

10 Karmarkars Algorithm

- ▶ inequalities $Ax \leq b$; $m \times n$ matrix A with rows a_i^T
- ▶ $P = \{x \mid Ax \leq b\}$; $P^\circ := \{x \mid Ax < b\}$
- ▶ interior point algorithm: $x \in P^\circ$ throughout the algorithm
- ▶ for $x \in P^\circ$ define

$$s_i(x) := b_i - a_i^T x$$

as the **slack** of the i -th constraint

logarithmic barrier function:

$$\phi(x) = - \sum_{i=1}^m \log(s_i(x))$$

Penalty for point x ; points close to the boundary have a very large penalty.

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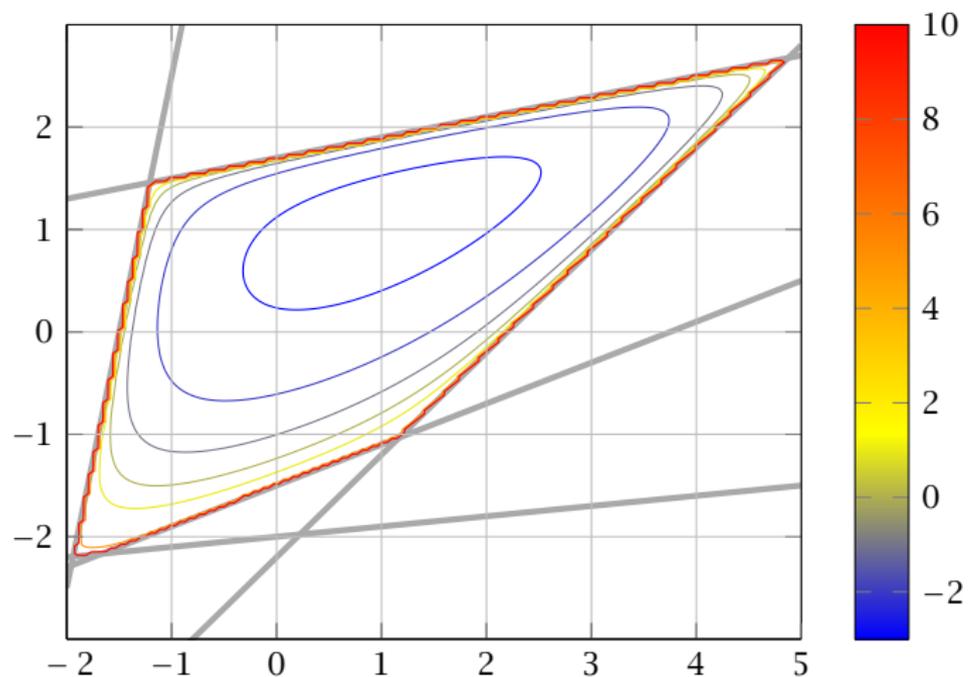
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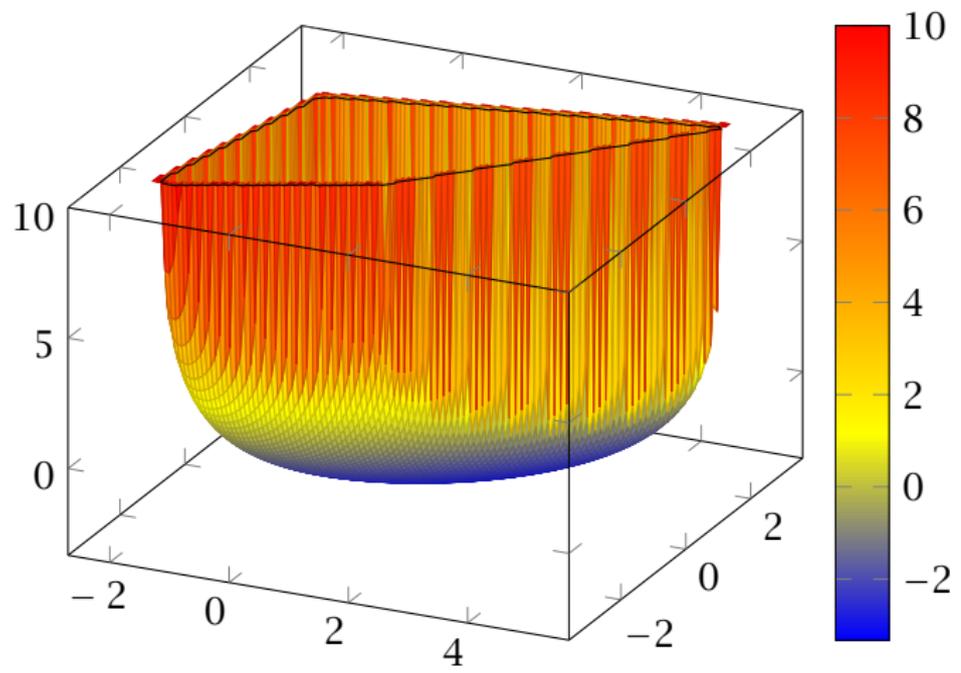
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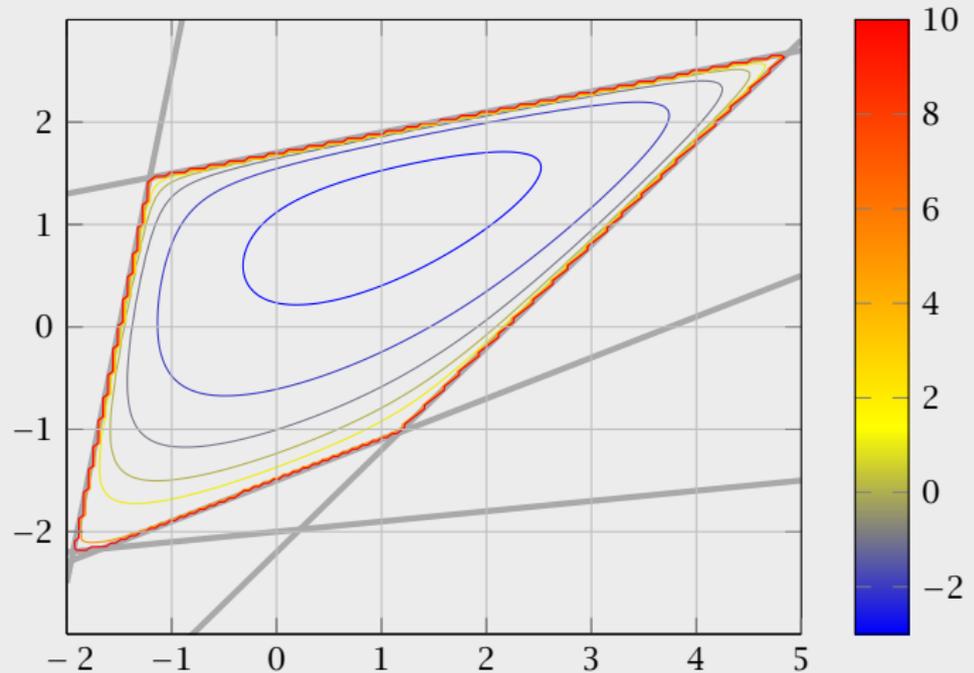
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Gradient and Hessian

Taylor approximation:

$$\phi(x + \epsilon) \approx \phi(x) + \nabla\phi(x)^T \epsilon + \frac{1}{2} \epsilon^T \nabla^2 \phi(x) \epsilon$$

Gradient:

$$\nabla\phi(x) = \sum_{i=1}^m \frac{1}{s_i(x)} \cdot a_i = A^T d_x$$

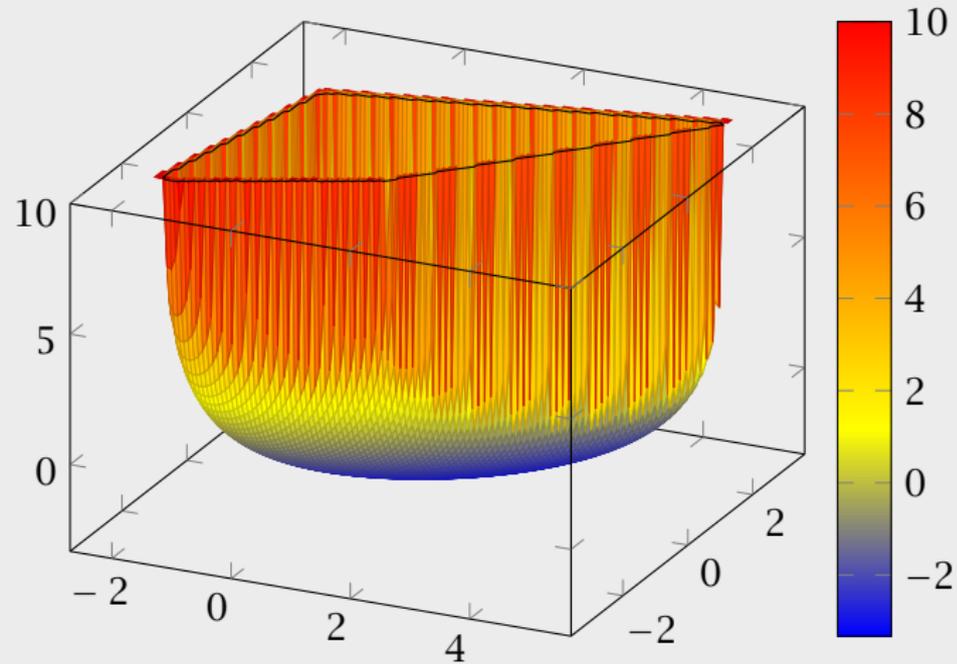
where $d_x^T = (1/s_1(x), \dots, 1/s_m(x))$. (d_x vector of inverse slacks)

Hessian:

$$H_x := \nabla^2 \phi(x) = \sum_{i=1}^m \frac{1}{s_i(x)^2} a_i a_i^T = A^T D_x^2 A$$

with $D_x = \text{diag}(d_x)$.

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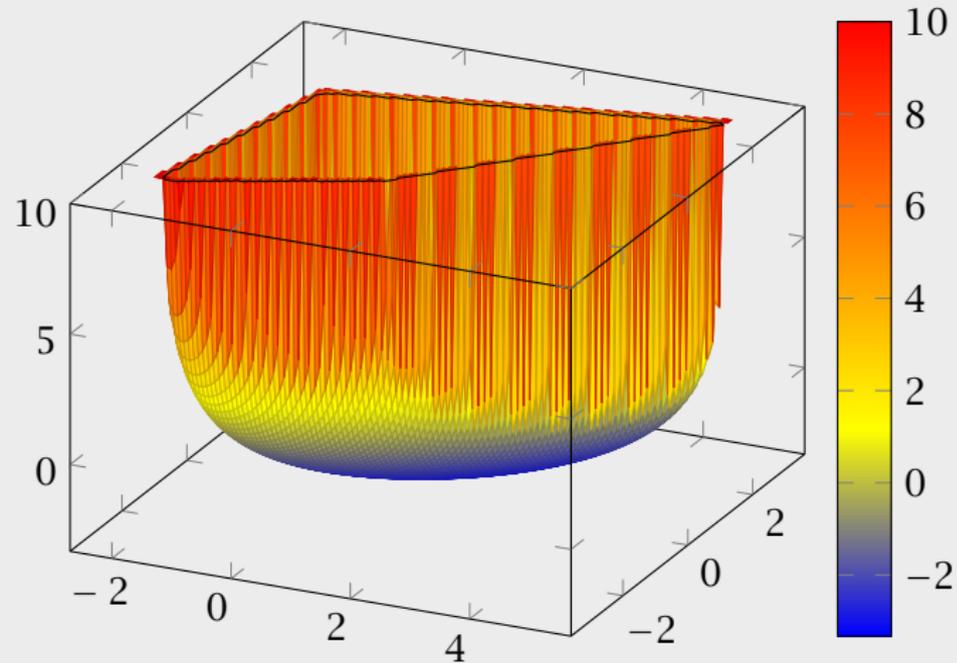
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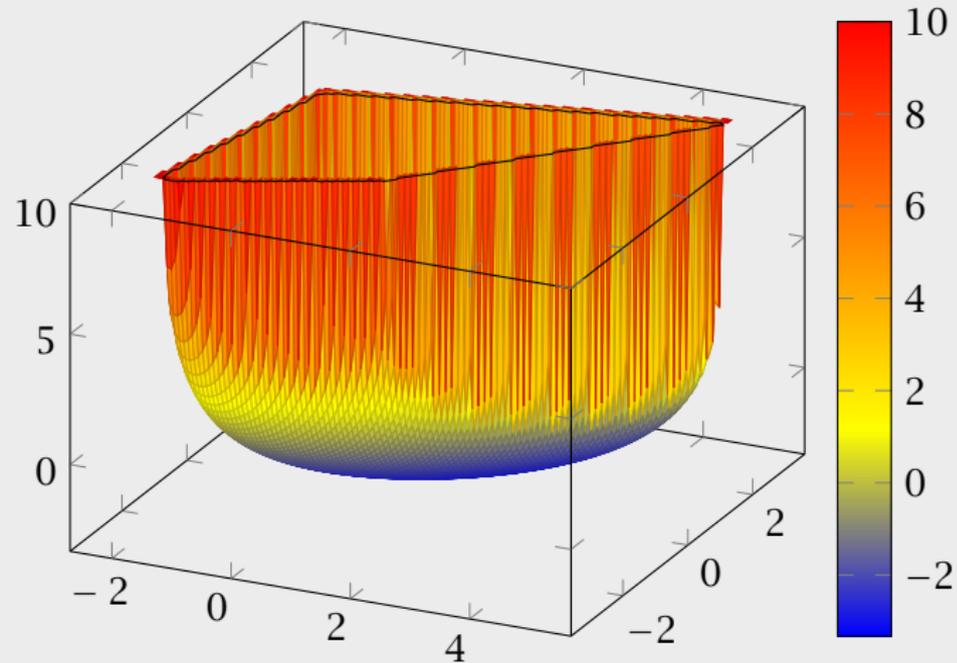
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Proof for Gradient

$$\begin{aligned}\frac{\partial \phi(x)}{\partial x_i} &= \frac{\partial}{\partial x_i} \left(- \sum_r \ln(s_r(x)) \right) \\ &= - \sum_r \frac{\partial}{\partial x_i} \left(\ln(s_r(x)) \right) = - \sum_r \frac{1}{s_r(x)} \frac{\partial}{\partial x_i} \left(s_r(x) \right) \\ &= - \sum_r \frac{1}{s_r(x)} \frac{\partial}{\partial x_i} \left(b_r - a_r^T x \right) = \sum_r \frac{1}{s_r(x)} \frac{\partial}{\partial x_i} \left(a_r^T x \right) \\ &= \sum_r \frac{1}{s_r(x)} A_{ri}\end{aligned}$$

The i -th entry of the gradient vector is $\sum_r 1/s_r(x) \cdot A_{ri}$. This gives that the gradient is

$$\nabla \phi(x) = \sum_r 1/s_r(x) a_r = A^T d_x$$

Gradient and Hessian

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Proof for Hessian

$$\begin{aligned}\frac{\partial}{\partial x_j} \left(\sum_r \frac{1}{s_r(x)} A_{ri} \right) &= \sum_r A_{ri} \left(-\frac{1}{s_r(x)^2} \right) \cdot \frac{\partial}{\partial x_j} (s_r(x)) \\ &= \sum_r A_{ri} \frac{1}{s_r(x)^2} A_{rj}\end{aligned}$$

Note that $\sum_r A_{ri} A_{rj} = (A^T A)_{ij}$. Adding the additional factors $1/s_r(x)^2$ can be done with a diagonal matrix.

Hence the Hessian is

$$H_x = A^T D^2 A$$

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Properties of the Hessian

H_x is positive semi-definite for $x \in P^\circ$

$$u^T H_x u = u^T A^T D_x^2 A u = \|D_x A u\|_2^2 \geq 0$$

This gives that $\phi(x)$ is convex.

If $\text{rank}(A) = n$, H_x is positive definite for $x \in P^\circ$

$$u^T H_x u = \|D_x A u\|_2^2 > 0 \text{ for } u \neq 0$$

This gives that $\phi(x)$ is **strictly** convex.

$\|u\|_{H_x} := \sqrt{u^T H_x u}$ is a (semi-)norm; the unit ball w.r.t. this norm is an ellipsoid.

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Dikin Ellipsoid

$$E_x = \{y \mid (y - x)^T H_x (y - x) \leq 1\} = \{y \mid \|y - x\|_{H_x} \leq 1\}$$

Points in E_x are feasible!!!

When x is a local minimum, the Hessian is positive semi-definite. If x is not a local minimum, the Hessian is not positive semi-definite and the Dikin Ellipsoid is not an ellipsoid.

In order to become infeasible when going from x to y one of the terms in the sum would need to be larger than 1.

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In order to become infeasible when going from x to y one of the terms in the sum would need to be larger than 1.

Properties of the Hessian

H_x is positive semi-definite for $x \in P^\circ$

$$u^T H_x u = u^T A^T D_x^2 A u = \|D_x A u\|_2^2 \geq 0$$

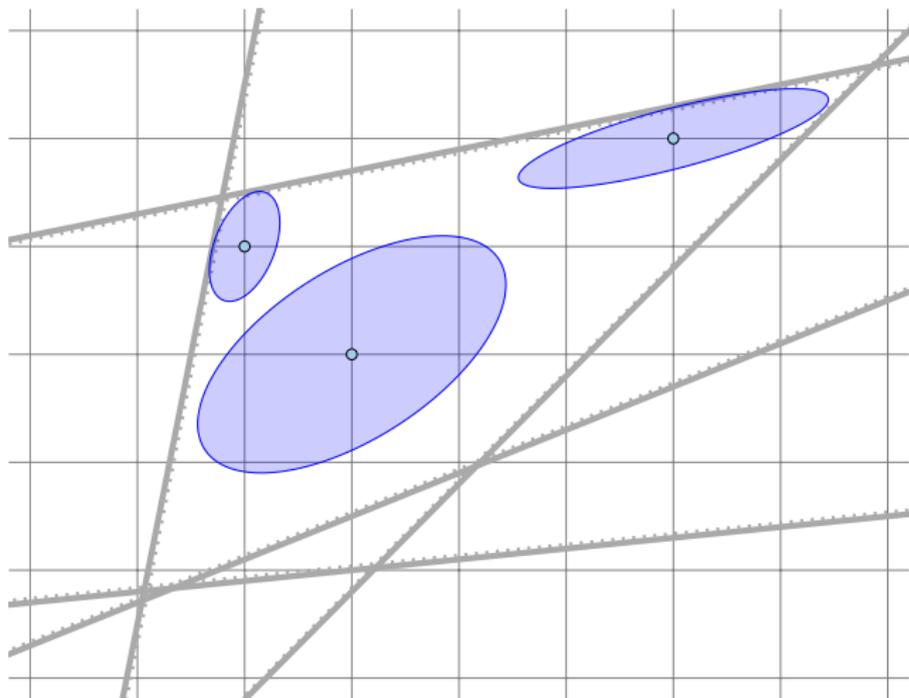
This gives that $\phi(x)$ is convex.

If $\text{rank}(A) = n$, H_x is positive definite for $x \in P^\circ$

$$u^T H_x u = \|D_x A u\|_2^2 > 0 \text{ for } u \neq 0$$

This gives that $\phi(x)$ is **strictly** convex.

$\|u\|_{H_x} := \sqrt{u^T H_x u}$ is a (semi-)norm; the unit ball w.r.t. this norm is an ellipsoid.



$$E_x = \{y \mid (y - x)^T H_x (y - x) \leq 1\} = \{y \mid \|y - x\|_{H_x} \leq 1\}$$

Points in E_x are feasible!!!

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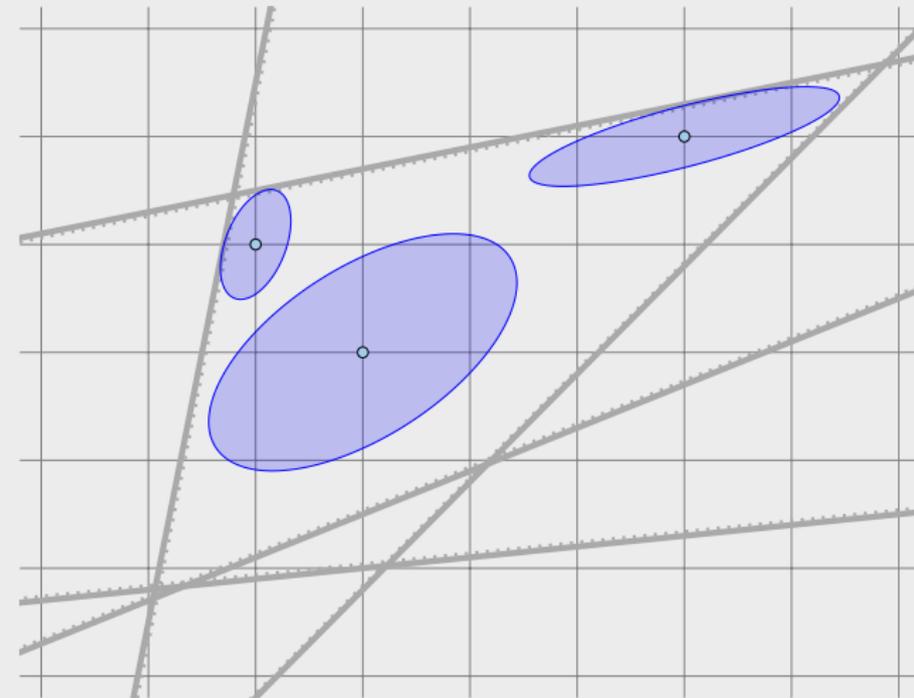
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$$x_{ac} := \arg \min_{x \in P^\circ} \phi(x)$$

- ▶ x_{ac} is solution to

$$\nabla \phi(x) = \sum_{i=1}^m \frac{1}{s_i(x)} a_i = 0$$

- ▶ depends on the **description** of the polytope
- ▶ x_{ac} exists and is unique iff P° is nonempty and bounded



Central Path

In the following we assume that the LP and its dual are **strictly feasible** and that $\text{rank}(A) = n$.

Central Path:

Set of points $\{x^*(t) \mid t > 0\}$ with

$$x^*(t) = \operatorname{argmin}_x \{tc^T x + \phi(x)\}$$

- ▶ $t = 0$: analytic center
- ▶ $t = \infty$: optimum solution

$x^*(t)$ exists and is unique for all $t \geq 0$.

Analytic Center

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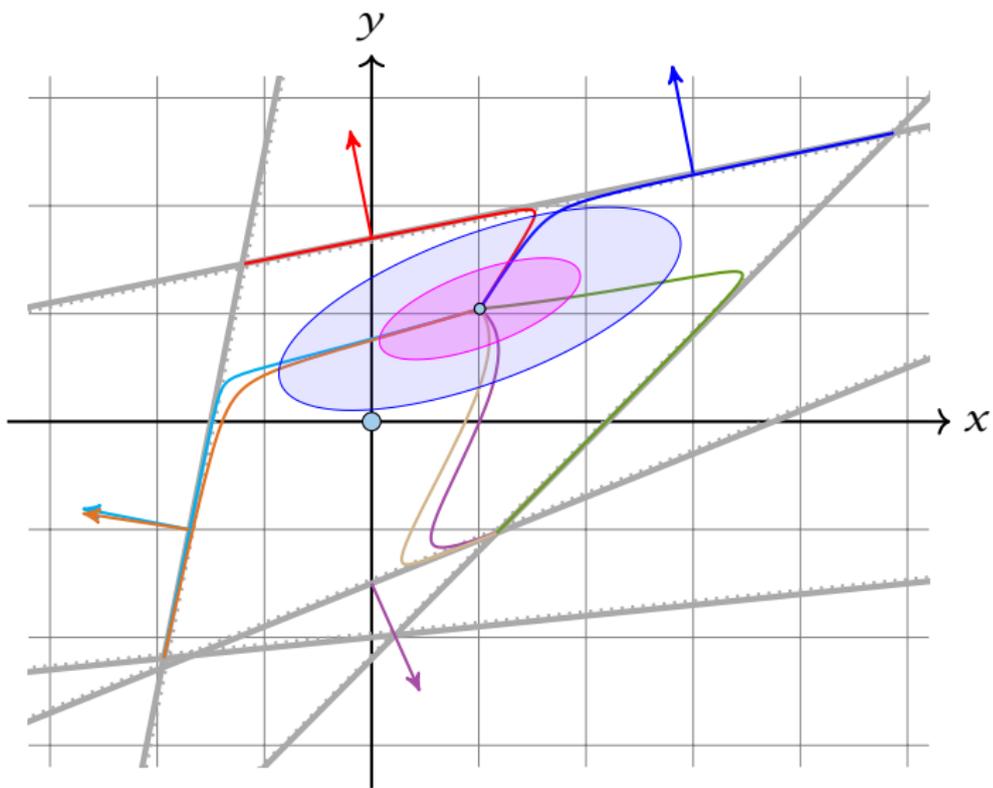
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Different Central Paths



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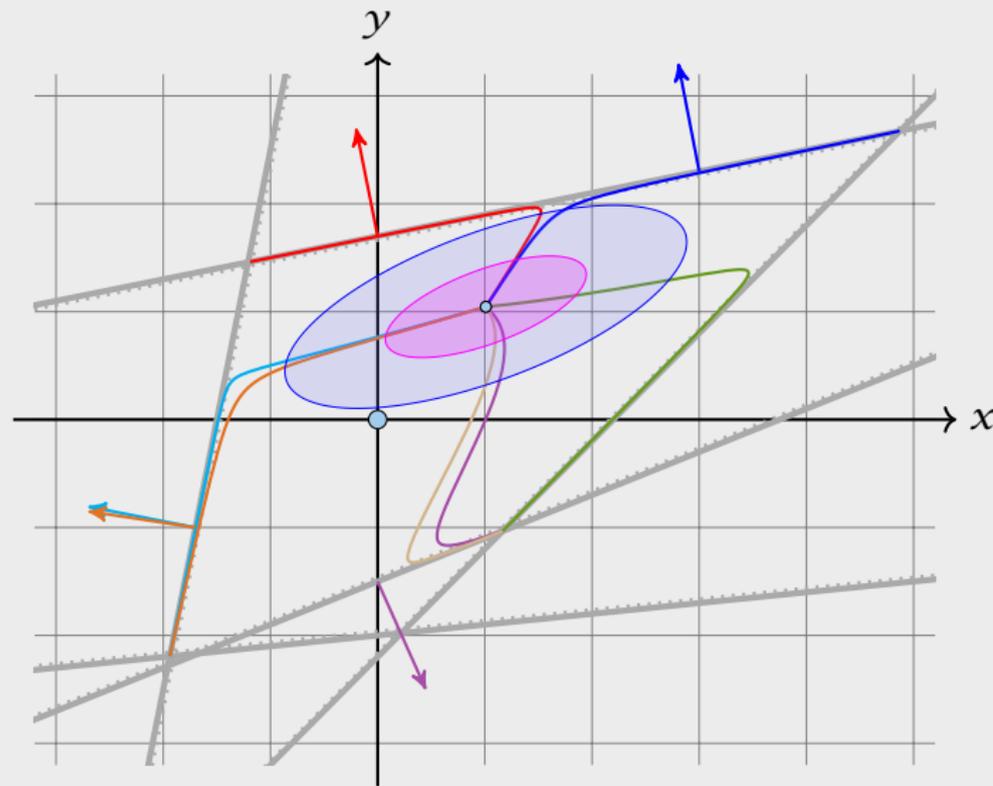
Intuitive Idea:

Find point on central path for large value of t . Should be close to optimum solution.

Questions:

- ▶ Is this really true? How large a t do we need?
- ▶ How do we find corresponding point $x^*(t)$ on central path?

Different Central Paths



The Dual

primal-dual pair:

$$\begin{array}{ll} \min & c^T x \\ \text{s.t.} & Ax \leq b \end{array}$$

$$\begin{array}{ll} \max & -b^T z \\ \text{s.t.} & A^T z + c = 0 \\ & z \geq 0 \end{array}$$

Assumptions

- ▶ primal and dual problems are strictly feasible;
- ▶ $\text{rank}(A) = n$.

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Force Field Interpretation

Point $x^*(t)$ on central path is solution to $tc + \nabla\phi(x) = 0$

- ▶ We can view each constraint as generating a repelling force. The combination of these forces is represented by $\nabla\phi(x)$.
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How large should t be?

Point $x^*(t)$ on central path is solution to $tc + \nabla\phi(x) = 0$.

This means

$$tc + \sum_{i=1}^m \frac{1}{s_i(x^*(t))} a_i = 0$$

or

$$c + \sum_{i=1}^m z_i^*(t) a_i = 0 \quad \text{with} \quad z_i^*(t) = \frac{1}{ts_i(x^*(t))}$$

It is not strictly true that feasibility is maintained for all t .
The quality gap between $x^*(t)$ and the optimum is

$\phi(x^*(t)) - \phi^* = \frac{1}{t} \sum_{i=1}^m \frac{1}{s_i(x^*(t))} \log s_i(x^*(t))$

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As t increases, the forces generated by the constraints are repelling and the force generated by the objective is pulling towards the optimum solution.

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▶ if gap is less than $1/2^{\Omega(L)}$ we can snap to optimum point

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First idea:

- ▶ start somewhere in the polytope
- ▶ use iterative method (Newtons method) to minimize $f_t(x) := tc^T x + \phi(x)$

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Newton Method

Quadratic approximation of f_t

$$f_t(x + \epsilon) \approx f_t(x) + \nabla f_t(x)^T \epsilon + \frac{1}{2} \epsilon^T H_{f_t}(x) \epsilon$$

Suppose this were exact:

$$f_t(x + \epsilon) = f_t(x) + \nabla f_t(x)^T \epsilon + \frac{1}{2} \epsilon^T H_{f_t}(x) \epsilon$$

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We want to move to a point where this gradient is 0:

Newton Step at $x \in P^\circ$

$$\begin{aligned}\Delta x_{\text{nt}} &= -H_{f_t}^{-1}(x) \nabla f_t(x) \\ &= -H_{f_t}^{-1}(x) (tc + \nabla \phi(x)) \\ &= -(A^T D_x^2 A)^{-1} (tc + A^T d_x)\end{aligned}$$

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$$x := x + \Delta x_{\text{nt}}$$

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Measuring Progress of Newton Step

Newton decrement:

$$\begin{aligned}\lambda_t(x) &= \|D_x A \Delta x_{nt}\| \\ &= \|\Delta x_{nt}\|_{H_x}\end{aligned}$$

Square of Newton decrement is linear estimate of reduction if we do a Newton step:

$$-\lambda_t(x)^2 = \nabla f_t(x)^T \Delta x_{nt}$$

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Convergence of Newtons Method

Theorem 2

If $\lambda_t(\mathbf{x}) < 1$ then

- ▶ $\mathbf{x}_+ := \mathbf{x} + \Delta\mathbf{x}_{nt} \in P^\circ$ (new point feasible)
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This means we have **quadratic convergence**. Very fast.

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feasibility:

- ▶ $\lambda_t(x) = \|\Delta x_{nt}\|_{H_x} < 1$; hence x_+ lies in the **Dikin ellipsoid** around x .

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bound on $\lambda_t(\mathbf{x}^+)$:

we use $D := D_x = \text{diag}(d_x)$ and $D_+ := D_{x^+} = \text{diag}(d_{x^+})$

To see the last equality we use Pythagoras

$$\|a\|^2 + \|a + b\|^2 = \|b\|^2$$

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$$\begin{aligned}a^T (a + b) &= \Delta \mathbf{x}_{\text{nt}}^{+T} A^T D_+ (D_+ A \Delta \mathbf{x}_{\text{nt}}^+ + (I - D_+^{-1} D) D A \Delta \mathbf{x}_{\text{nt}}) \\ &= \Delta \mathbf{x}_{\text{nt}}^{+T} (A^T D_+^2 A \Delta \mathbf{x}_{\text{nt}}^+ - A^T D^2 A \Delta \mathbf{x}_{\text{nt}} + A^T D_+ D A \Delta \mathbf{x}_{\text{nt}}) \\ &= \Delta \mathbf{x}_{\text{nt}}^{+T} (H_+ \Delta \mathbf{x}_{\text{nt}}^+ - H \Delta \mathbf{x}_{\text{nt}} + A^T D_+ \vec{\mathbf{1}} - A^T D \vec{\mathbf{1}}) \\ &= \Delta \mathbf{x}_{\text{nt}}^{+T} (-\nabla f_t(\mathbf{x}^+) + \nabla f_t(\mathbf{x}) + \nabla \phi(\mathbf{x}^+) - \nabla \phi(\mathbf{x})) \\ &= 0\end{aligned}$$

Convergence of Newtons Method

bound on $\lambda_t(\mathbf{x}^+)$:

we use $D := D_x = \text{diag}(d_x)$ and $D_+ := D_{x^+} = \text{diag}(d_{x^+})$

$$\begin{aligned}\lambda_t(\mathbf{x}^+)^2 &= \|D_+ A \Delta \mathbf{x}_{\text{nt}}^+\|^2 \\ &\leq \|D_+ A \Delta \mathbf{x}_{\text{nt}}^+\|^2 + \|D_+ A \Delta \mathbf{x}_{\text{nt}}^+ + (I - D_+^{-1} D) D A \Delta \mathbf{x}_{\text{nt}}\|^2 \\ &= \|(I - D_+^{-1} D) D A \Delta \mathbf{x}_{\text{nt}}\|^2 \\ &= \|(I - D_+^{-1} D)^2 \vec{\mathbf{1}}\|^2 \\ &\leq \|(I - D_+^{-1} D) \vec{\mathbf{1}}\|^4 \\ &= \|D A \Delta \mathbf{x}_{\text{nt}}\|^4 \\ &= \lambda_t(\mathbf{x})^4\end{aligned}$$

The second inequality follows from $\sum_i y_i^4 \leq (\sum_i y_i^2)^2$

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If $\lambda_t(x)$ is large we do not have a guarantee.

Try to avoid this case!!!

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Path-following Methods

Try to slowly travel along the central path.

Algorithm 1 PathFollowing

- 1: start at analytic center
- 2: **while** solution not good enough **do**
- 3: make step to improve objective function
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Short Step Barrier Method

simplifying assumptions:

- ▶ a first central point $x^*(t_0)$ is given
- ▶ $x^*(t)$ is computed exactly in each iteration

ϵ is approximation we are aiming for

start at $t = t_0$, repeat until $m/t \leq \epsilon$

- ▶ compute $x^*(\mu t)$ using Newton starting from $x^*(t)$
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where $\mu = 1 + 1/(2\sqrt{m})$

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gradient of f_{t+} at $(x = x^*(t))$

$$\begin{aligned}\nabla f_{t+}(x) &= \nabla f_t(x) + (\mu - 1)tc \\ &= -(\mu - 1)A^T D_x \vec{1}\end{aligned}$$

This holds because $0 = \nabla f_t(x) = tc + A^T D_x \vec{1}$.

The Newton decrement is

$$\begin{aligned}\lambda_{t+}(x)^2 &= \nabla f_{t+}(x)^T H^{-1} \nabla f_{t+}(x) \\ &= (\mu - 1)^2 \vec{1}^T B (B^T B)^{-1} B^T \vec{1} \quad B = D_x^T A \\ &\leq (\mu - 1)^2 m \\ &= 1/4\end{aligned}$$

This means we are in the range of quadratic convergence!!!

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Number of Iterations

the number of Newton iterations per outer iteration is very small; in practise only 1 or 2

Number of outer iterations:

We need $t_k = \mu^k t_0 \geq m/\epsilon$. This holds when

$$k \geq \frac{\log(m/(\epsilon t_0))}{\log(\mu)}$$

We get a bound of

$$\mathcal{O}\left(\sqrt{m} \log \frac{m}{\epsilon t_0}\right)$$

We show how to get a starting point with $t_0 = 1/2^L$. Together with $\epsilon \approx 2^{-L}$ we get $\mathcal{O}(L\sqrt{m})$ iterations.

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Damped Newton Method

For $x \in P^\circ$ and direction $v \neq 0$ define

$$\sigma_x(v) := \max_i \frac{a_i^T v}{s_i(x)}$$

Observation:

$$x + \alpha v \in P \quad \text{for } \alpha \in \{0, 1/\sigma_x(v)\}$$

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Suppose that we move from x to $x + \alpha v$. The linear estimate says that $f_t(x)$ should change by $\nabla f_t(x)^T \alpha v$.

The following argument shows that f_t is well behaved. For small α the reduction of $f_t(x)$ is close to linear estimate.

$$f_t(x + \alpha v) - f_t(x) = \alpha \nabla f_t(x)^T v + \phi(x + \alpha v) - \phi(x)$$

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Define $w_i = a_i^T v / s_i(x)$ and $\sigma = \max_i w_i$. Then

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$$\begin{aligned} &\leq -\sum_i \frac{w_i^2}{\sigma^2} (\alpha\sigma + \log(1 - \alpha\sigma)) \\ &= -\frac{1}{\sigma^2} \|v\|_{H_x}^2 (\alpha\sigma + \log(1 - \alpha\sigma)) \end{aligned}$$

Damped Newton Iteration:

In a damped Newton step we choose

$$x_+ = x + \frac{1}{1 + \sigma_x(\Delta x_{nt})} \Delta x_{nt}$$

Damped Newton Method

Define $w_i = a_i^T v / s_i(x)$ and $\sigma = \max_i w_i$. Then

$$\begin{aligned} &f_t(x + \alpha v) - f_t(x) - \nabla f_t(x)^T \alpha v \\ &= -\sum_i (\alpha w_i + \log(1 - \alpha w_i)) \\ &\leq -\sum_{w_i > 0} (\alpha w_i + \log(1 - \alpha w_i)) + \sum_{w_i \leq 0} \frac{\alpha^2 w_i^2}{2} \\ &\leq -\sum_{w_i > 0} \frac{w_i^2}{\sigma^2} (\alpha\sigma + \log(1 - \alpha\sigma)) + \frac{(\alpha\sigma)^2}{2} \sum_{w_i \leq 0} \frac{w_i^2}{\sigma^2} \end{aligned}$$

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Theorem:

In a damped Newton step the cost decreases by at least

$$\lambda_t(x) - \log(1 + \lambda_t(x))$$

Proof: The decrease in cost is

$$-\alpha \nabla f_t(x)^T v + \frac{1}{\sigma^2} \|v\|_{H_x}^2 (\alpha\sigma + \log(1 - \alpha\sigma))$$

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$$\begin{aligned} &\geq \lambda_t(\mathbf{x}) - \log(1 + \lambda_t(\mathbf{x})) \\ &\geq 0.09 \end{aligned}$$

for $\lambda_t(\mathbf{x}) \geq 0.5$

Centering Algorithm:

Input: precision δ ; starting point \mathbf{x}

1. compute $\Delta\mathbf{x}_{nt}$ and $\lambda_t(\mathbf{x})$
2. if $\lambda_t(\mathbf{x}) \leq \delta$ return \mathbf{x}
3. set $\mathbf{x} := \mathbf{x} + \alpha\Delta\mathbf{x}_{nt}$ with

$$\alpha = \begin{cases} \frac{1}{1 + \sigma_x(\Delta\mathbf{x}_{nt})} & \lambda_t \geq 1/2 \\ 1 & \text{otw.} \end{cases}$$

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Centering

Lemma 3

The centering algorithm starting at x_0 reaches a point with $\lambda_t(x) \leq \delta$ after

$$\frac{f_t(x_0) - \min_y f_t(y)}{0.09} + \mathcal{O}(\log \log(1/\delta))$$

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This can be very, very slow...

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How to get close to analytic center?

Let $P = \{Ax \leq b\}$ be our (**feasible**) polyhedron, and x_0 a feasible point.

We change $b \rightarrow b + \frac{1}{\lambda} \cdot \vec{1}$, where $L = \langle A \rangle + \langle b \rangle + \langle c \rangle$ (**encoding length**) and $\lambda = 2^{2L}$. Recall that a basis is feasible in the old LP iff it is feasible in the new LP.

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Lemma [without proof]

The inverse of a matrix M can be represented with rational numbers that have denominators $z_{ij} = \det(M)$.

For two basis solutions $x_B, x_{\bar{B}}$, the cost-difference $c^T x_B - c^T x_{\bar{B}}$ can be represented by a rational number that has denominator $z = \det(A_B) \cdot \det(A_{\bar{B}}) \cdot \lambda$.

This means that in the perturbed LP it is sufficient to decrease the duality gap to $1/2^{4L}$ (i.e., $t \approx 2^{4L}$). This means the previous analysis essentially also works for the perturbed LP.

For a point x from the polytope (not necessarily BFS) the objective value $\bar{c}^T x$ is at most $n2^M 2^L$, where $M \leq L$ is the encoding length of the largest entry in \bar{c} .

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Start at x_0 .

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$x_0 = x^*(1)$ is point on central path for \hat{c} and $t = 1$.

You can travel the central path in both directions. Go towards 0 until $t \approx 1/2^{\Omega(L)}$. This requires $O(\sqrt{m}L)$ outer iterations.

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Clearly,

$$t \cdot \hat{c}^T x_{\hat{c}} + \phi(x_{\hat{c}}) \leq t \cdot \hat{c}^T x_c + \phi(x_c)$$

The different between $f_t(x_{\hat{c}})$ and $f_t(x_c)$ is

$$\begin{aligned} & t c^T x_{\hat{c}} + \phi(x_{\hat{c}}) - t c^T x_c - \phi(x_c) \\ & \leq t(c^T x_{\hat{c}} + \hat{c}^T x_c - \hat{c}^T x_{\hat{c}} - c^T x_c) \\ & \leq 4tn2^{3L} \end{aligned}$$

For $t = 1/2^{\Omega(L)}$ the last term becomes constant. Hence, using damped Newton we can move from $x_{\hat{c}}$ to x_c quickly.

In total for this analysis we require $\mathcal{O}(\sqrt{mL})$ outer iterations for the whole algorithm.

One iteration can be implemented in $\tilde{\mathcal{O}}(m^3)$ time.

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Clearly,

$$t \cdot \hat{c}^T x_{\hat{c}} + \phi(x_{\hat{c}}) \leq t \cdot \hat{c}^T x_c + \phi(x_c)$$

The different between $f_t(x_{\hat{c}})$ and $f_t(x_c)$ is

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For $t = 1/2^{\Omega(L)}$ the last term becomes constant. Hence, using damped Newton we can move from $x_{\hat{c}}$ to x_c quickly.

In total for this analysis we require $\mathcal{O}(\sqrt{mL})$ outer iterations for the whole algorithm.

One iteration can be implemented in $\tilde{\mathcal{O}}(m^3)$ time.

How to get close to analytic center?

Start at x_0 .

Choose $\hat{c} := -\nabla \phi(x)$.

$x_0 = x^*(1)$ is point on central path for \hat{c} and $t = 1$.

You can travel the central path in both directions. Go towards 0 until $t \approx 1/2^{\Omega(L)}$. This requires $\mathcal{O}(\sqrt{mL})$ outer iterations.

Let $x_{\hat{c}}$ denote this point.

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