : x \in W\}$$

- The lifting operator $L_I^W: \mathbb{R}^I \rightarrow \mathbb{R}^n$ is defined as

$$L_I^W(z) = \arg \min \{\|x\|_2 : x \in W, x_I = z\}$$

- LEMMA:** $\kappa_A = \max \left\{ \frac{\|L_I^W(z)\|_\infty}{\|z\|_1} : z \in \pi_I(W) \right\}$

- For every $z \in \pi_I(W)$, $x = L_I^W(z) \in W$ s.t.

$$x_I = z, \text{ and } \|x\|_\infty \leq \kappa_A \|z\|_1$$

Motivating the layering idea: final rounding step in standard IPM

$$\begin{array}{ll} \min c^T x & \max b^T y \\ Ax = b & A^T y + s = c \\ x \geq 0 & s \geq 0 \end{array}$$

- Limit optimal solution (x^*, y^*, s^*) , and optimal partition $[n] = B \cup N$ s.t. $B = \text{supp}(x^*)$, $N = \text{supp}(s^*)$.
- Given (x, y, s) near central path with ‘small enough’ $\mu = s^T x / n$ such that for every $i \in [n]$, either x_i or s_i very small.
- Assume that we can correctly guess
 $B = \{i: x_i > M\sqrt{\mu}\}, \quad N = \{i: s_i > M\sqrt{\mu}\}$

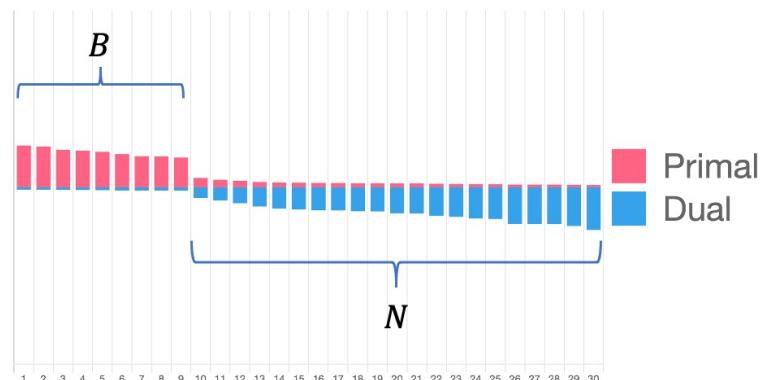
- Assume we have a partition B, N , we have

$$i \in B: x_i > M\sqrt{\mu}, \quad s_i < \sqrt{\mu}/M$$

$$i \in N: x_i < \sqrt{\mu}/M, \quad s_i > M\sqrt{\mu}$$

- **Goal:** move to $\bar{x} = x + \Delta x$, $\bar{y} = y + \Delta y$, $\bar{s} = s + \Delta s$
s.t. $\text{supp}(\bar{x}) \subseteq B$, $\text{supp}(\bar{s}) \subseteq N$. Then, $\bar{x}^\top \bar{s} = 0$: optimal solution.
- Choice:

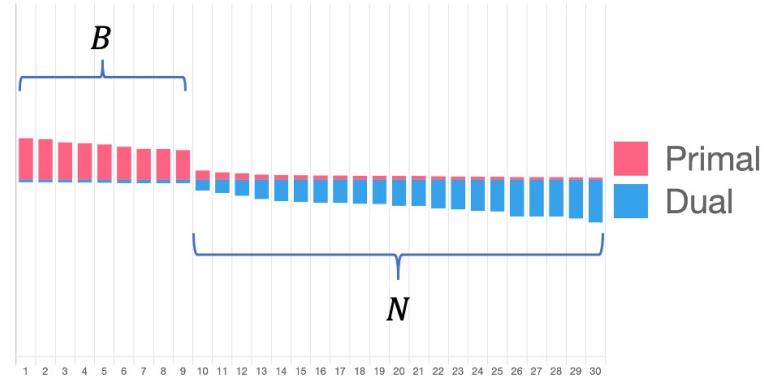
$$\Delta x = -L_N^W(x_N), \quad \Delta s = -L_B^W(s_B)$$



Layered-least-squares step

Assume we have a partition B, N , with

$$\begin{aligned} i \in B: x_i &> M\sqrt{\mu}, & s_i &< \sqrt{\mu}/M \\ i \in N: x_i &< \sqrt{\mu}/M, & s_i &> M\sqrt{\mu} \end{aligned}$$



Standard primal predictor step:

$$\Delta x = \arg \min \sum_{i=1}^n \left(\frac{x_i + \Delta x_i}{x_i} \right)^2$$

s. t. $\Delta x \in W$

Vavasis-Ye LLS step with layers
 (B, N) :

$$\Delta x_N = \arg \min \sum_{i \in N} \left(\frac{x_i + \Delta x_i}{x_i} \right)^2$$

s. t. $\Delta x \in W$

$$\Delta x_B = \arg \min \sum_{i \in B} \left(\frac{x_i + \Delta x_i}{x_i} \right)^2$$

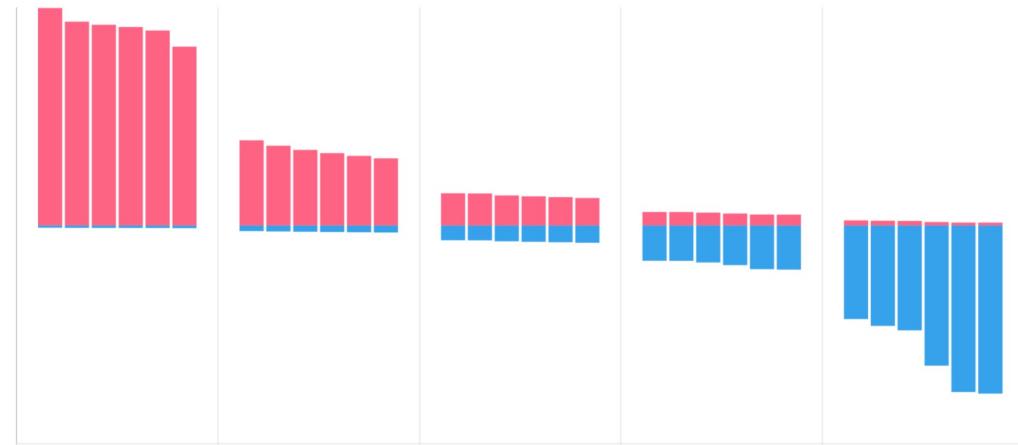
s. t. $(\Delta x_B, \Delta x_N) \in W$

Layered-least-squares step

Vavasis-Ye '96

- Order variables decreasingly as $x_1 \geq x_2 \geq \dots \geq x_n$
- Arrange variables into layers (J_1, J_2, \dots, J_t) ; start a new layer when
$$x_i > O(n^c) \bar{\chi}_A x_{i+1}$$
- Primal step direction by least squares problems from backwards, layer-by-layer
- Lifting costs from lower layers low
- Dual step in the opposite direction

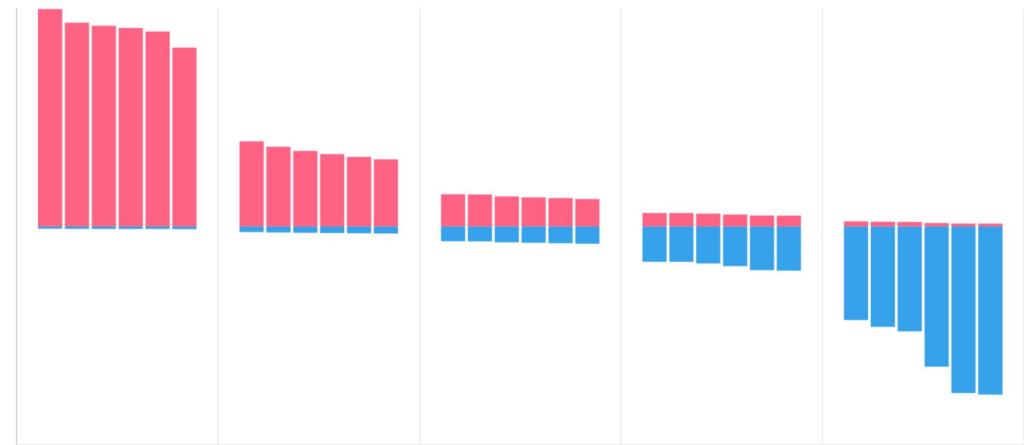
Not scaling invariant!



Progress measure: crossover events

Vavasis-Ye'96

- **DEFINITION:** The variables x_i and x_j cross over between μ and μ' , $\mu > \mu'$, if
 - $O(n^c)(\bar{\chi}_A)^n x_j(\mu) \geq x_i(\mu)$
 - $O(n^c)(\bar{\chi}_A)^n x_j(\mu'') < x_i(\mu'')$ for any $\mu'' \leq \mu'$
- **LEMMA:** In the Vavasis-Ye algorithm, a crossover event happens every $O(n^{1.5} \log(\bar{\chi}_A + n))$ iterations, totalling to $O(n^{3.5} \log(\bar{\chi}_A + n))$.



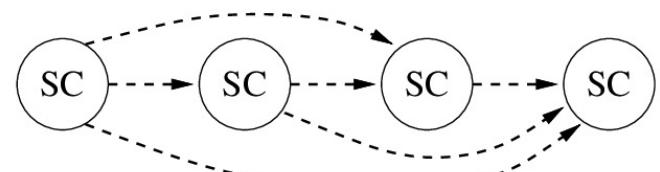
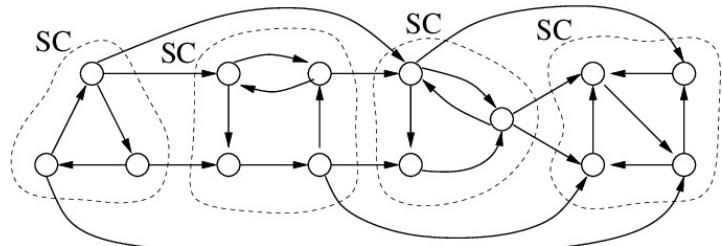
Scaling invariant layering

DNV'20

- Instead of the ratios x_i/x_j , we consider the rescaled circuit imbalance measures $\kappa_{ij}x_i/x_j$
- Layers: strongly connected components of the arcs

$$(i,j): \frac{\kappa_{ij}x_i}{x_j} > \frac{1}{\text{poly}(n)}$$

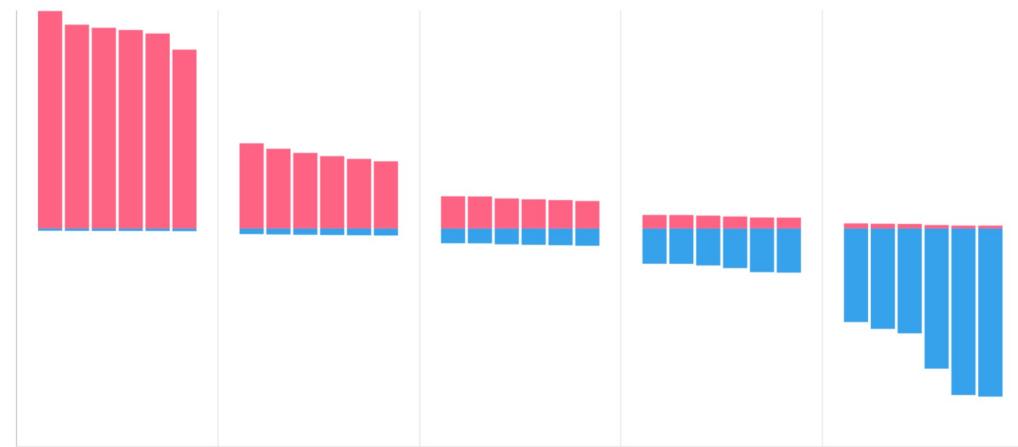
The κ_{ij} values are not known: increasingly improving estimates.



Scaling invariant crossover events

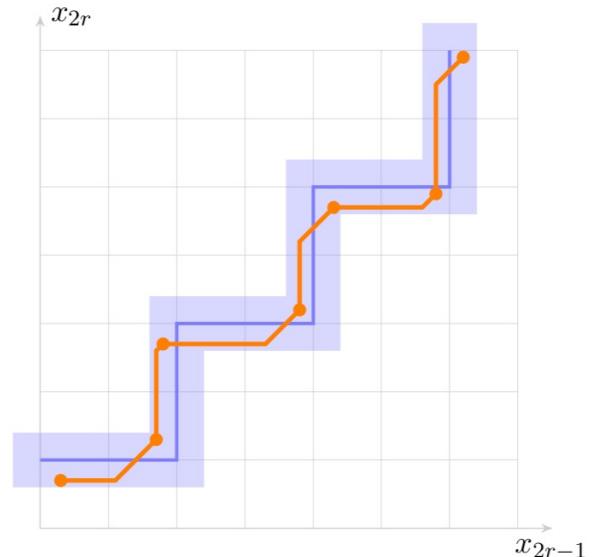
Vavasis-Ye'96

- **DEFINITION:** The variables x_i and x_j cross over between μ and μ' , $\mu > \mu'$, if
 - $O(n^c)(\bar{\chi}_A)^n x_j(\mu) \geq \kappa_{ij} x_i(\mu)$
 - $O(n^c)(\bar{\chi}_A)^n x_j(\mu'') < \kappa_{ij} x_i(\mu'')$ for any $\mu'' \leq \mu'$
- Amortized analysis, resulting in improved $O(n^{2.5} \log(n) \log(\bar{\chi}_A + n))$ iteration bound.



Limitation of IPMs

- **THEOREM (Allamigeon–Benchimol–Gaubert–Joswig ‘18):** No standard path following method can be strongly polynomial.
- Proof using **tropical geometry**: studies the tropical limit of a family of parametrized linear programs.



Future directions

- Circuit imbalance measure: key parameter for strongly polynomial solvability.
- LP classes with existence of strongly polynomial algorithms open:
 - LPs with 2 nonzeros per column in the constraint matrix, equivalently: min cost generalized flows
 - Undiscounted Markov Decision Processes
- Extend the theory of circuit imbalances more generally, to convex programming and integer programming.

Thank you!

Postdoc position open



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