BUNDLE METHODS FOR STOCHASTIC PROGRAMS PROXIMAL BUNDLE METHOD

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GENERAL FORMULATION

In this part of the course we will focus on efficient optimization methods to solve convex programs of the form

$$\min f(x)$$
 s.t. $x \in X$,

with

- $\blacktriangleright \ f: \Re^n \to \Re$ a convex but nonsmooth function
- $X \subset \Re^n$ a convex set (e.g. $X = \{x \in \Re^n_+ : Ax = b\}, X = \Re^n$)

This formulation covers many practical optimization problems, for instance

- ▶ Two-stage stochastic programming problems
- Multistage stochastic programming problems

TWO-STAGE STOCHASTIC LINEAR PROGRAMMING

In two-stage stochastic linear programming problems with finitely many scenarios $\xi^i = (q^i, T^i, W^i, h^i)$ we wish to solve the high dimensional LP

$$\begin{cases} \min & c^{\top}x + \sum_{i=1}^{N} p_i[q^{i^{\top}}y^i] \\ \text{s.t.} & Ax = b, \ x \ge 0 \\ & T^ix + W^iy^i = h^i, \ y^i \ge 0, \quad i = 1, \dots, N \end{cases}$$



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TWO-STAGE DECOMPOSITION

min
$$f(x)$$
 s.t. $x \in X$, with $f(x) := c^{\top} x + \sum_{i=1}^{N} p_i Q(x, \xi^i)$,

$$Q(x,\xi) = \begin{cases} \min & q^{\top}y \\ \text{s.t.} & Wy = h - Tx \\ & y \ge 0 \,. \end{cases} \text{ and } X := \{x \in \Re_{+}^{n} : Ax = b\}$$

We know that $g = c - \sum_{i=1}^{N} p_i T^i \pi^i \in \partial f(x)$, where π^i is a dual solution of $Q(x,\xi^i)$

Multistage stochastic linear programs

$$\min_{\substack{A_1x_1=b_1\\x_1\geq 0}} c_1^{\top}x_1 + \mathbb{E}\left[\min_{\substack{B_2x_1+A_2x_2=b_2\\x_2\geq 0}} c_2^{\top}x_2 + \mathbb{E}\left[\cdots + \mathbb{E}[\min_{\substack{B_Tx_T-1+A_Tx_T=b_T\\x_T\geq 0}} c_T^{\top}x_T]\right]\right]$$

▶ Some elements of the data $\xi = (c_t, B_t, A_t, b_t)$ depend on uncertainties.

By assuming finitely many scenarios and dualizing the nonantecipativity constraints (that can be written as Gx = 0) we get



Multistage stochastic linear programs

(See Lecture 17)

DUAL PROBLEM

$$\min_u f(u), \quad ext{with} \quad f(u) := -\sum_{i=1}^N D^i(u)$$

$$D^{i}(u) := \begin{cases} \min_{x^{i}} & p_{i} \sum_{t=1}^{T} (c_{t}^{i})^{\top} x_{t}^{i} + u^{\top} G^{i} x^{i} \\ \text{s.t.} & A_{1} x_{1} = b_{1} \\ & B_{t}^{i} x_{t-1}^{i} + A_{t}^{i} x_{t}^{i} = b_{t}^{k}, \ t = 2, \dots, T \\ & x_{t}^{i} \ge 0. \end{cases}$$

Computing f(u) for each given u amounts to solving N LPs.

We know that $g = -Gx(u) \in \partial f(u)$, where $x(u) = (x^1(u), \dots, x^N(u))$ and $x^i(u)$ is a solution of $D^i(u)$

Let's stick with the more compact and general formulation

$$\min f(x) \quad \text{s.t.} \quad x \in X \,,$$

with $f:\Re^n\to\Re$ a convex but nonsmooth function and $X\subset\Re^n$ a convex set.

We'll assume the availability of an oracle providing us with first-order information on f:

In stochastic programming, the oracle should be smart enough to use parallel computing:

- \blacktriangleright the oracle consists of solving N optimization subproblems to compute f(x) and a subgradient g
- most of time dedicate to minimize f is spent in the oracle!

Therefore, subgradient and (pure) cutting-plane methods are not very efficient $^1\ldots$

¹These methods require, in general, many oracle calls. \triangleleft $\rightarrow \triangleleft$

Consider the problem

$$\min_{x \in X} f(x)$$

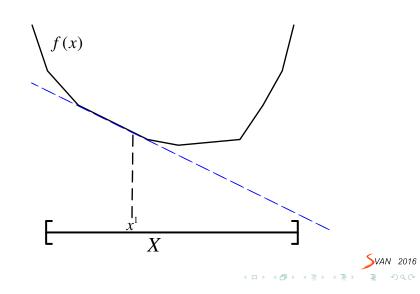
and suppose that X is a compact set.

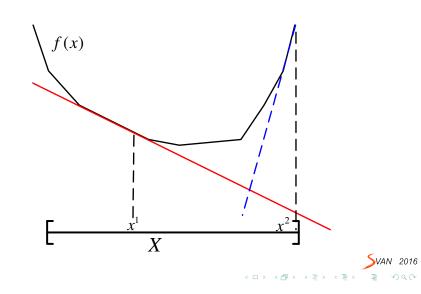
Algorithm

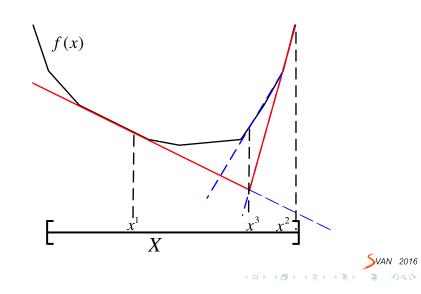
- 1. Given $x_0 \in X$, call the oracle to compute $f(x_0)$ and $g_0 \in \partial f(x_0)$. Set $f_0^{\text{up}} = f(x_0)$ and k = 0
- 2. (iterate) Find $x_{k+1} = \arg\min_{x \in X} \check{f}_k(x)$. Let $f_k^{\text{low}} = \check{f}_k(x_{k+1})$.
- 3. (stopping test) If $f_k^{\text{up}} f^{\text{low}}$ is small enough, stop.
- 4. (oracle) Compute $f(x_{k+1}), g_{k+1} \in \partial f(x_k)$ and set $f_{k+1}^{\text{up}} = \min\{f(x_{k+1}), f_k^{\text{up}}\}.$
- 5. (loop) Set $k \leftarrow k + 1$ and go back to Step 2.

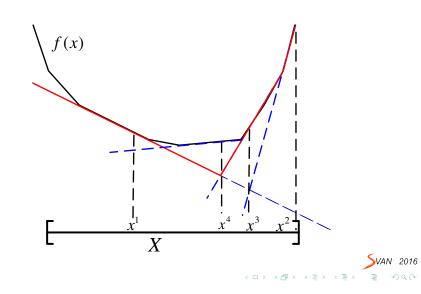
CUTTING-PLANE MODEL

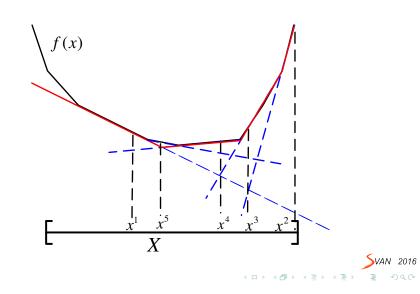
$$\check{f}_{k}(\cdot) = \max_{j=1,\dots,k} \{f(x_{j}) + g_{j}^{\top}(\cdot - x_{j})\}$$
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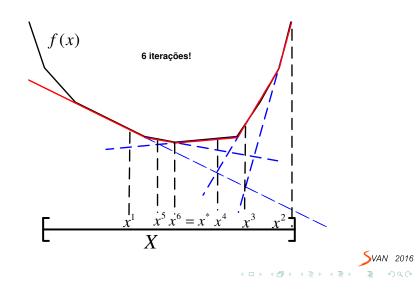












The method requires solving a LP at each iteration

$$x_{k+1} = \arg\min_{x \in X} \check{f}_k(x), \quad \check{f}_k(\cdot) = \max_{j=1,...,k} \{f(x_j) + g_j^{\top}(\cdot - x_j)\}$$

that is equivalent to

$$\begin{cases} \min_{x,r} & r \\ s.t. & f(x_j) + g_j^{\top}(x - x_j) \le r, \quad j = 1, \dots, k \\ & x \in X, \ r \in \Re. \end{cases}$$

A new constraint is added at each iteration!

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 $PROS \times CONS$

- only computes a single subgradient per iteration
- easy to code
- \square easy and reliable stopping test
- $f(x_{k+1}) \not\leq f(x_k)$ (it is not a descent method)
- \square instable and has low convergence rate
- \mathbb{R} requires compactness of the feasible set
- \square doesn't exploit good starting points
- **k** subproblem becomes heavier and heavier...

The *Regularized Decomposition Method* (1986) for 2-SLP address some of the above drawbacks.

Regularized Decomposition Method is just a particular case of (proximal) Bundle Methods!

BUNDLE METHODS

MAIN INGREDIENTS

- (I) a convex model $f_k^M \leq f$ (eg. cutting-plane model)
- (II) a stability center \hat{x}_k (eg.: the best point so far)
- (III) a parameter t_k (or f_k^{lev}) to be updated at every iteration

The next trial point x_{k+1} of a bundle method depends on the above 3 ingredients, whose organization define different methods:

PROXIMAL BUNDLE METHOD $(t_k > 0)$

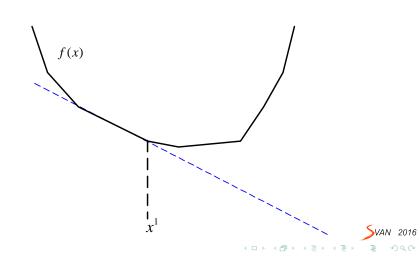
$$x_{k+1} := \arg\min\left\{f_k^M(x) + \frac{1}{2t_k} \|x - \hat{x}_k\|^2 : x \in X\right\}.$$

LEVEL BUNDLE METHOD $(f_k^{\text{lev}} \in \Re)$

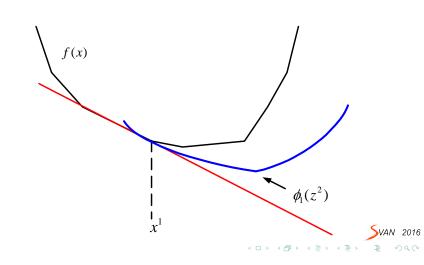
$$x_{k+1} := \arg\min\left\{\frac{1}{2}\|x - \hat{x}_k\|^2 : f_k^M(x) \le f_k^{\text{lev}}, x \in X\right\}.$$

Today we focus on proximal bundle method!

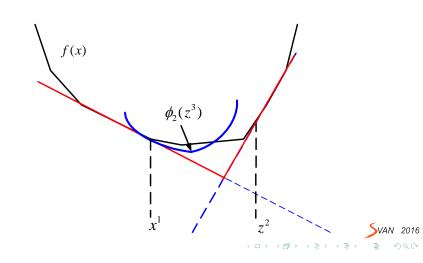
$$f^{M} \equiv \check{f}, \qquad x_{k+1} := \arg\min\left\{\check{f}_{k}(x) + \frac{1}{2t_{k}}\|x - \hat{x}_{k}\|^{2} : x \in X\right\}$$



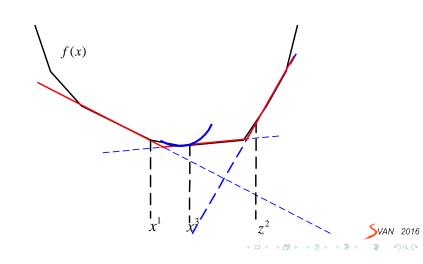
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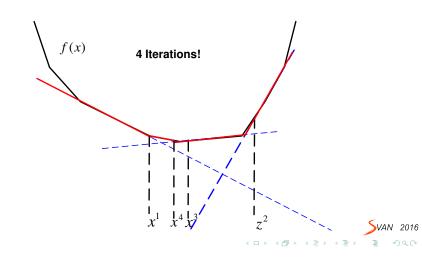
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$\mathrm{Pros}\,\times\,\mathrm{Cons}$

- \square only computes a single subgradient per iteration
- \square easy and reliable stopping test
- 🕼 stable
- does not require X to be compact
- \square exploit good-quality initial points
- subproblem defining x_{k+1} can be kept small



$\mathrm{Pros}\,\times\,\mathrm{Cons}$

- \square only computes a single subgradient per iteration
- \square easy and reliable stopping test
- stable
- does not require X to be compact
- it is a descent method
- exploit good-quality initial points
- subproblem defining x_{k+1} can be kept small
- **↓** convergence analysis is more involving...



Let's consider a more economical model:

$$f_k^M(x) := \max_{j \in \mathcal{B}_k} \{ f(x_j) + g_j^{\top} (x - x_j) \}$$

- ▶ The cutting-plane method takes $\mathcal{B}_k := \{1, 2, ..., k\}$. We will consider $\mathcal{B}_k \subset \{1, 2, ..., k\}$ (or something a bit different)
- ▶ The method generates a sequence of trial points $\{x_k\} \subset X$ by solving a QP:

$$x_{k+1} := \arg\min\left\{f_k^M(x) + \frac{1}{2t_k}\|x - \hat{x}_k\|^2 : x \in X\right\}.$$



Solving the QP subproblem

The QP

$$\min\left\{f_{k}^{M}(x) + \frac{1}{2t_{k}}\|x - \hat{x}_{k}\|^{2} : x \in X\right\}$$

can be rewritten as

$$\begin{cases} \min_{x,r} & r + \frac{1}{2t_k} \|x - \hat{x}_k\|^2 \\ \text{s.a} & f(x_j) + g_j^\top (x - x_j) \le r, \quad j \in \mathcal{B}_k \\ & x \in X, \ r \in \Re. \end{cases}$$

We can apply specialized softwares.



$$x_{k+1} := \arg\min\left\{f_k^M(x) + \frac{1}{2t_k} \|x - \hat{x}_k\|^2 : x \in X\right\}.$$

A rule decide when to update the stability center \hat{x}_k . Such rule depends on the predicted decrease by the model f_k^M

$$v_k = f(\hat{x}_k) - f_k^M(x_{k+1})$$

and a constant $\kappa \in (0, 1)$:

• Serious step: if $f(x_{k+1}) \leq f(\hat{x}_k) - \kappa v_k$, then

$$\hat{x}_{k+1} \leftarrow x_{k+1}$$

• Null step: if $f(x_{k+1}) > f(\hat{x}_k) - \kappa v_k$, then

$$\hat{x}_{k+1} \leftarrow \hat{x}_k$$

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The serious-step sequence $\{\hat{x}_k\}$ is a subsequence of $\{x_k\}$

NEXT ITERATE

LEMMA

Suppose that X is a polyhedron or $ri(X) \neq \emptyset$. Then

$$x_{k+1} = \hat{x}_k - t_k \hat{g}_k \quad com \quad \hat{g}_k = p_f^k + p_X^k ,$$

where $p_f^k \in \partial f_k^M(x_{k+1})$ and $p_X^k \in \partial i_X(x_{k+1})$. (*i_X* is the indicator function of *X*.)

Furthermore, the affine function

$$f_{k^a}^L(x) := f_k^M(x_{k+1}) + \langle \hat{g}_k, x - x_{k+1} \rangle$$

is a lower approximation for the model f_k^M :

$$f_{k^a}^L(x) \le f_k^M(x) \quad \forall \ x \in X.$$



OPTIMALITY MEASURE

Propositon

Let the predicted decrease and aggregate linearization error defined by

$$v_k := f(\hat{x}_k) - f_k^M(x_{k+1}) \text{ and } \hat{e}_k := f(\hat{x}_k) - f_{k^a}^L(\hat{x}_k).$$

Then,

$$\hat{e}_k \ge 0$$
, $\hat{e}_k + t_k \|\hat{g}_k\|^2 = v_k \ge 0$ for all k.

Furthermore

$$f(\hat{x}_k) \le f(x) + \hat{e}_k + \|\hat{g}_k\| \|\hat{x}_k - x\|$$
 for all $x \in X$ and k .

If $(\hat{e}_k, \hat{g}_k) = 0$, then \hat{x}_k is solution to the problem

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Algorithm: proximal bundle method

$$f_k^M(x) = \max_{j \in \mathcal{B}_k} \{ f(x_j) + g_j^\top (x - x_j) \}, \quad x_{k+1} = \arg\min\left\{ f_k^M(x) + \frac{1}{2t_k} \|x - \hat{x}_k\|^2 : x \in X \right\}$$

- Step 0. Choose $\kappa \in (0, 1)$, $t_1 \ge t_{\min} > 0$, $x_1 \in X$ and tolerance tol > 0. Call the oracle to compute $(f(x_1), g_1)$. Define $\hat{x}_1 \leftarrow x_1, k \leftarrow 1, \mathcal{B}_1 \leftarrow \{1\},$
- Step 1. Solve the QP to obtain x_{k+1} . Define $\hat{g}_k \leftarrow (\hat{x}_k x_{k+1})/t_k$, $v_k \leftarrow f(\hat{x}_k) - \check{f}_k(x_{k+1})$, and $\hat{e}_k \leftarrow v_k - t_k \|\hat{g}_k\|^2$
- **Step 2.** If $\hat{e}_k \leq \text{tol}$ and $\|\hat{g}_k\| \leq \text{tol}$, stop: \hat{x}_k is an approximate solution
- Step 3. Call the oracle to obtain $(f(x_{k+1}), g_{k+1})$ Serious step. If $f(x_{k+1}) \leq f(\hat{x}_k) - \kappa v_k$, then $\hat{x}_{k+1} \leftarrow x_{k+1}$ and choose $t_{k+1} \geq t_k$ Null step. Otherwise, define $\hat{x}_{k+1} \leftarrow \hat{x}_k$ and choose $t_{k+1} \in [t_{\min}, t_k]$

Step 4. Choose
$$\mathcal{B}_{k+1} \supset \{k+1, k^a\}$$

Set $k \leftarrow k+1$ and go back to Step 1

Some comments

▶ Only 2 linearizations are required: f_k^L and $f_{k^a}^L$, i.e.,

$$\mathcal{B}_{k+1} = \{k+1, k^a\} \quad \text{suffices!}$$

- the prox-parameter t_k is non-increasing along null steps
- ▶ a simple heuristic to update the prox-parameter is the following
 - $\blacktriangleright \text{ compute } t_{\mathtt{aux}} := t_k \left(1 + \frac{(g_{k+1} g_k)^\top (x_{k+1} x_k)}{\|g_{k+1} g_k\|^2} \right)$
 - ▶ if null step: $t_{k+1} \leftarrow \min\{t_k, \max\{t_{aux}, t_k/2, t_{\min}\}\}$
 - ▶ if serious step: $t_{k+1} \leftarrow \max\{t_k, \min\{t_{aux}, 10t_k\}\}$
- \blacktriangleright it is advisable to consider different tolerances for the measures \hat{e}_k and \hat{g}_k

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- the sequence $\{f(\hat{x}_k)\}$ is non-increasing
- any accumulation point of $\{\hat{x}_k\}$ is a solution to the problem