Alternating Direction Method of Multipliers

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source:

Distributed Optimization and Statistical Learning via the Alternating Direction Method of Multipliers (Boyd, Parikh, Chu, Peleato, Eckstein)

Goals

robust methods for

- ► arbitrary-scale optimization
 - machine learning/statistics with huge data-sets
 - dynamic optimization on large-scale network
- decentralized optimization
 - devices/processors/agents coordinate to solve large problem, by passing relatively small messages

Outline

Dual decomposition

Method of multipliers

Alternating direction method of multipliers

Common patterns

Examples

Consensus and exchange

Conclusions

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Dual problem

convex equality constrained optimization problem

 $\begin{array}{ll} \text{minimize} & f(x) \\ \text{subject to} & Ax = b \end{array}$

- ▶ Lagrangian: $L(x,y) = f(x) + y^T(Ax b)$
- dual function: $g(y) = \inf_x L(x, y)$
- dual problem: maximize g(y)
- recover $x^{\star} = \operatorname{argmin}_{x} L(x, y^{\star})$

Dual decomposition

Dual ascent

- \blacktriangleright gradient method for dual problem: $y^{k+1} = y^k + \alpha^k \nabla g(y^k)$
- $\nabla g(y^k) = A\tilde{x} b$, where $\tilde{x} = \operatorname{argmin}_x L(x, y^k)$
- dual ascent method is

$$egin{array}{rcl} x^{k+1} &:= & \mathrm{argmin}_x \, L(x,y^k) & // \, x\mbox{-minimization} \ y^{k+1} &:= & y^k + lpha^k (Ax^{k+1} - b) & // \, dual \, update \end{array}$$

works, with lots of strong assumptions

Dual decomposition

► suppose *f* is separable:

$$f(x) = f_1(x_1) + \dots + f_N(x_N), \quad x = (x_1, \dots, x_N)$$

▶ then L is separable in x: $L(x,y) = L_1(x_1,y) + \cdots + L_N(x_N,y) - y^T b$,

$$L_i(x_i, y) = f_i(x_i) + y^T A_i x_i$$

 \blacktriangleright x-minimization in dual ascent splits into N separate minimizations

$$x_i^{k+1} := \operatorname*{argmin}_{x_i} L_i(x_i, y^k)$$

which can be carried out in parallel

Dual decomposition

Dual decomposition

▶ dual decomposition (Everett, Dantzig, Wolfe, Benders 1960–65)

$$\begin{aligned} x_i^{k+1} &:= & \operatorname{argmin}_{x_i} L_i(x_i, y^k), \quad i = 1, \dots, N \\ y^{k+1} &:= & y^k + \alpha^k (\sum_{i=1}^N A_i x_i^{k+1} - b) \end{aligned}$$

- ▶ scatter y^k ; update x_i in parallel; gather $A_i x_i^{k+1}$
- ► solve a large problem
 - by iteratively solving subproblems (in parallel)
 - dual variable update provides coordination
- ▶ works, with lots of assumptions; often slow

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Method of multipliers

Method of multipliers

- a method to robustify dual ascent
- use augmented Lagrangian (Hestenes, Powell 1969), $\rho > 0$

$$L_{\rho}(x,y) = f(x) + y^{T}(Ax - b) + (\rho/2) ||Ax - b||_{2}^{2}$$

▶ method of multipliers (Hestenes, Powell; analysis in Bertsekas 1982)

$$x^{k+1} := \operatorname{argmin}_{x} L_{\rho}(x, y^{k})$$
$$y^{k+1} := y^{k} + \rho(Ax^{k+1} - b)$$

(note specific dual update step length ρ)

Method of multipliers dual update step

• optimality conditions (for differentiable *f*):

$$Ax^{\star} - b = 0, \qquad \nabla f(x^{\star}) + A^T y^{\star} = 0$$

(primal and dual feasibility)

• since x^{k+1} minimizes $L_{\rho}(x, y^k)$

$$0 = \nabla_x L_{\rho}(x^{k+1}, y^k) = \nabla_x f(x^{k+1}) + A^T (y^k + \rho(Ax^{k+1} - b)) = \nabla_x f(x^{k+1}) + A^T y^{k+1}$$

▶ dual update $y^{k+1} = y^k + \rho(x^{k+1} - b)$ makes (x^{k+1}, y^{k+1}) dual feasible

• primal feasibility achieved in limit: $Ax^{k+1} - b \rightarrow 0$

Method of multipliers

Method of multipliers

(compared to dual decomposition)

- ▶ good news: converges under much more relaxed conditions (f can be nondifferentiable, take on value +∞, ...)
- ► bad news: quadratic penalty destroys splitting of the x-update, so can't do decomposition

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- a method
 - with good robustness of method of multipliers
 - which can support decomposition
- "robust dual decomposition" or "decomposable method of multipliers"
- ▶ proposed by Gabay, Mercier, Glowinski, Marrocco in 1976

Alternating direction method of multipliers

• ADMM problem form (with f, g convex)

 $\begin{array}{ll} \mbox{minimize} & f(x) + g(z) \\ \mbox{subject to} & Ax + Bz = c \end{array}$

- two sets of variables, with separable objective

•
$$L_{\rho}(x,z,y) = f(x) + g(z) + y^T (Ax + Bz - c) + (\rho/2) \|Ax + Bz - c\|_2^2$$

► ADMM:

- \blacktriangleright if we minimized over x and z jointly, reduces to method of multipliers
- ▶ instead, we do one pass of a Gauss-Seidel method
- \blacktriangleright we get splitting since we minimize over x with z fixed, and vice versa

ADMM and optimality conditions

optimality conditions (for differentiable case):

- primal feasibility: Ax + Bz c = 0
- dual feasibility: $\nabla f(x) + A^T y = 0$, $\nabla g(z) + B^T y = 0$

 \blacktriangleright since z^{k+1} minimizes $L_{\rho}(x^{k+1},z,y^k)$ we have

$$0 = \nabla g(z^{k+1}) + B^T y^k + \rho B^T (Ax^{k+1} + Bz^{k+1} - c)$$

= $\nabla g(z^{k+1}) + B^T y^{k+1}$

- \blacktriangleright so with ADMM dual variable update, $(x^{k+1},z^{k+1},y^{k+1})$ satisfies second dual feasibility condition
- \blacktriangleright primal and first dual feasibility are achieved as $k \rightarrow \infty$

ADMM with scaled dual variables

► combine linear and quadratic terms in augmented Lagrangian

$$\begin{split} L_{\rho}(x,z,y) &= f(x) + g(z) + y^{T}(Ax + Bz - c) + (\rho/2) \|Ax + Bz - c\|_{2}^{2} \\ &= f(x) + g(z) + (\rho/2) \|Ax + Bz - c + u\|_{2}^{2} + \text{const.} \end{split}$$

with $u^{k} = (1/\rho)y^{k}$

► ADMM (scaled dual form):

$$\begin{aligned} x^{k+1} &:= \arg \min_{x} \left(f(x) + (\rho/2) \| Ax + Bz^{k} - c + u^{k} \|_{2}^{2} \right) \\ z^{k+1} &:= \arg \min_{z} \left(g(z) + (\rho/2) \| Ax^{k+1} + Bz - c + u^{k} \|_{2}^{2} \right) \\ u^{k+1} &:= u^{k} + (Ax^{k+1} + Bz^{k+1} - c) \end{aligned}$$

Convergence

- ► assume (very little!)
 - $f,\ g$ convex, closed, proper
 - L_0 has a saddle point
- ► then ADMM converges:
 - iterates approach feasibility: $Ax^k + Bz^k c \rightarrow 0$
 - objective approaches optimal value: $f(x^k) + g(z^k) \rightarrow p^\star$

Related algorithms

- operator splitting methods (Douglas, Peaceman, Rachford, Lions, Mercier, ... 1950s, 1979)
- proximal point algorithm (Rockafellar 1976)
- ▶ Dykstra's alternating projections algorithm (1983)
- ► Spingarn's method of partial inverses (1985)
- ► Rockafellar-Wets progressive hedging (1991)
- ▶ proximal methods (Rockafellar, many others, 1976–present)
- Bregman iterative methods (2008–present)
- ▶ most of these are special cases of the proximal point algorithm

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- x-update step requires minimizing f(x) + (ρ/2) ||Ax v||₂² (with v = Bz^k − c + u^k, which is constant during x-update)
- ▶ similar for *z*-update
- several special cases come up often
- can simplify update by exploiting structure in these cases

Decomposition

► suppose *f* is block-separable,

$$f(x) = f_1(x_1) + \dots + f_N(x_N), \qquad x = (x_1, \dots, x_N)$$

- A is conformably block separable: $A^T A$ is block diagonal
- ▶ then x-update splits into N parallel updates of x_i

Proximal operator

• consider x-update when A = I

$$x^{+} = \underset{x}{\operatorname{argmin}} \left(f(x) + (\rho/2) \|x - v\|_{2}^{2} \right) = \mathbf{prox}_{f,\rho}(v)$$

► some special cases:

$$\begin{split} f &= I_C \text{ (indicator fct. of set } C \text{)} \quad x^+ := \Pi_C(v) \text{ (projection onto } C \text{)} \\ f &= \lambda \| \cdot \|_1 \text{ (} \ell_1 \text{ norm} \text{)} \qquad x_i^+ := S_{\lambda/\rho}(v_i) \text{ (soft thresholding)} \\ (S_a(v) &= (v-a)_+ - (-v-a)_+ \text{)} \end{split}$$

Common patterns

Quadratic objective

•
$$f(x) = (1/2)x^T P x + q^T x + r$$

•
$$x^+ := (P + \rho A^T A)^{-1} (\rho A^T v - q)$$

use matrix inversion lemma when computationally advantageous

$$(P + \rho A^T A)^{-1} = P^{-1} - \rho P^{-1} A^T (I + \rho A P^{-1} A^T)^{-1} A P^{-1}$$

- (direct method) cache factorization of $P + \rho A^T A$ (or $I + \rho A P^{-1} A^T$)
- ▶ (iterative method) warm start, early stopping, reducing tolerances

Smooth objective

- ► *f* smooth
- can use standard methods for smooth minimization
 - gradient, Newton, or quasi-Newton
 - preconditionned CG, limited-memory BFGS (scale to very large problems)
- ► can exploit
 - warm start
 - early stopping, with tolerances decreasing as ADMM proceeds

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Constrained convex optimization

consider ADMM for generic problem

 $\begin{array}{ll} \text{minimize} & f(x) \\ \text{subject to} & x \in \mathcal{C} \end{array}$

 \blacktriangleright ADMM form: take g to be indicator of ${\mathcal C}$

 $\begin{array}{ll} \mbox{minimize} & f(x) + g(z) \\ \mbox{subject to} & x-z = 0 \end{array}$

► algorithm:

$$\begin{aligned} x^{k+1} &:= & \operatorname*{argmin}_{x} \left(f(x) + (\rho/2) \| x - z^k + u^k \|_2^2 \right) \\ z^{k+1} &:= & \Pi_{\mathcal{C}} (x^{k+1} + u^k) \\ u^{k+1} &:= & u^k + x^{k+1} - z^{k+1} \end{aligned}$$

Lasso

► lasso problem:

minimize
$$(1/2) ||Ax - b||_2^2 + \lambda ||x||_1$$

► ADMM form:

minimize
$$(1/2) \|Ax - b\|_2^2 + \lambda \|z\|_1$$

subject to $x - z = 0$

► ADMM:

$$\begin{aligned} x^{k+1} &:= (A^T A + \rho I)^{-1} (A^T b + \rho z^k - y^k) \\ z^{k+1} &:= S_{\lambda/\rho} (x^{k+1} + y^k/\rho) \\ y^{k+1} &:= y^k + \rho (x^{k+1} - z^{k+1}) \end{aligned}$$

Lasso example

- ► example with dense A ∈ R^{1500×5000} (1500 measurements; 5000 regressors)
- computation times

factorization (same as ridge regression)1.3ssubsequent ADMM iterations0.03slasso solve (about 50 ADMM iterations)2.9sfull regularization path $(30 \lambda's)$ 4.4s

not bad for a very short Matlab script

Sparse inverse covariance selection

- S: empirical covariance of samples from N(0, Σ), with Σ⁻¹ sparse (*i.e.*, Gaussian Markov random field)
- estimate Σ^{-1} via ℓ_1 regularized maximum likelihood

minimize $\operatorname{Tr}(SX) - \log \det X + \lambda \|X\|_1$

▶ methods: COVSEL (Banerjee et al 2008), graphical lasso (FHT 2008)

Sparse inverse covariance selection via ADMM

► ADMM form:

minimize $\operatorname{Tr}(SX) - \log \det X + \lambda \|Z\|_1$ subject to X - Z = 0

► ADMM:

$$\begin{aligned} X^{k+1} &:= & \operatorname*{argmin}_{X} \left(\mathbf{Tr}(SX) - \log \det X + (\rho/2) \| X - Z^k + U^k \|_F^2 \right) \\ Z^{k+1} &:= & S_{\lambda/\rho}(X^{k+1} + U^k) \\ U^{k+1} &:= & U^k + (X^{k+1} - Z^{k+1}) \end{aligned}$$

Analytical solution for X-update

- \blacktriangleright compute eigendecomposition $\rho(Z^k-U^k)-S=Q\Lambda Q^T$
- form diagonal matrix \tilde{X} with

$$\tilde{X}_{ii} = \frac{\lambda_i + \sqrt{\lambda_i^2 + 4\rho}}{2\rho}$$

$$\blacktriangleright \ \text{let} \ X^{k+1} := Q \tilde{X} Q^T$$

► cost of *X*-update is an eigendecomposition

Sparse inverse covariance selection example

$\blacktriangleright~\Sigma^{-1}~{\rm is}~1000\times1000$ with $10^4~{\rm nonzeros}$

- graphical lasso (Fortran): 20 seconds 3 minutes
- ADMM (Matlab): 3 10 minutes
- (depends on choice of λ)
- very rough experiment, but with no special tuning, ADMM is in ballpark of recent specialized methods
- ► (for comparison, COVSEL takes 25+ min when ∑⁻¹ is a 400 × 400 tridiagonal matrix)

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Consensus optimization

 \blacktriangleright want to solve problem with N objective terms

minimize
$$\sum_{i=1}^{N} f_i(x)$$

- e.g., f_i is the loss function for *i*th block of training data

ADMM form:

minimize
$$\sum_{i=1}^{N} f_i(x_i)$$

subject to $x_i - z = 0$

- x_i are local variables
- -z is the global variable
- $x_i z = 0$ are *consistency* or *consensus* constraints
- can add regularization using a g(z) term

Consensus and exchange

Consensus optimization via ADMM

•
$$L_{\rho}(x,z,y) = \sum_{i=1}^{N} \left(f_i(x_i) + y_i^T(x_i-z) + (\rho/2) \|x_i-z\|_2^2 \right)$$

► ADMM:

$$\begin{aligned} x_i^{k+1} &:= & \operatorname*{argmin}_{x_i} \left(f_i(x_i) + y_i^{kT}(x_i - z^k) + (\rho/2) \|x_i - z^k\|_2^2 \right) \\ z^{k+1} &:= & \frac{1}{N} \sum_{i=1}^N \left(x_i^{k+1} + (1/\rho) y_i^k \right) \\ y_i^{k+1} &:= & y_i^k + \rho(x_i^{k+1} - z^{k+1}) \end{aligned}$$

• with regularization, averaging in z update is followed by $\mathbf{prox}_{a,\rho}$

Consensus optimization via ADMM

• using $\sum_{i=1}^{N} y_i^k = 0$, algorithm simplifies to

$$\begin{aligned} x_i^{k+1} &:= \arg\min_{x_i} \left(f_i(x_i) + y_i^{kT}(x_i - \overline{x}^k) + (\rho/2) \|x_i - \overline{x}^k\|_2^2 \right) \\ y_i^{k+1} &:= y_i^k + \rho(x_i^{k+1} - \overline{x}^{k+1}) \end{aligned}$$

where
$$\overline{x}^k = (1/N) \sum_{i=1}^N x_i^k$$

- ▶ in each iteration
 - gather \boldsymbol{x}_i^k and average to get $\overline{\boldsymbol{x}}^k$
 - scatter the average \overline{x}^k to processors
 - update y_i^k locally (in each processor, in parallel)
 - update x_i locally

Statistical interpretation

- ▶ f_i is negative log-likelihood for parameter x given *i*th data block
- x_i^{k+1} is MAP estimate under prior $\mathcal{N}(\overline{x}^k + (1/\rho)y_i^k, \rho I)$
- prior mean is previous iteration's consensus shifted by 'price' of processor i disagreeing with previous consensus
- ▶ processors only need to support a Gaussian MAP method
 - type or number of data in each block not relevant
 - consensus protocol yields global maximum-likelihood estimate

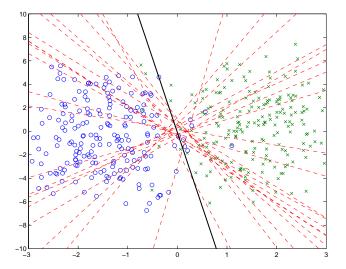
Consensus classification

- ▶ data (examples) (a_i, b_i) , i = 1, ..., N, $a_i \in \mathbf{R}^n$, $b_i \in \{-1, +1\}$
- ▶ linear classifier $sign(a^Tw + v)$, with weight w, offset v
- margin for *i*th example is $b_i(a_i^T w + v)$; want margin to be positive
- ▶ loss for *i*th example is $l(b_i(a_i^Tw + v))$ - *l* is loss function (hinge, logistic, probit, exponential, ...)
- ► choose w, v to minimize $\frac{1}{N} \sum_{i=1}^{N} l(b_i(a_i^T w + v)) + r(w)$ - r(w) is regularization term $(\ell_2, \ell_1, ...)$
- split data and use ADMM consensus to solve

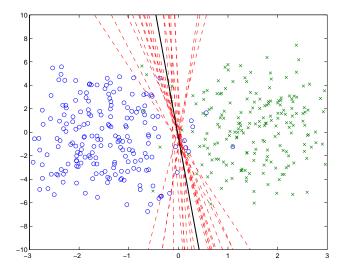
Consensus SVM example

- ▶ hinge loss $l(u) = (1 u)_+$ with ℓ_2 regularization
- ▶ baby problem with n = 2, N = 400 to illustrate
- examples split into 20 groups, in worst possible way: each group contains only positive or negative examples

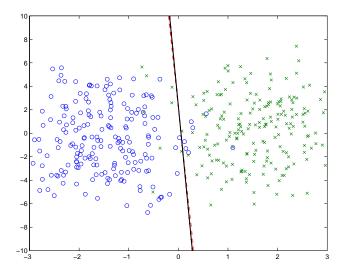
Iteration 1



Iteration 5



Iteration 40



Distributed lasso example

- ▶ example with **dense** $A \in \mathbf{R}^{400000 \times 8000}$ (roughly 30 GB of data)
 - distributed solver written in C using MPI and GSL
 - no optimization or tuned libraries (like ATLAS, MKL)
 - split into 80 subsystems across 10 (8-core) machines on Amazon EC2
- computation times

loading data	30s
factorization	5m
subsequent ADMM iterations	0.5–2s
lasso solve (about 15 ADMM iterations)	5–6m

Exchange problem

minimize
$$\sum_{i=1}^{N} f_i(x_i)$$

subject to $\sum_{i=1}^{N} x_i = 0$

- another canonical problem, like consensus
- ▶ in fact, it's the dual of consensus
- \blacktriangleright can interpret as N agents exchanging n goods to minimize a total cost
- ▶ $(x_i)_j \ge 0$ means agent *i* receives $(x_i)_j$ of good *j* from exchange
- ▶ $(x_i)_j < 0$ means agent *i* contributes $|(x_i)_j|$ of good *j* to exchange
- constraint $\sum_{i=1}^{N} x_i = 0$ is equilibrium or market clearing constraint
- \blacktriangleright optimal dual variable y^{\star} is a set of valid prices for the goods
- suggests real or virtual cash payment $(y^*)^T x_i$ by agent i

Exchange ADMM

▶ solve as a generic constrained convex problem with constraint set

$$\mathcal{C} = \{ x \in \mathbf{R}^{nN} \mid x_1 + x_2 + \dots + x_N = 0 \}$$

► scaled form:

$$\begin{aligned} x_i^{k+1} &:= \arg\min_{x_i} \left(f_i(x_i) + (\rho/2) \| x_i - x_i^k + \overline{x}^k + u^k \|_2^2 \right) \\ u^{k+1} &:= u^k + \overline{x}^{k+1} \end{aligned}$$

► unscaled form:

$$\begin{aligned} x_i^{k+1} &:= \arg\min_{x_i} \left(f_i(x_i) + y^{kT} x_i + (\rho/2) \| x_i - (x_i^k - \overline{x}^k) \|_2^2 \right) \\ y^{k+1} &:= y^k + \rho \overline{x}^{k+1} \end{aligned}$$

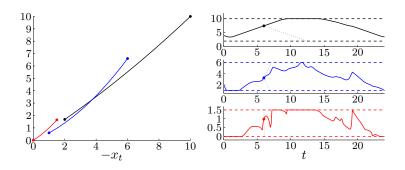
Interpretation as tâtonnement process

- ► tâtonnement process: iteratively update prices to clear market
- work towards equilibrium by increasing/decreasing prices of goods based on excess demand/supply
- dual decomposition is the simplest tâtonnement algorithm
- ADMM adds proximal regularization
 - incorporate agents' prior commitment to help clear market
 - convergence far more robust convergence than dual decomposition

Distributed dynamic energy management

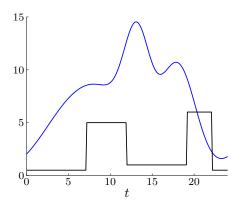
- $\blacktriangleright~N$ devices exchange power in time periods $t=1,\ldots,T$
- $x_i \in \mathbf{R}^T$ is power flow *profile* for device i
- $f_i(x_i)$ is cost of profile x_i (and encodes constraints)
- $x_1 + \cdots + x_N = 0$ is energy balance (in each time period)
- dynamic energy management problem is exchange problem
- exchange ADMM gives distributed method for dynamic energy management
- each device optimizes its own profile, with quadratic regularization for coordination
- residual (energy imbalance) is driven to zero

Generators



- ► 3 example generators
- ▶ left: generator costs/limits; right: ramp constraints
- can add cost for power changes

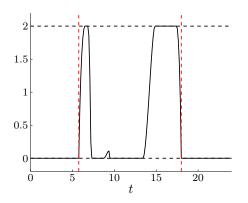
Fixed loads



► 2 example fixed loads

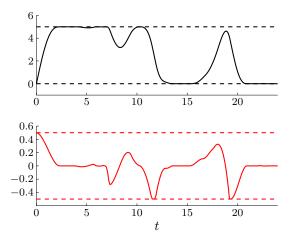
 \blacktriangleright cost is $+\infty$ for not supplying load; zero otherwise

Shiftable load



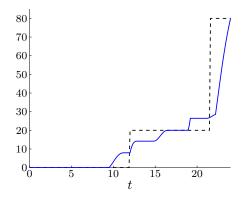
- ► total energy consumed over an interval must exceed given minimum level
- limits on energy consumed in each period
- \blacktriangleright cost is $+\infty$ for violating constraints; zero otherwise

Battery energy storage system



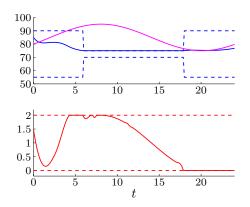
- energy store with maximum capacity, charge/discharge limits
- ► black: battery charge, red: charge/discharge profile
- \blacktriangleright cost is $+\infty$ for violating constraints; zero otherwise

Electric vehicle charging system



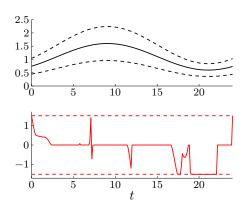
- ► black: desired charge profile; blue: charge profile
- shortfall cost for not meeting desired charge

HVAC



- ▶ thermal load (e.g., room, refrigerator) with temperature limits
- ▶ magenta: ambient temperature; blue: load temperature
- ► red: cooling energy profile
- \blacktriangleright cost is $+\infty$ for violating constraints; zero otherwise

External tie



▶ buy/sell energy from/to external grid at price $p^{\text{ext}}(t) \pm \gamma(t)$

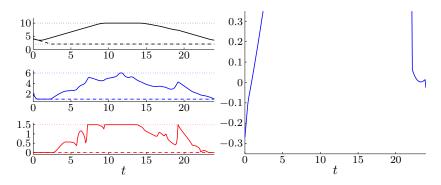
▶ solid: $p^{\text{ext}}(t)$; dashed: $p^{\text{ext}}(t) \pm \gamma(t)$

Smart grid example

10 devices

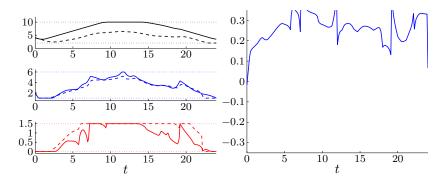
- ► 3 generators
- ► 2 fixed loads
- ▶ 1 shiftable load
- ▶ 1 EV charging systems
- ▶ 1 battery
- ► 1 HVAC system
- ▶ 1 external tie





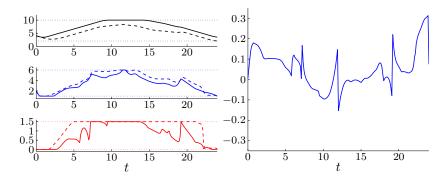
- \blacktriangleright left: solid: optimal generator profile, dashed: profile at kth iteration
- ▶ right: residual vector \bar{x}^k

iteration: k = 3



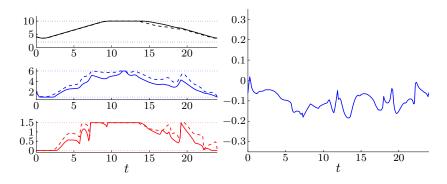
- \blacktriangleright left: solid: optimal generator profile, dashed: profile at kth iteration
- ▶ right: residual vector \bar{x}^k

iteration: k = 5



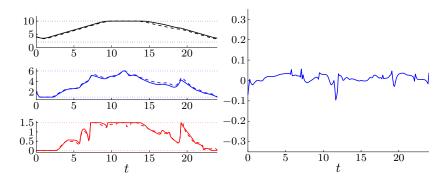
- \blacktriangleright left: solid: optimal generator profile, dashed: profile at kth iteration
- \blacktriangleright right: residual vector \bar{x}^k

iteration: k = 10



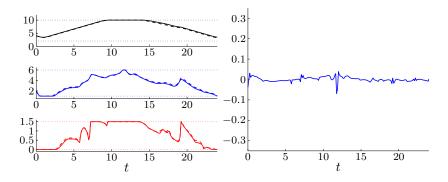
- ▶ left: solid: optimal generator profile, dashed: profile at *k*th iteration
- \blacktriangleright right: residual vector \bar{x}^k

iteration: k = 15



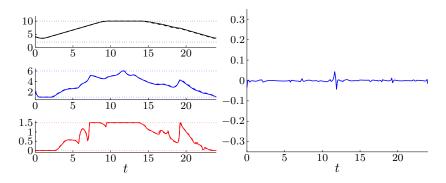
- ▶ left: solid: optimal generator profile, dashed: profile at *k*th iteration
- \blacktriangleright right: residual vector \bar{x}^k

iteration: k = 20



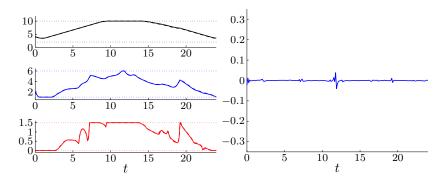
- \blacktriangleright left: solid: optimal generator profile, dashed: profile at kth iteration
- \blacktriangleright right: residual vector \bar{x}^k

iteration: k = 25



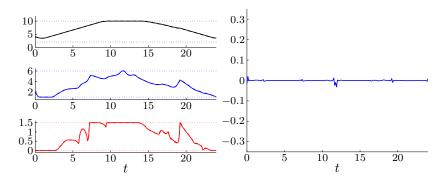
- ▶ left: solid: optimal generator profile, dashed: profile at *k*th iteration
- \blacktriangleright right: residual vector \bar{x}^k

iteration: k = 30



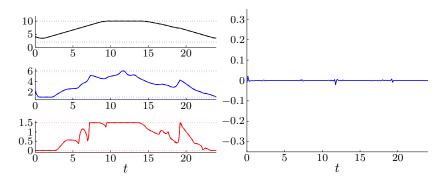
- \blacktriangleright left: solid: optimal generator profile, dashed: profile at kth iteration
- \blacktriangleright right: residual vector \bar{x}^k

iteration: k = 35



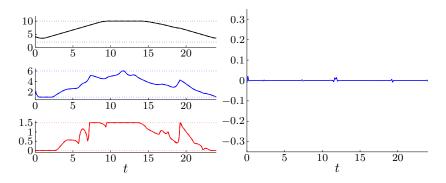
- \blacktriangleright left: solid: optimal generator profile, dashed: profile at kth iteration
- \blacktriangleright right: residual vector \bar{x}^k

iteration: k = 40



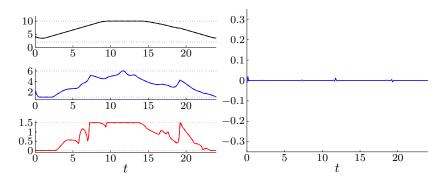
- ▶ left: solid: optimal generator profile, dashed: profile at *k*th iteration
- \blacktriangleright right: residual vector \bar{x}^k

iteration: k = 45



- \blacktriangleright left: solid: optimal generator profile, dashed: profile at kth iteration
- \blacktriangleright right: residual vector \bar{x}^k

iteration: k = 50



- \blacktriangleright left: solid: optimal generator profile, dashed: profile at kth iteration
- \blacktriangleright right: residual vector \bar{x}^k

Outline

Dual decomposition

Method of multipliers

Alternating direction method of multipliers

Common patterns

Examples

Consensus and exchange

Conclusions

Conclusions

Summary and conclusions

ADMM

- ▶ is the same as, or closely related to, many methods with other names
- has been around since the 1970s
- gives simple single-processor algorithms that can be competitive with state-of-the-art
- can be used to coordinate many processors, each solving a substantial problem, to solve a very large problem