

# ECE 8201: Low-dimensional Signal Models for High-dimensional Data Analysis

Lecture 9: Atomic norm for low-complexity signal models

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## Main Reference

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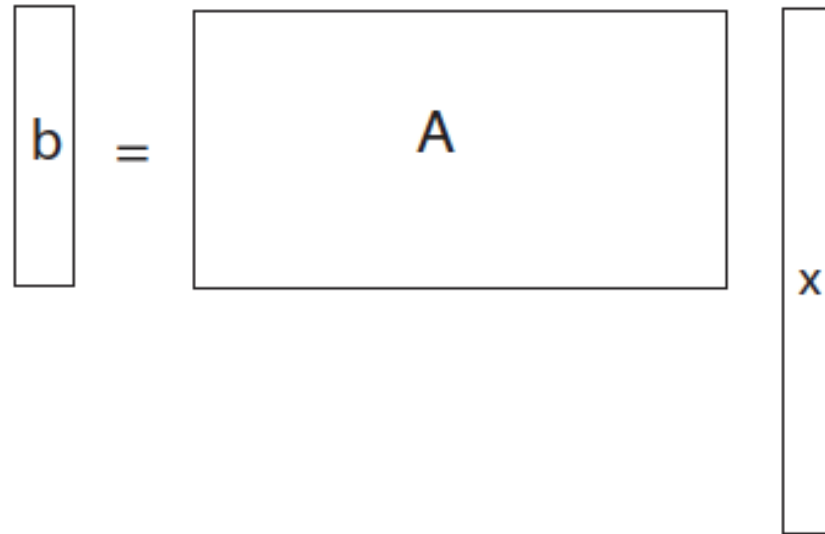
- Chandrasekaran, V., B. Recht, P. A. Parrilo, and A. S. Willsky. "The convex geometry of linear inverse problems." *Foundations of Computational Mathematics* 12, no. 6 (2012): 805-849.

# Recap

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- We start with an under-determined linear system:

$$b = Ax$$



- Estimate  $x \in \mathbb{R}^n$  from linear measurements  $b = Ax \in \mathbb{R}^m$ , where  $m \ll n$ .

# Recap

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Hope?

- Constrain the set of interesting signals: looking for  $x$  with additional structures:
  - sparsity:  $\|x\|_0$  is small;
  - low-rankness:  $\text{rank}(\text{reshape}(x))$  is small;
- Constrain the measurement operator:
  - The measurement matrix  $A$  has to be “incoherent” to the postulated structure: for example,  $A$  cannot be a partial identity matrix for sparse  $x$ ;
  - More generically, this incoherence can be provided if we choose “random” measurement matrix, e.g.  $A$  composed of iid Gaussian entries provides “universal” guarantees.

- Efficient algorithms based on convex relaxations:

$$\min \|\mathbf{x}\|_1, \quad \text{s.t.} \quad \mathbf{b} = \mathbf{A}\mathbf{x},$$

and nuclear norm minimization for rank minimization;

- advantages:
  - computationally efficient, many solvers have been developed to handle large-scale convex problems;
  - provable near-optimal performance in terms of sample complexity through the machinery of convex analysis;
- Question: can we extend this framework to other low-dimensional structures?

# The Atomic Norm Approach

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- The **atomic norm** is proposed by Chandrasekaran et.al. to find tightest convex relaxations of general parsimonious models.
- “Parsimony” refers to the fact that the signal of interest  $x$  can be described by a much smaller number of parameters than its ambient dimension;
  - For a known room, the image of the room is determined by the location and orientation of the camera, rather than the number of pixels of the image;
  - For an  $K$ -sparse vector  $x$ , it can be described by  $2K$  parameters;
- This models the signal of interest  $x$  as composed of atoms in an *atomic* set:

$$\mathcal{A} = \{\mathbf{a}_i\}$$

which could be infinite;

- The signal  $x$  can be written as a superposition of a small number of *atoms* in an atomic set  $\mathcal{A}$ :

$$\mathbf{x} = \sum_{i=1}^r c_i \mathbf{a}_i, \quad \mathbf{a}_i \in \mathcal{A}, c_i > 0.$$

Known examples:

- Sparse case:  $\mathcal{A}$  is composed of normalized vectors of sparsity one;
- Low-rank case:  $\mathcal{A}$  is composed of normalized matrices of rank one;
- Define the atomic norm (a.k.a. the “guage of  $\mathcal{A}$ ”) as

$$\begin{aligned} \|x\|_{\mathcal{A}} &= \inf \{t > 0 : x \in t\text{conv}(\mathcal{A})\} \\ &= \inf \left\{ \sum_i c_i \mid x = \sum_i c_i a_i, a_i \in \mathcal{A}, c_i > 0 \right\}. \end{aligned}$$

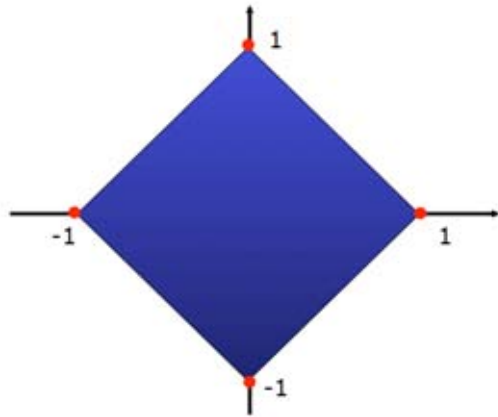
if  $\mathcal{A}$  is centrally symmetric about the origin (i.e.,  $\mathbf{a} \in \mathcal{A}$  if and only if  $-\mathbf{a} \in \mathcal{A}$ ) we have that  $\|\cdot\|_{\mathcal{A}}$  is a norm. It is also a convex function.

## Special cases of atomic norm

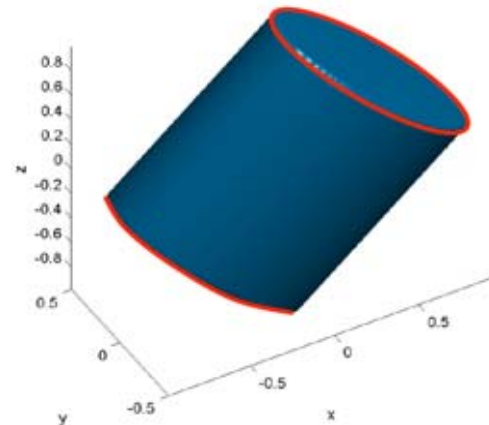
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The atomic norm minimization recovers the  $\ell_1$  and nuclear norm:

- **Sparse signals:** *an atom for sparse signals* is a normalized vector of sparsity one, and the atomic norm is  $\ell_1$  norm;
- **Low-rank matrices:** *an atom for low-rank matrices* is a normalized rank-one matrix; and the atomic norm is nuclear norm;



unit ball of  $\ell_1$  norm



unit ball of nuclear norm



# Atomic norm minimization

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... for underdetermined linear systems:

- For a given atomic set  $\mathcal{A}$ , define the atomic norm  $\|\mathbf{x}\|_{\mathcal{A}}$ ;
- Run the convex program:

$$\min_{\mathbf{x}} \|\mathbf{x}\|_{\mathcal{A}} \quad \text{s.t.} \quad \mathbf{b} = \mathbf{A}\mathbf{x}$$

- For noisy measurements:

$$\min_{\mathbf{x}} \|\mathbf{x}\|_{\mathcal{A}} \quad \text{s.t.} \quad \|\mathbf{b} - \mathbf{A}\mathbf{x}\|_2 \leq \epsilon$$

## Examples of atomic norm

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- “democratic” signal representations via  $l_\infty$  norm;
- “joint sparsity” or “group sparsity” via  $l_1/l_2$  norm;

# Democratic representations

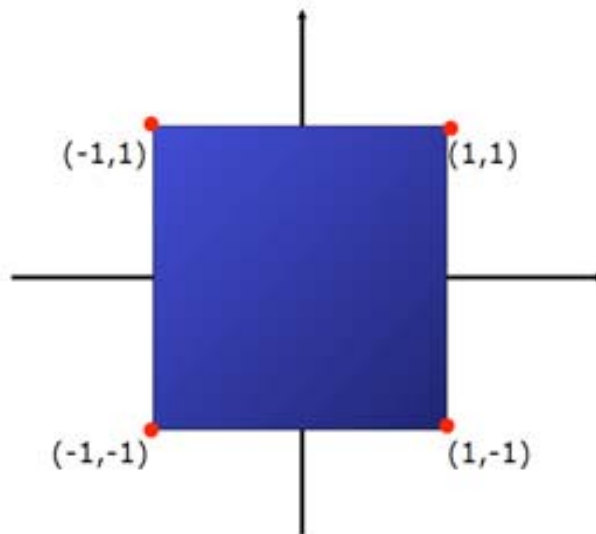
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- Signal representations  $\mathbf{x}$  that have the same amplitudes for every entry;



- motivated by integer programming  $\mathbf{x} \in \{+1, -1\}^n$ ;
- peak-to-average power ratio reduction in OFDM communication;

- The atomic set contains all sign vectors  $\mathcal{A} = \{\{+1, -1\}^n\}$ ;
- The atomic norm becomes  $\|\mathbf{x}\|_{\mathcal{A}} = \|\mathbf{x}\|_{\infty}$



# Joint sparsity

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Consider sparse recovery with multiple snapshots:

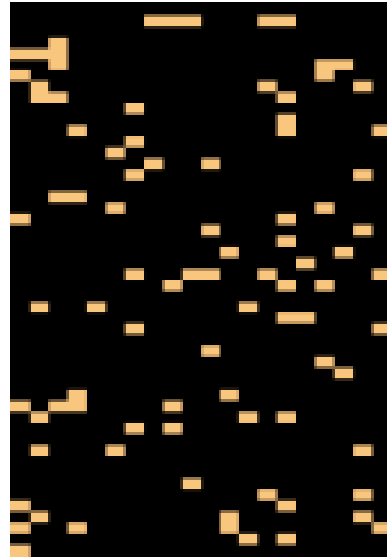
$$B = AX$$

where  $X \in \mathbb{R}^{n \times T}$ , where  $T$  is the number of snapshots.



- Different snapshots  $X = [\mathbf{x}_1, \dots, \mathbf{x}_T]$  share the same support but have different coefficients;

- motivated by multi-task learning, array processing ,etc...



(a) Sparse



(b) Group sparse

- The atoms in  $\mathcal{A}$  can be written as rank-one matrix as

$$\mathcal{A} = \{\mathbf{e}_i \mathbf{u}_i^T \mid \|\mathbf{u}_i\|_2 = 1\}$$

where  $\mathbf{e}_i$  is the  $i$ th standard basis vector, and  $\|\mathbf{u}_i\|_2 = 1$ .

- The atomic norm becomes

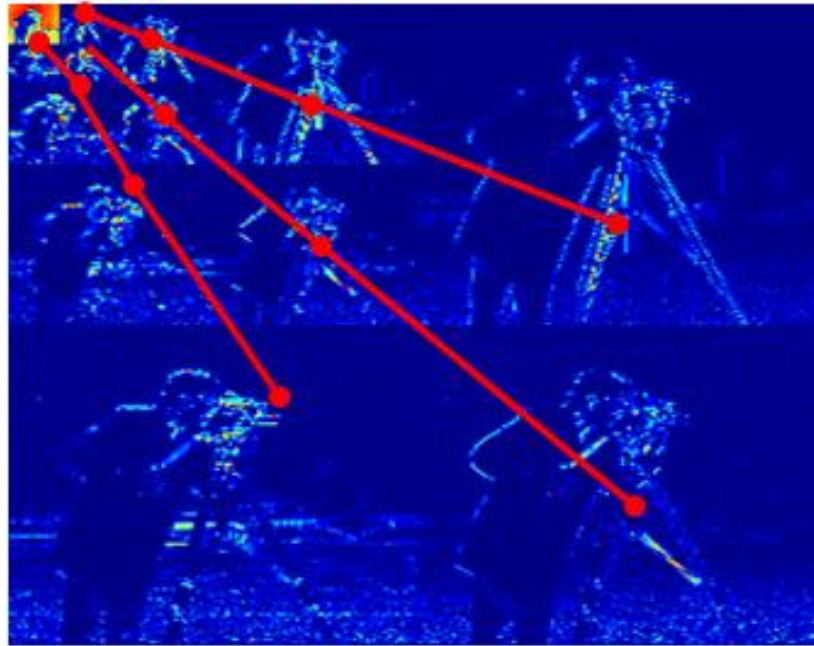
$$\begin{aligned} \|\mathbf{X}\|_{\mathcal{A}} &= \inf \{t > 0 : \mathbf{X} \in t\text{conv}(\mathcal{A})\} \\ &= \inf \left\{ \sum_i c_i \mid \mathbf{X} = \sum_i c_i \mathbf{e}_i \mathbf{u}_i^T, \|\mathbf{u}_i\|_2 = 1, c_i > 0 \right\} \\ &= \sum_{i=1}^n \left( \sum_{j=1}^T |x_{ij}|^2 \right)^{1/2} := \|\mathbf{X}\|_{1,2} \end{aligned}$$

which is the  $\ell_1/\ell_2$  norm of  $\mathbf{X}$ .

## Group sparsity

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The joint sparsity for multiple snapshots problem is related to the so-called group sparsity/structured sparsity.



We can group the coefficients into overlapping/non-overlapping groups, such that we motivate sparsity between groups, but not within the groups.



## Performance with Gaussian measurements

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If  $\mathbf{A}$  is composed of i.i.d. Gaussian entries, how many measurements do we need to guarantee success recovery?

$$\hat{\mathbf{x}} = \underset{\mathbf{x}}{\operatorname{argmin}} \|\mathbf{x}\|_{\mathcal{A}} \quad \text{s.t.} \quad \mathbf{b} = \mathbf{A}\mathbf{x}$$

Surprisingly, this can be answered with a single notion called “Gaussian width”

- A convex set is a cone if it is closed under positive linear combinations.
- The tangent cone at  $\mathbf{x}$  with respect to the scaled unit ball  $\|\mathbf{x}\|_{\mathcal{A}}\text{conv}(\mathcal{A})$  is

$$T_{\mathcal{A}}(\mathbf{x}) = \text{cone}\{\mathbf{z} - \mathbf{x} : \|\mathbf{z}\|_{\mathcal{A}} \leq \|\mathbf{x}\|_{\mathcal{A}}\}$$

which is the set of descent directions of the atomic norm at the point  $\mathbf{x}$ ,

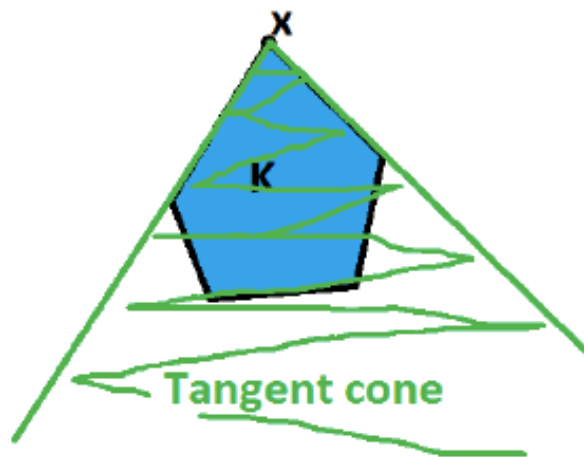


FIGURE 1. The tangent cone

- **Proposition:** We have that  $\hat{x} = x^*$  is the unique optimal solution if and only if  $\text{null}(A) \cap T_{\mathcal{A}}(x^*) = \{0\}$ .

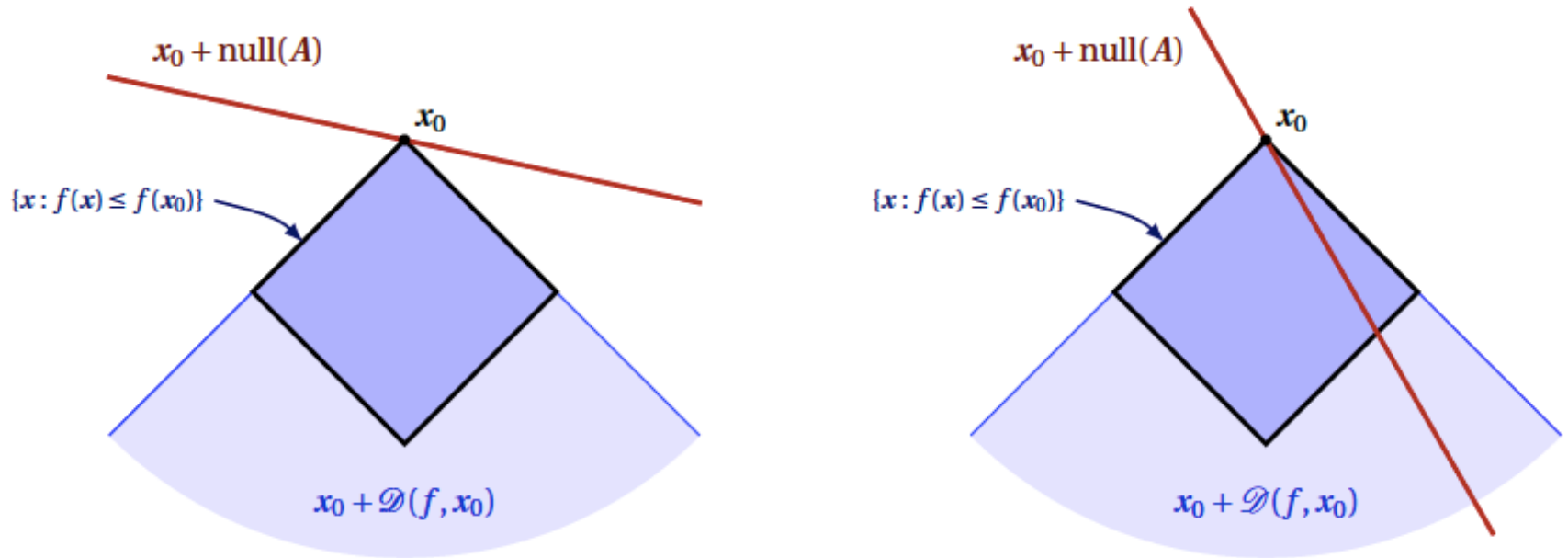


FIGURE 2.3: The optimality condition for a regularized inverse problem. The condition for the regularized linear inverse problem (2.4) to succeed requires that the descent cone  $\mathcal{D}(f, x_0)$  and the null space  $\text{null}(A)$  do not share a ray. [left] The regularized linear inverse problem succeeds. [right] The regularized linear inverse problem fails.

# Gaussian width

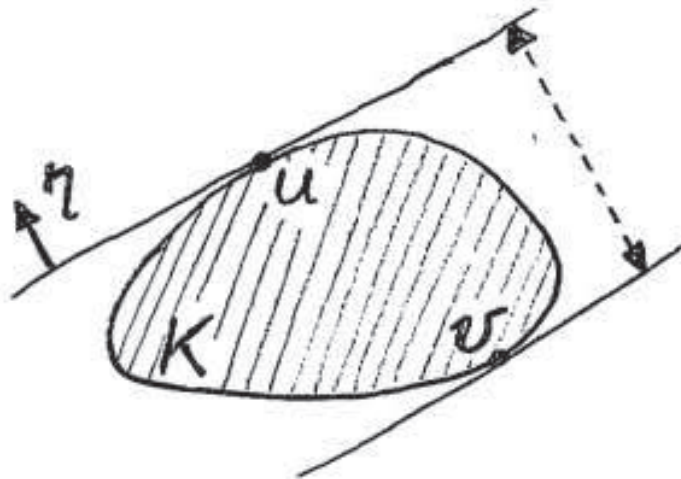
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- Definition: The Gaussian width of a set  $S \subset \mathbb{R}^n$  is defined as:

$$w(S) := \mathbb{E}_g \left[ \sup_{z \in S} g^T z \right], \quad g \sim \mathcal{N}(0, I)$$

- Related to the “mean width”:

$$w(S) = \frac{\lambda_p}{2} \int_{\mathbb{S}^{p-1}} \left( \max_{z \in S} u^T z - \min_{z \in S} u^T z \right) du = \frac{\lambda_p}{2} b(S)$$



- Gordon's escape through the mesh theorem: Let  $\Omega$  be a closed subset of  $\mathbb{S}^{n-1}$ . Let  $\mathbf{A} \in \mathbb{R}^{m \times n}$  be composed of i.i.d. standard Gaussian entries. Then

$$\mathbb{E} \left[ \min_{\mathbf{z} \in \Omega} \|\mathbf{A}\mathbf{z}\|_2 \right] \geq \lambda_m - w(\Omega)$$

where  $\lambda_m = \mathbb{E}[\|\mathbf{g}\|_2] \leq \sqrt{m}$ .

- Immediately, we have the number of measurements we need is  $m \gtrsim w(\Omega)^2 + 1$ :  
to guarantee

$$\text{null}(\mathbf{A}) \cap T_{\mathcal{A}}(\mathbf{x}^*) = \{0\}$$

by setting  $\Omega = T_{\mathcal{A}}(\mathbf{x}^*) \cap \mathbb{S}^{n-1}$ .

# Measurement bound with Gaussian width

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**Corollary 3.3.** *Let  $\Phi : \mathbb{R}^p \rightarrow \mathbb{R}^n$  be a random map with i.i.d. zero-mean Gaussian entries having variance  $1/n$ . Further let  $\Omega = T_{\mathcal{A}}(\mathbf{x}^*) \cap \mathbb{S}^{p-1}$  denote the spherical part of the tangent cone  $T_{\mathcal{A}}(\mathbf{x}^*)$ .*

1. *Suppose that we have measurements  $\mathbf{y} = \Phi \mathbf{x}^*$  and solve the convex program (5). Then  $\mathbf{x}^*$  is the unique optimum of (5) with probability at least  $1 - \exp\left(-\frac{1}{2} [\lambda_n - w(\Omega)]^2\right)$  provided*

$$n \geq w(\Omega)^2 + 1.$$

Underlying model	Convex heuristic	# Gaussian measurements
$s$ -sparse vector in $\mathbb{R}^p$	$\ell_1$ norm	$2s \log(p/s) + 5s/4$
$m \times m$ rank- $r$ matrix	nuclear norm	$3r(2m - r)$
sign-vector $\{-1, +1\}^p$	$\ell_\infty$ norm	$p/2$
$m \times m$ permutation matrix	norm induced by Birkhoff polytope	$9m \log(m)$
$m \times m$ orthogonal matrix	spectral norm	$(3m^2 - m)/4$

Table 1: A summary of the recovery bounds obtained using Gaussian width arguments.