

# Fast Energy Minimization with Label Costs and Applications in Model Fitting

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*presented by* Anton Osokin

*co-authors:*

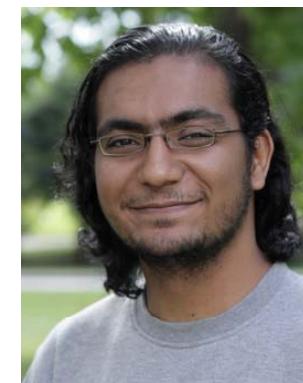
Andrew Delong



Yuri Boykov



Hossam Isack



some slides taken from Yuri Boykov and Andrew Delong

## Popular energy in vision

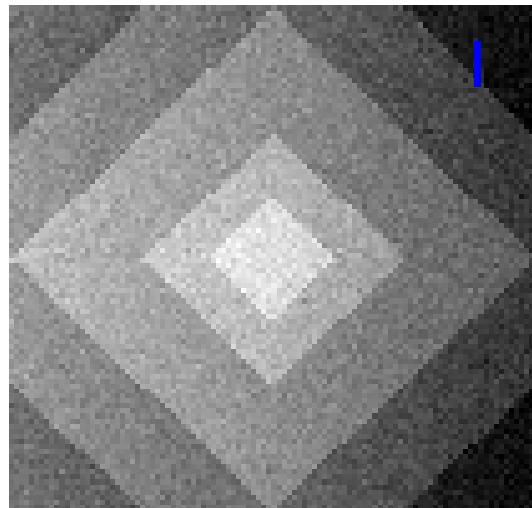
$$f_p \in \{1, \dots, K\}$$

$$E(f) = \sum_p D_p(f_p) + \sum_{(p,q) \in N} V_{pq}(f_p, f_q) \rightarrow \min_f$$

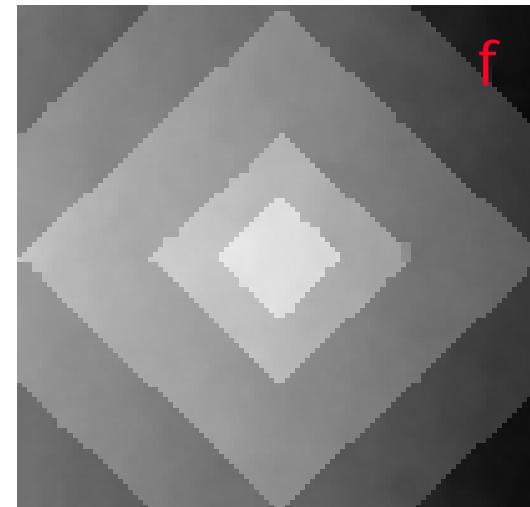
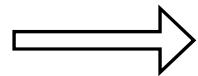
$D_p(f_p)$  – unary potentials

$V_{pq}(f_p, f_q)$  – pairwise potentials

# Image restoration



*observed noisy image  $I$*



*image labeling  $f$   
(restored intensities)*

$$\mathbf{I} = \{ I_1, I_2, \dots, I_n \}$$

$$\mathbf{F} = \{ f_1, f_2, \dots, f_n \}$$

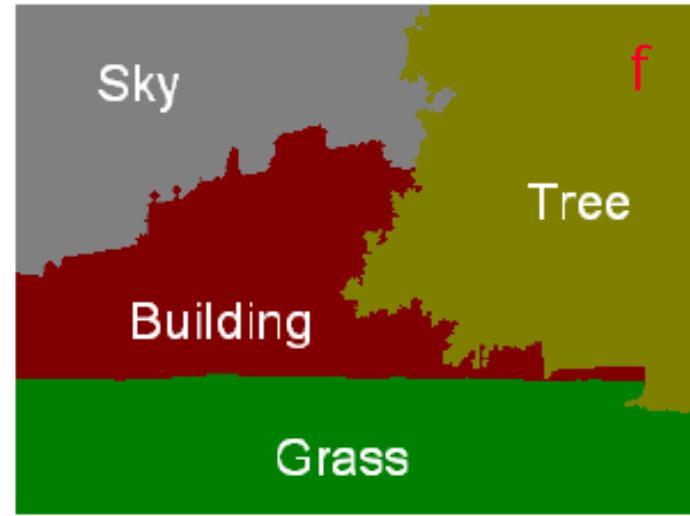
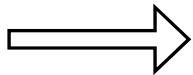
$$E(f) = \sum_p (f_p - I_p)^2 + \sum_{(p,q) \in N} (f_p - f_q)^2$$

*data fidelity*      *spatial regularization*

# Image segmentation



*Observed image  $I$*

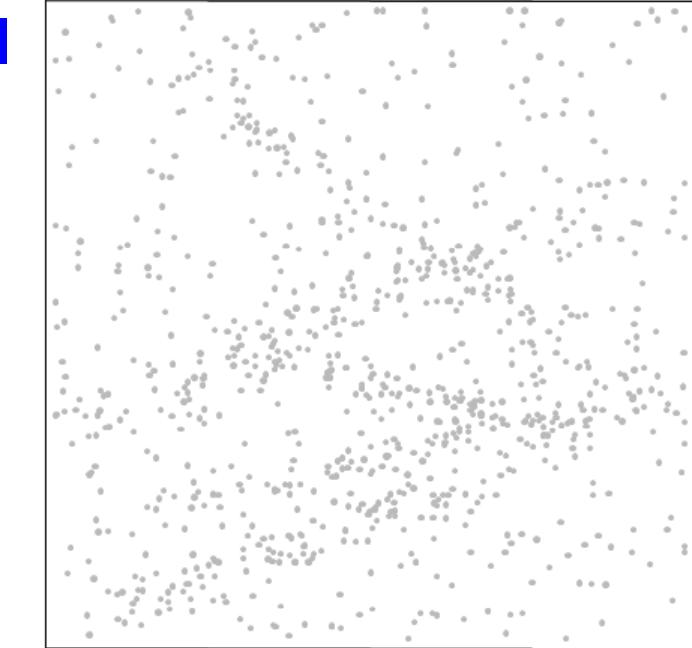


*image labeling  $f$   
(segments)*

$$E(f) = \sum_p -\log P(f_p | I_p) + \theta \sum_{(p,q) \in N} [f_p \neq f_q]$$

*data fidelity*      *spatial regularization*

# Geometric model fitting

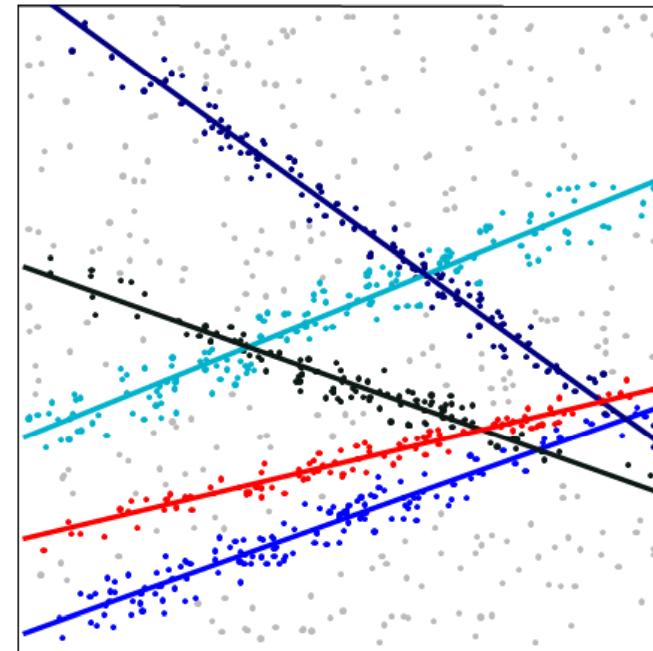


Sampled points  $I$

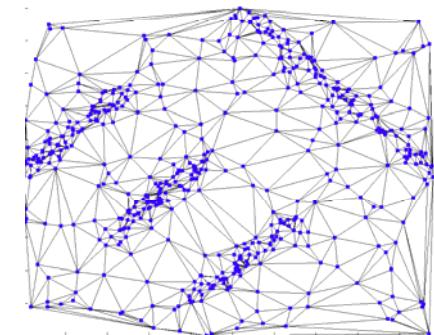
*data fidelity*

$$E(f) = \sum_p -\log dist(p, f_p) + \theta \sum_{(p,q) \in N} \frac{[f_p \neq f_q]}{dist(p, q)}$$

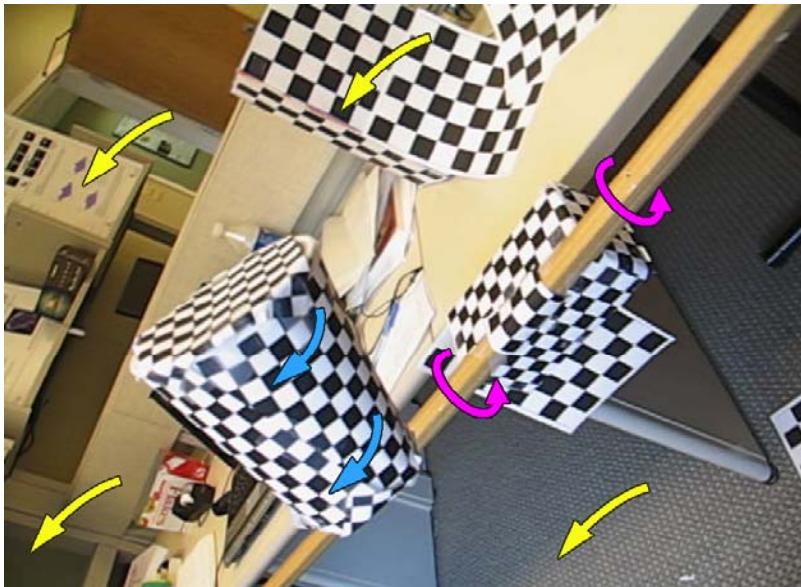
*spatial regularization*



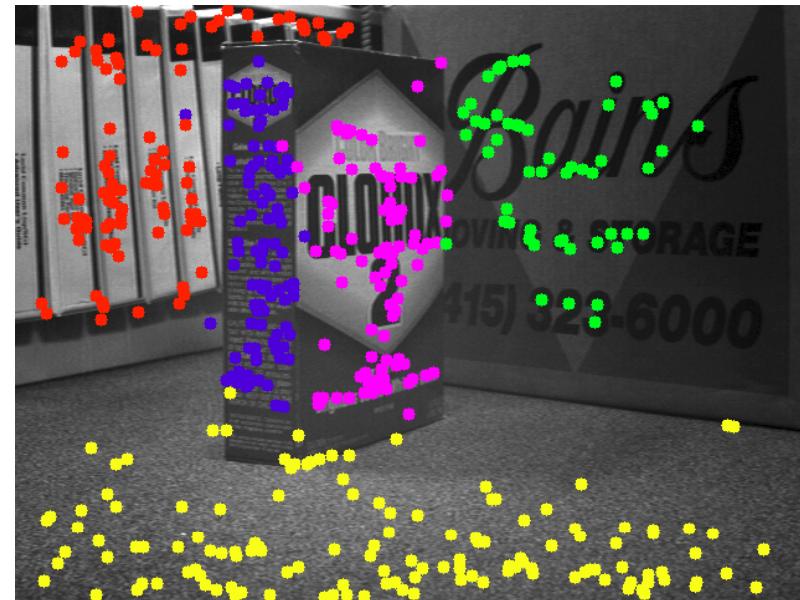
Points clustering  $f$



## More geometric model fitting



Motion estimation



Plane fitting

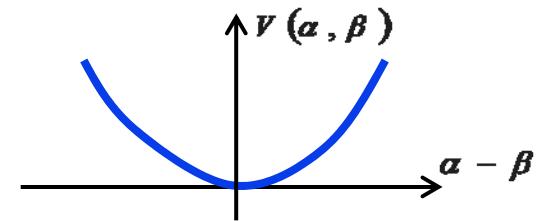
$$E(f) = \sum_p D_p(f_p) + \sum_{(p,q) \in N} V_{pq}(f_p, f_q)$$

# Optimization

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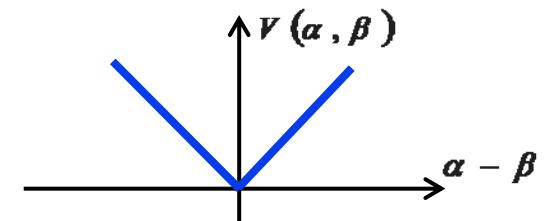
## ■ Convex regularization

- gradient descent works
- exact polynomial algorithms



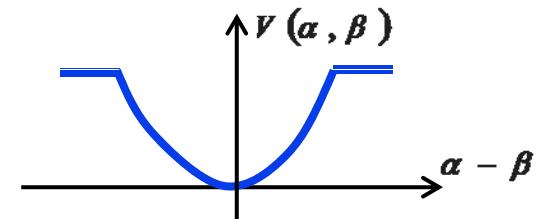
## ■ TV regularization

- a bit harder (non-differentiable)
- global minima algorithms (Ishikawa, etc.)



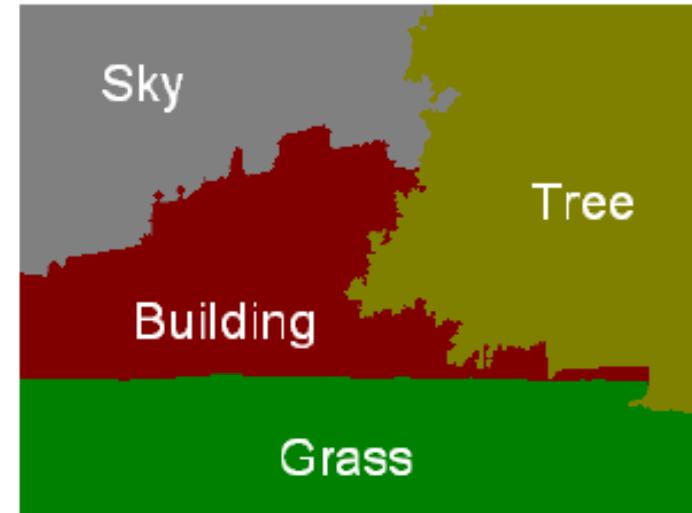
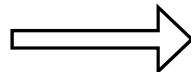
## ■ Robust regularization

- NP-hard, many local minima
- good approximations (message passing, a-expansion, a/b-swap)



Potts model  $E(f) = \sum_p -\log P(f_p | I_p) + \theta \sum_{(p,q) \in N} [f_p \neq f_q]$   
 (piece-wise constant labeling)

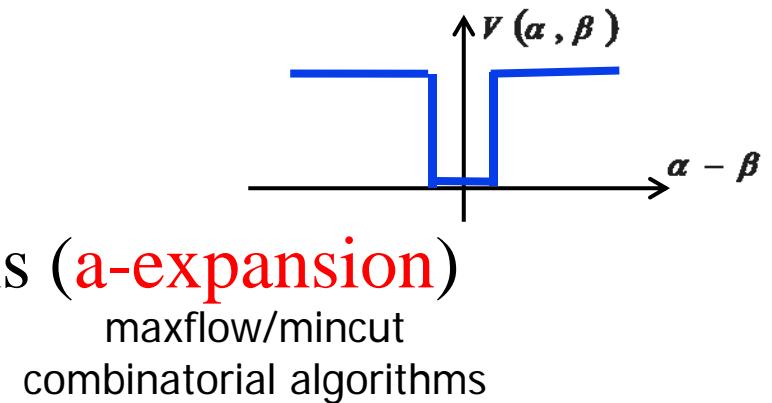
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$$V(\alpha, \beta) = \theta \cdot [\alpha \neq \beta]$$

## ■ Robust regularization

- NP-hard, many local minima
- provably good approximations (**a-expansion**)



## Adding label costs

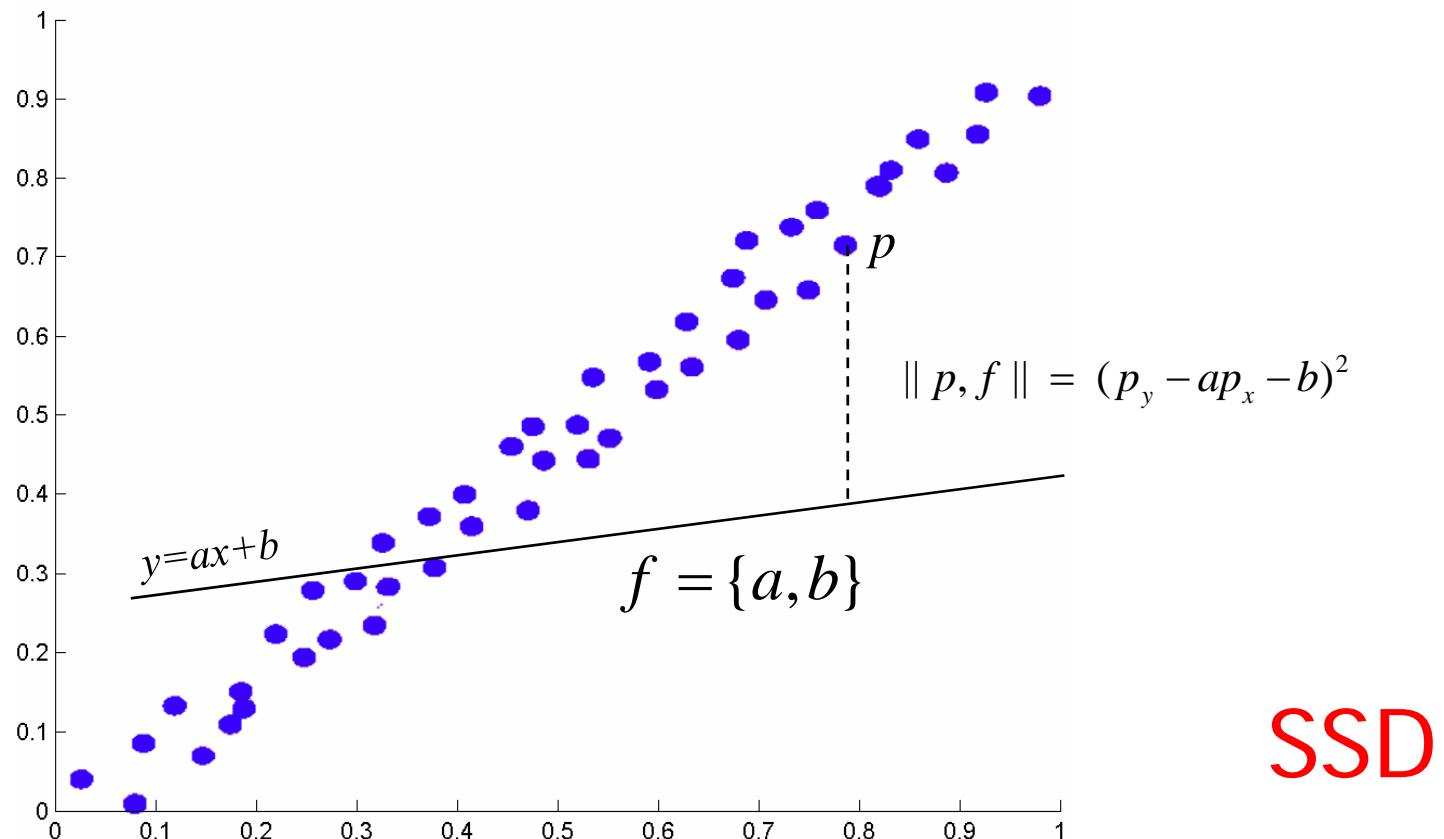
$$E(f) = \sum_p D_p(f_p) + \sum_{(p,q) \in N} V(f_p, f_q) + \sum_{f \in \Lambda} h_f \cdot \delta_f(f)$$

$\Lambda$  - set of labels  
allowed at each point p

$$\delta_f(f) = \begin{cases} 1, & \exists p : f_p = f \\ 0, & otherwise \end{cases}$$

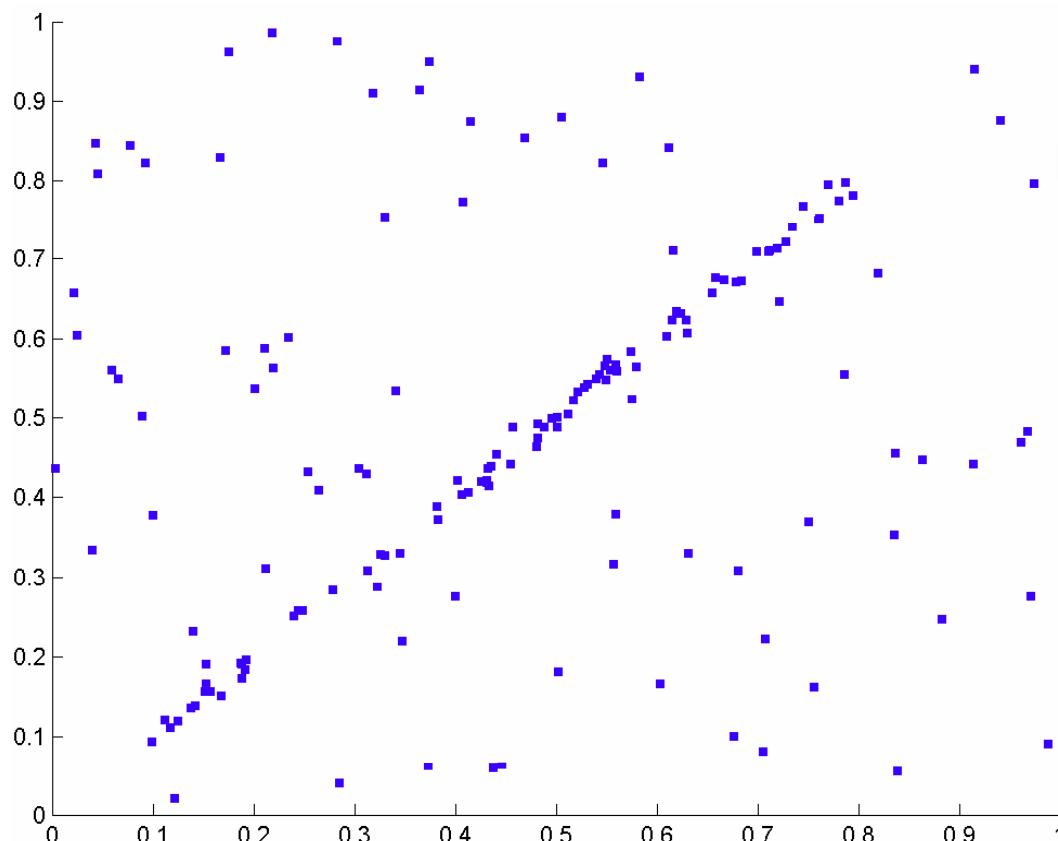
# Model fitting

$$f^* = \arg \min_f \sum_p \| p, f \|$$



# Many outliers

quadratic errors fail

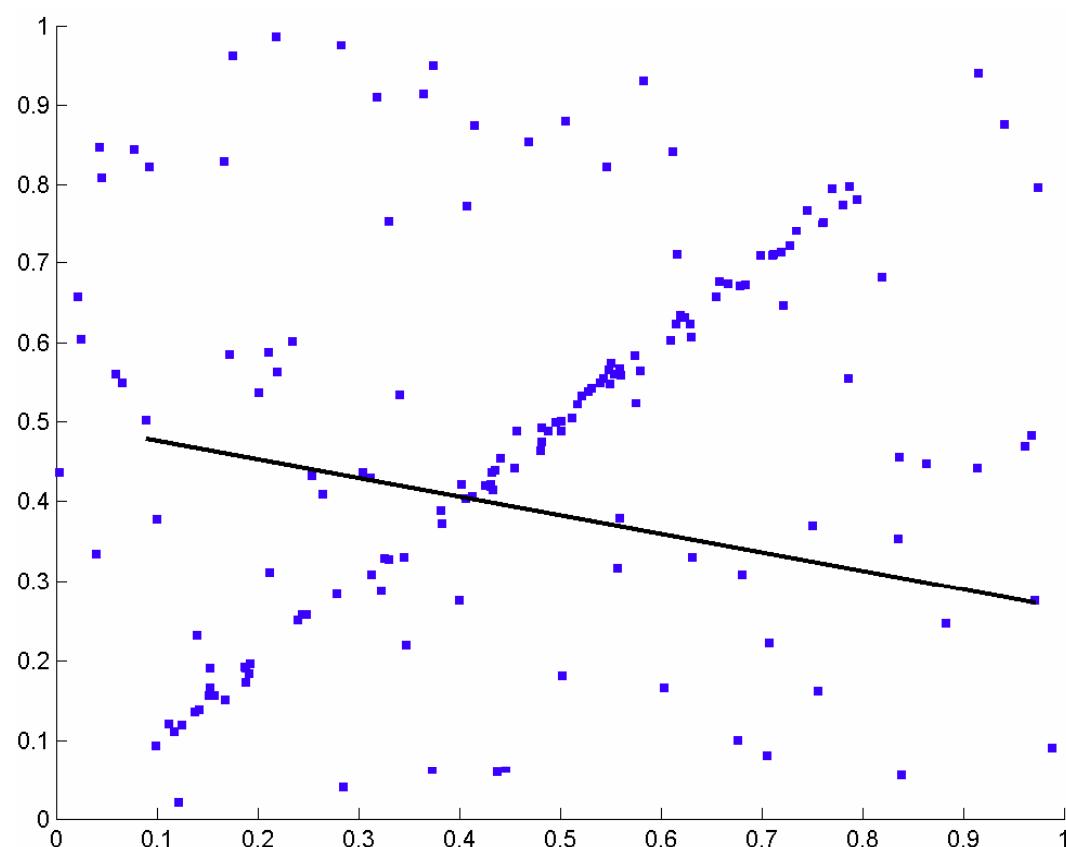


use more robust  
error measures, e.g.  
 $\| p, f \| = | p_y - ap_x - b |$   
gives “MEDIAN” line  

- more expensive  
computations  
(non-differentiable)
- **still fails if  
outliers exceed  
50%**

**RANSAC**

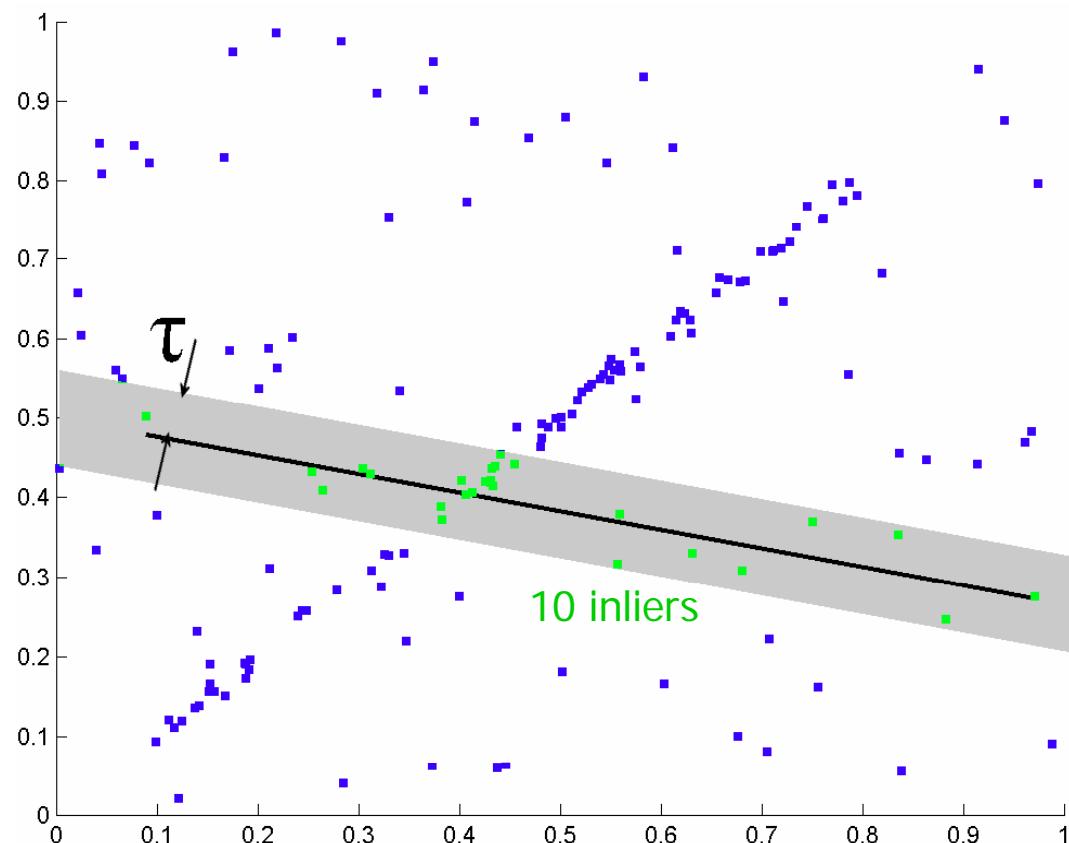
# Many outliers



1. sample randomly two points, get a line

**RANSAC**

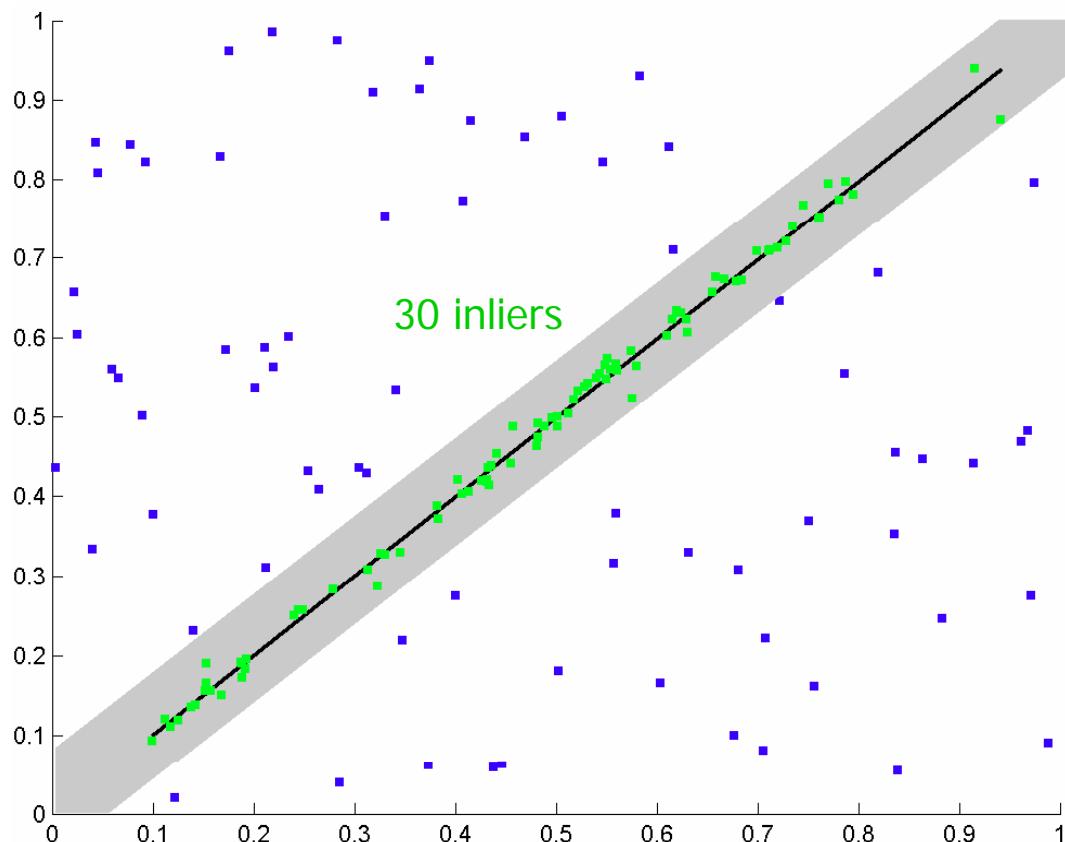
# Many outliers



1. sample randomly two points, get a line
2. count inliers for threshold  $T$

RANSAC

# Many outliers

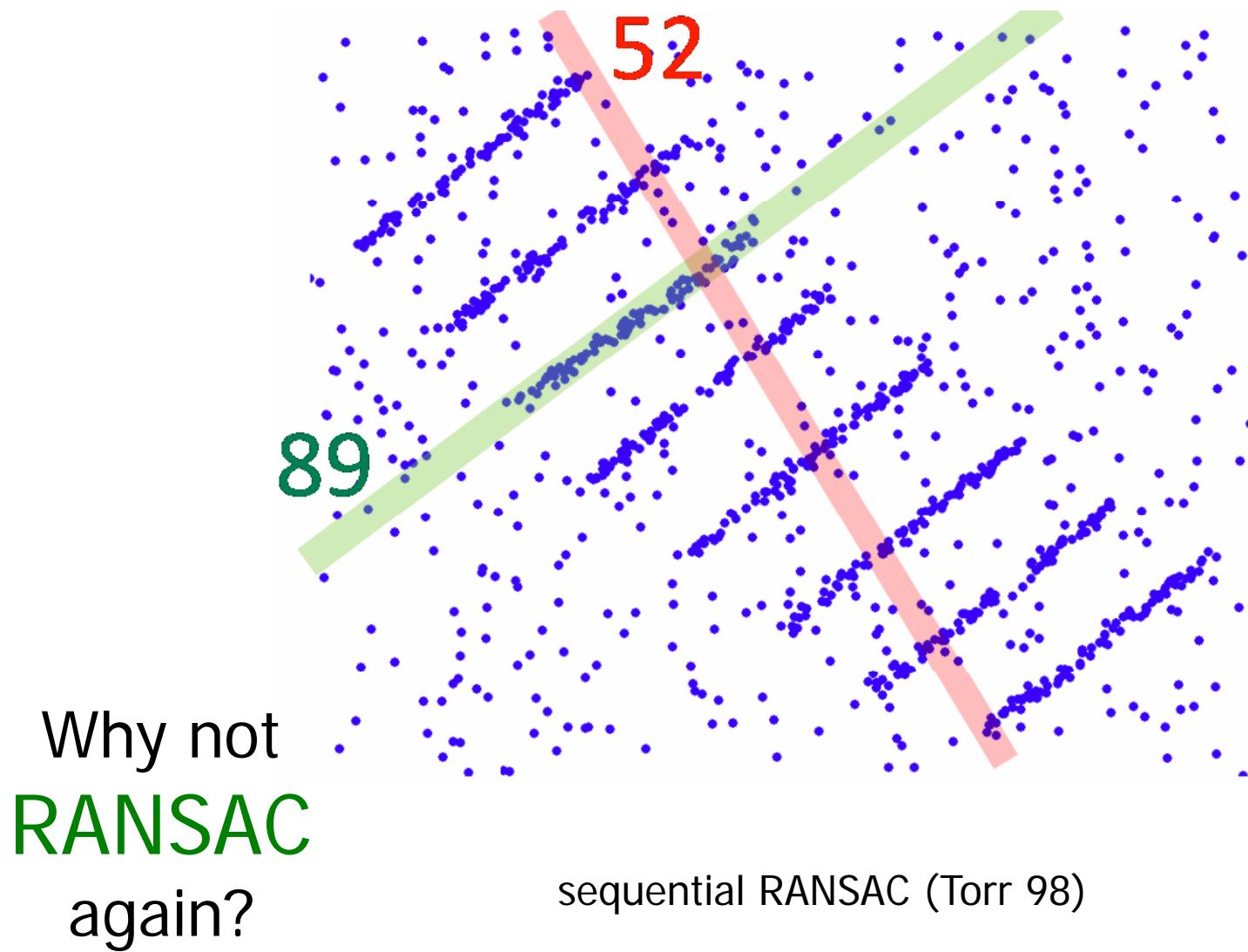


1. sample randomly two points, get a line
2. count inliers for threshold  $T$
3. repeat N times and select model with most inliers

**RANSAC**

# Multiple models and many outliers

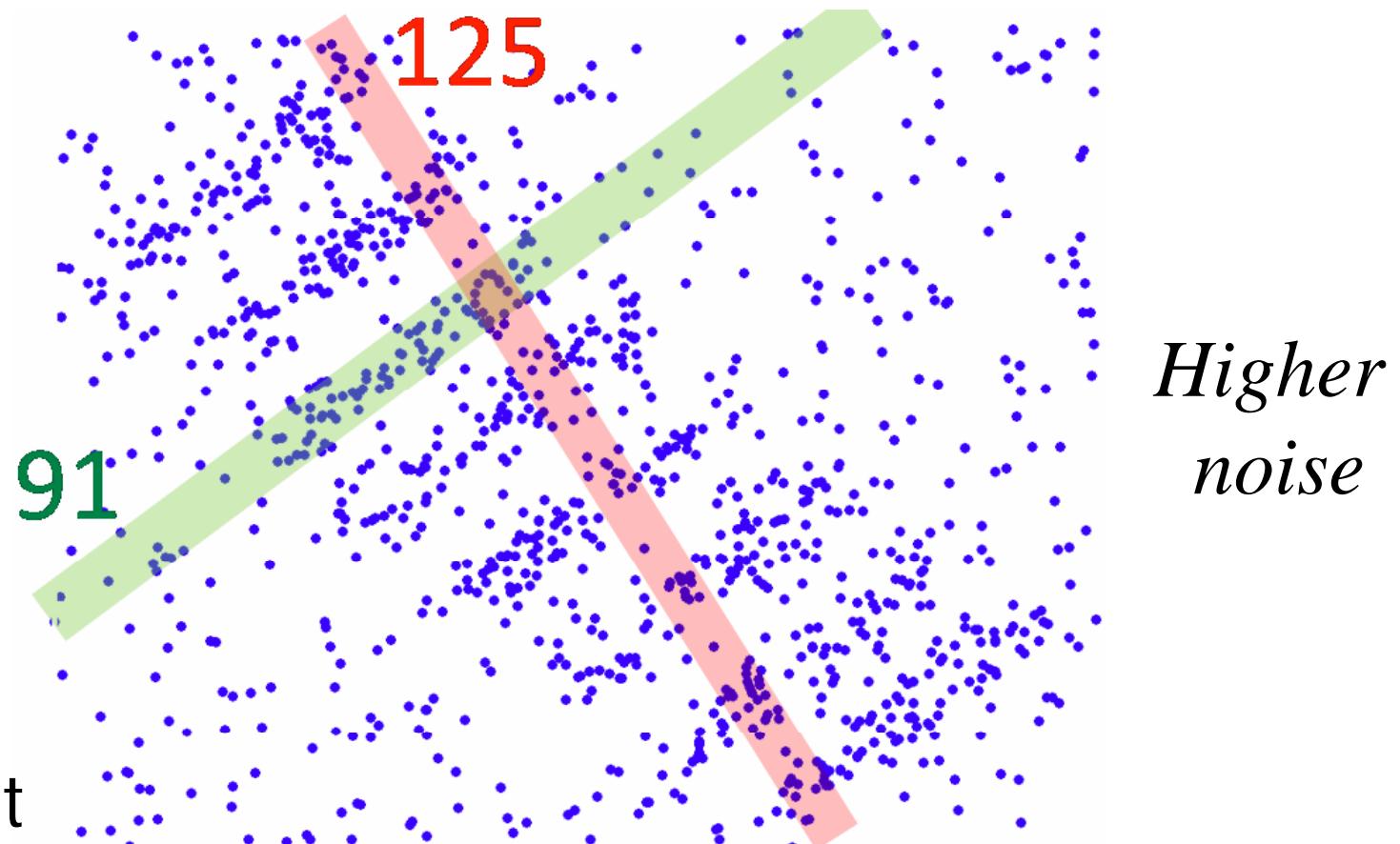
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# Multiple models and many outliers

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Why not  
**RANSAC**  
again?



In general, maximization of inliers  
does not work for  
**outliers + multiple models**

# Energy-based approach

$$E(f) = \sum_p \| p, f \|$$

energy-based interpretation  
of RANSAC criteria for  
**single** model fitting:

- find optimal **label**  $f$   
for one very specific  
error measure

$$\| p, f \| = \begin{cases} 0, & \text{if } dist(p, f) \leq T \\ 1, & \text{if } dist(p, f) > T \end{cases}$$

# Energy-based approach

$$E(f) = \sum_p \| p, f_p \|$$

If **multiple** models

- assign different models (labels  $f_p$ ) to every point p

Need regularization!

-find optimal **labeling**

$$f = \{ f_1, f_2, \dots, f_n \}$$

# Spatial regularization

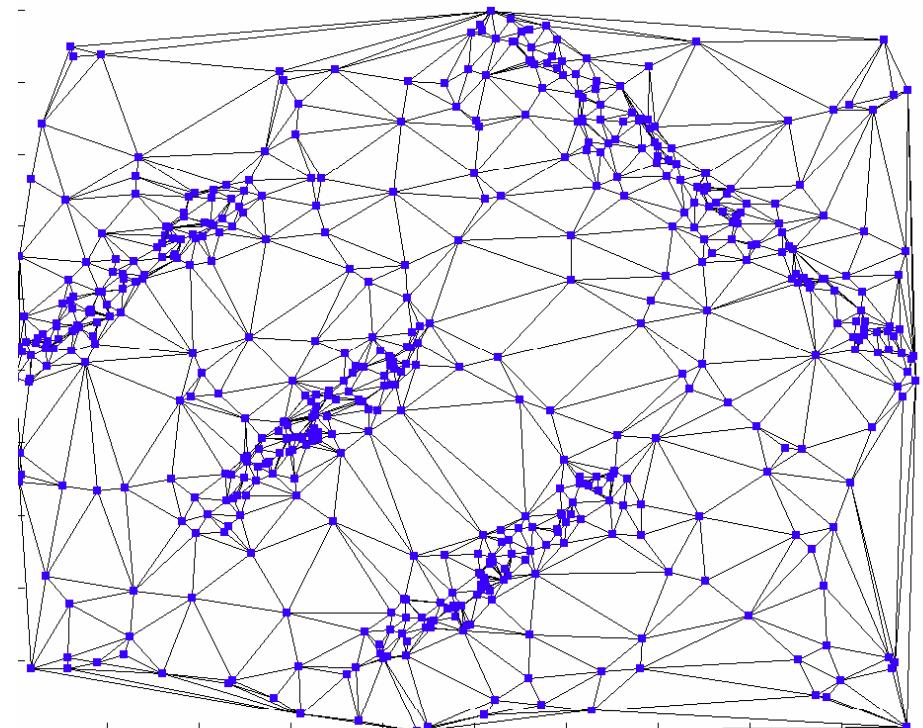
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$$E(f) = \sum_p \| p, f_p \| + \sum_{pq \in N} \theta_{pq} \cdot [f_p \neq f_q]$$

If **multiple** models

- assign different models (labels  $f_p$ ) to every point p

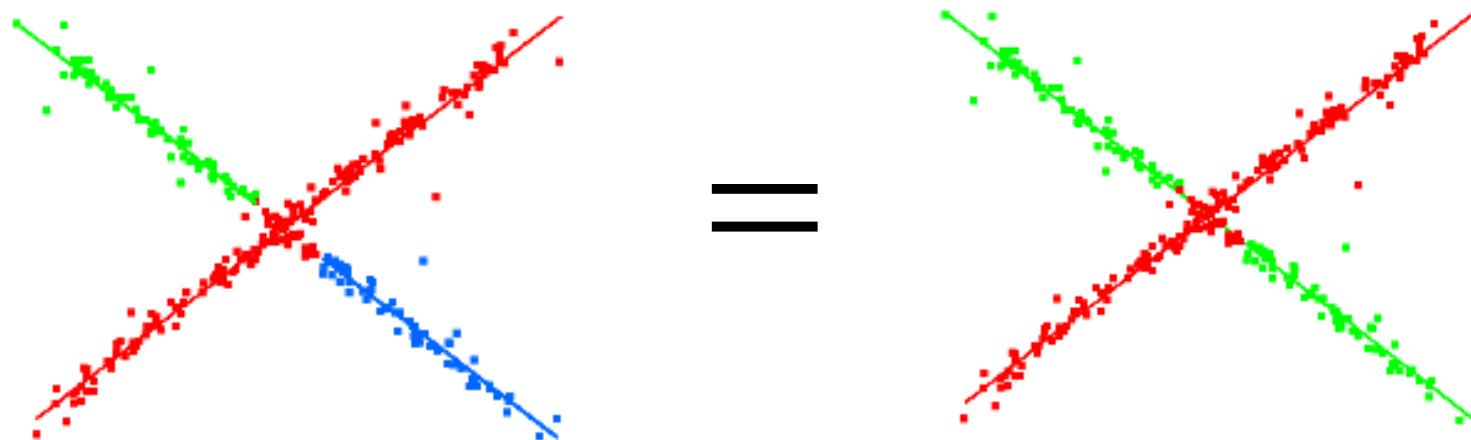
- find optimal labeling  
 $f = \{ f_1, f_2, \dots, f_n \}$



# Spatial regularization

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$$E(f) = \sum_p \| p, f_p \| + \sum_{pq \in N} \theta_{pq} \cdot [f_p \neq f_q]$$



Not enough!!!

# Energy-based approach

$$E(f) = \sum_p \| p, f_p \| + \sum_{f \in \Lambda} h_f \cdot \delta_f(f)$$

If **multiple** models

- assign different models (labels  $f_p$ ) to every point p

- find optimal labeling  
 $f = \{f_1, f_2, \dots, f_n\}$

$\Lambda$  - set of labels allowed at each point p

$$\delta_f(f) = \begin{cases} 1, & \exists p : f_p = f \\ 0, & \text{otherwise} \end{cases}$$

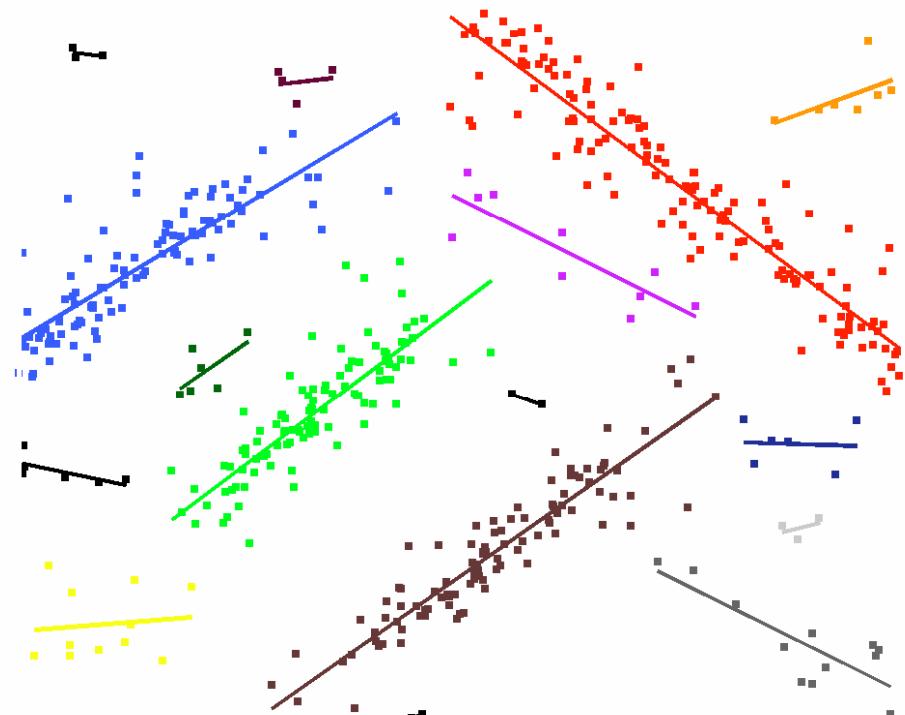
# Energy-based approach

$$E(f) = \sum_p \| p, f_p \| + \sum_{(p,q) \in N} \theta_{pq} \cdot [f_p \neq f_q] + \sum_{f \in \Lambda} h_f \cdot \delta_f(f)$$

If **multiple** models

- assign different models (labels  $f_p$ ) to every point  $p$

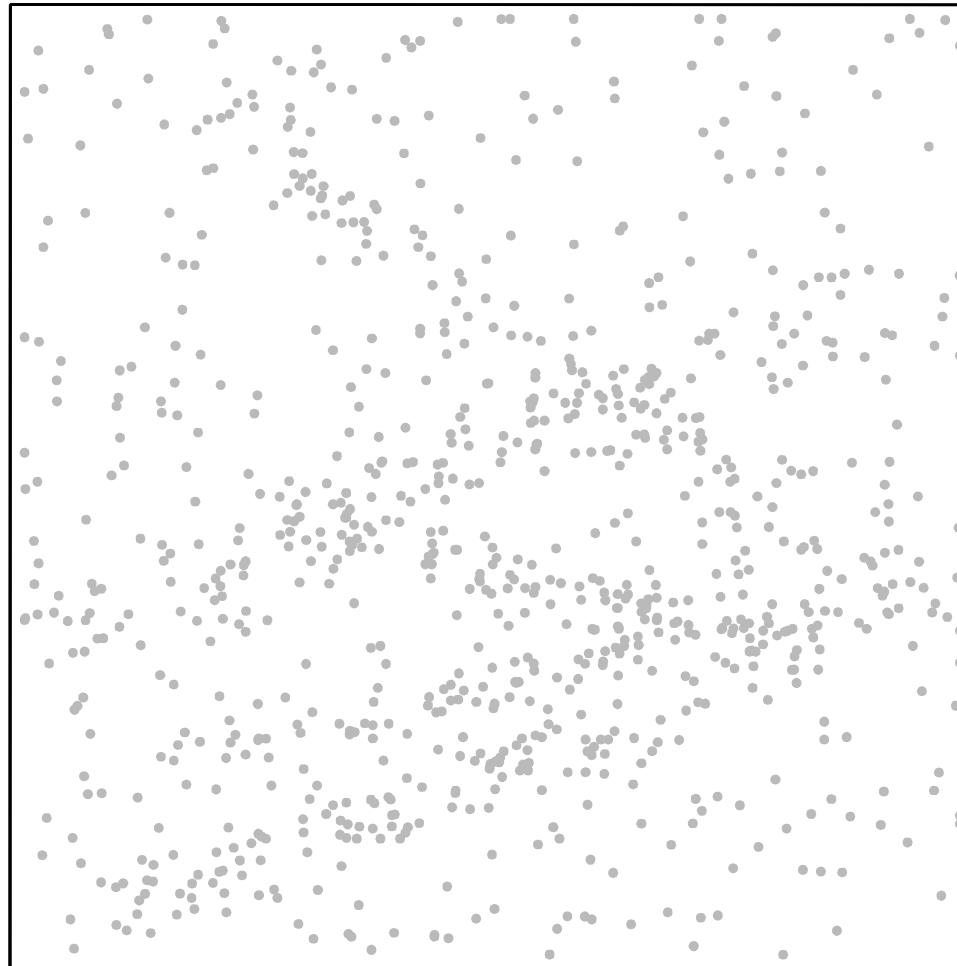
- find optimal labeling  
 $f = \{ f_1, f_2, \dots, f_n \}$



**Practical problem:** number of potential labels (models) is huge,  
how are we going to use a-expansion?

# PEARL

Propose  
Expand  
And  
Reestimate  
Labels

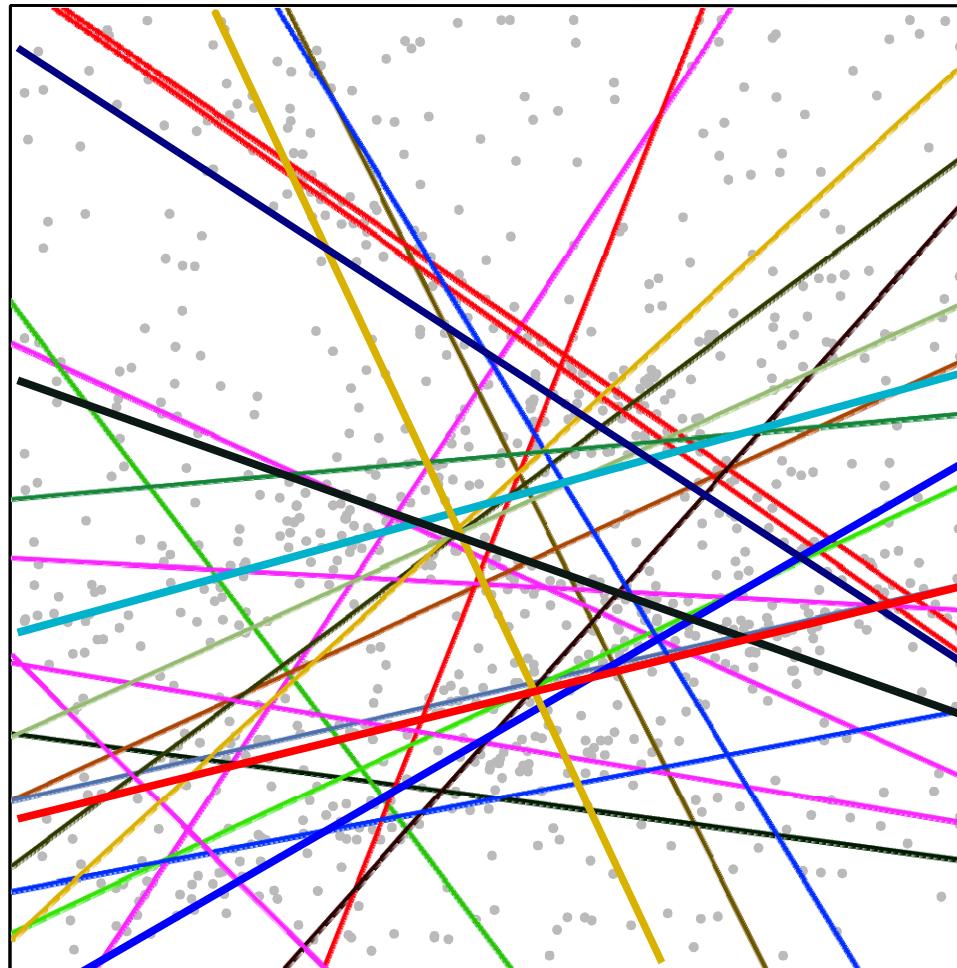


data points

# PEARL

---

Propose  
Expand  
And  
Reestimate  
Labels



data points + randomly sampled models

sample data  
to generate  
a **finite set**  
**of initial**  
**labels**

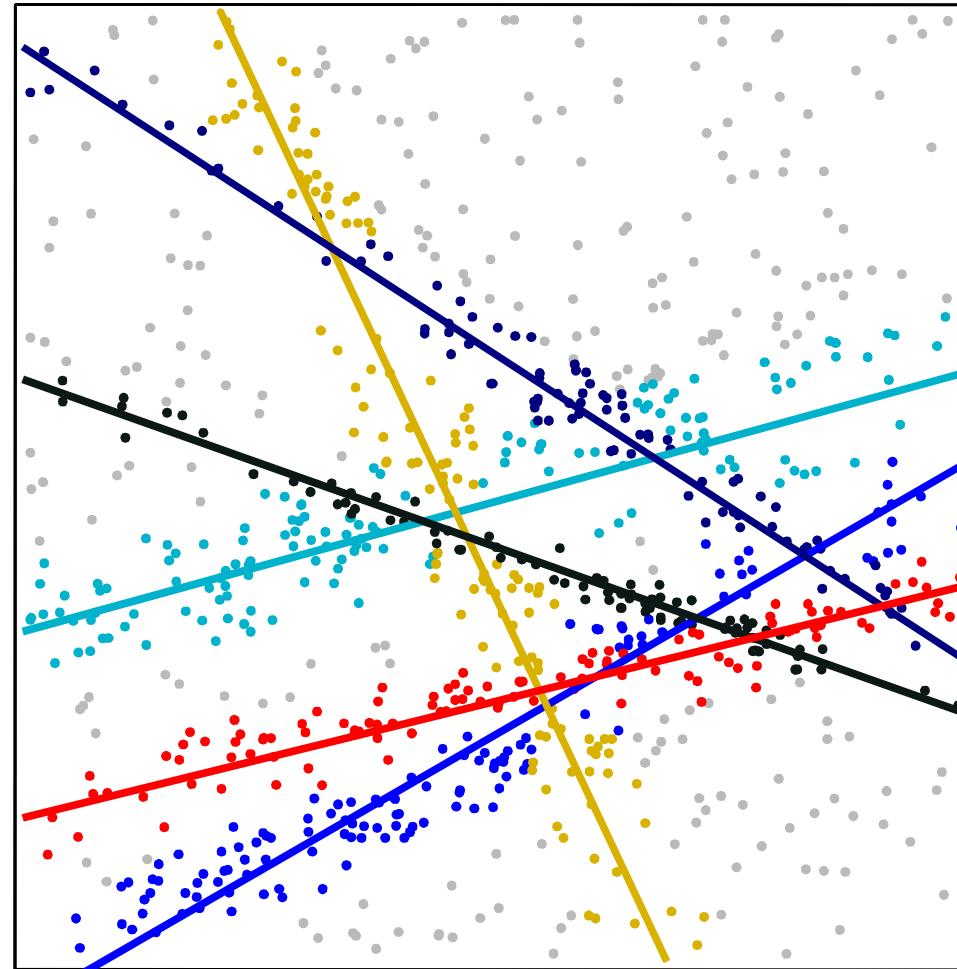
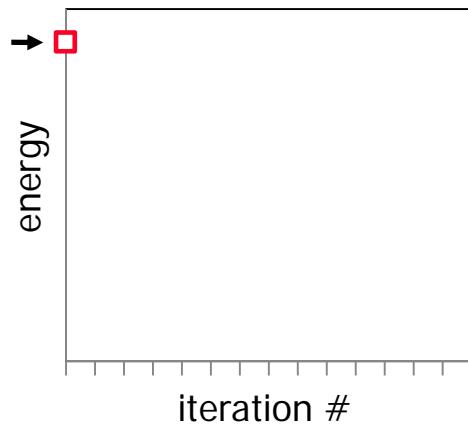
$\Lambda$

$$E(f) = \sum_p \| p, f_p \| + \sum_{pq \in N} \theta_{pq} \cdot [f_p \neq f_q]$$

# PEARL

---

Propose  
Expand  
And  
Reestimate  
Labels



**a-expansion:**  
minimize  $E(f)$   
over a fixed  
set of labels

$\Lambda$

$$f_p \in \Lambda$$

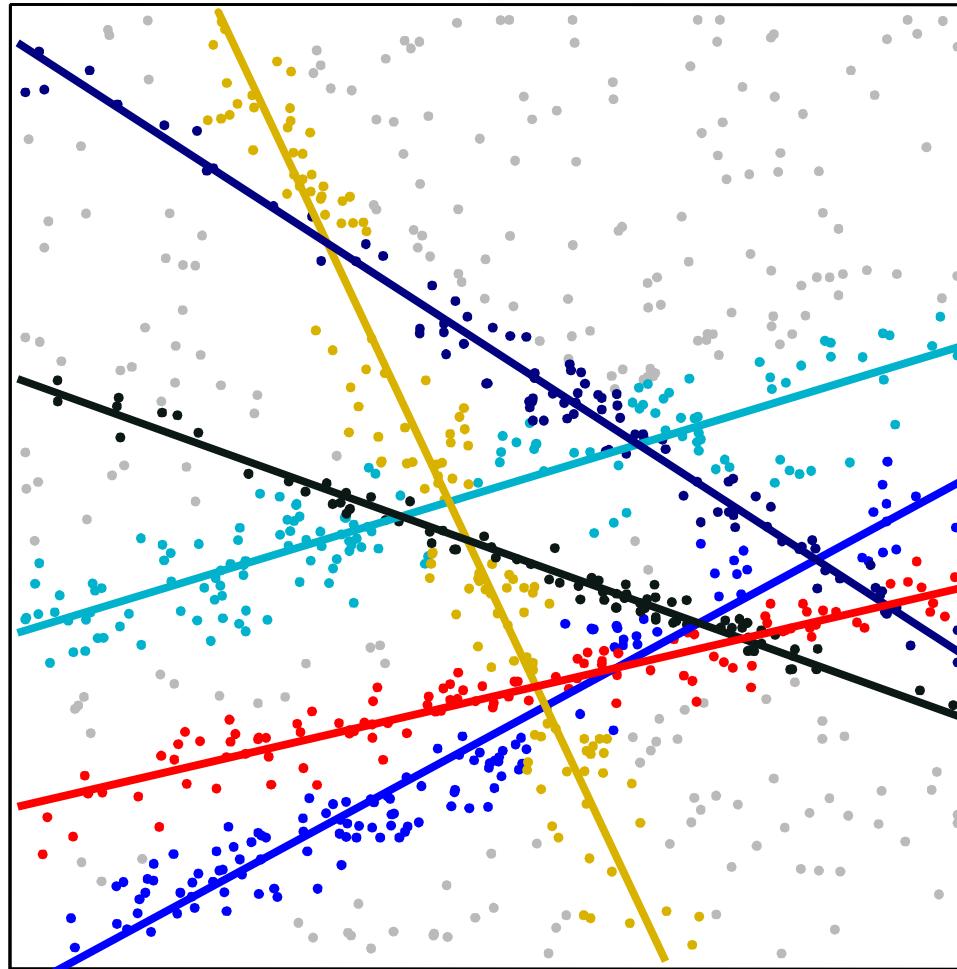
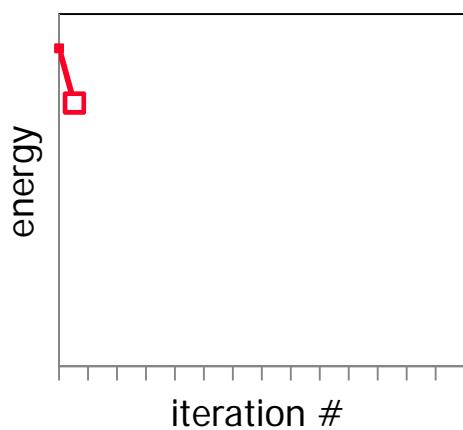
$$E(f) = \sum_p \| p, f_p \| + \sum_{pq \in N} \theta_{pq} \cdot [f_p \neq f_q]$$

fixed

# PEARL

---

Propose  
Expand  
And  
**Reestimate**  
Labels



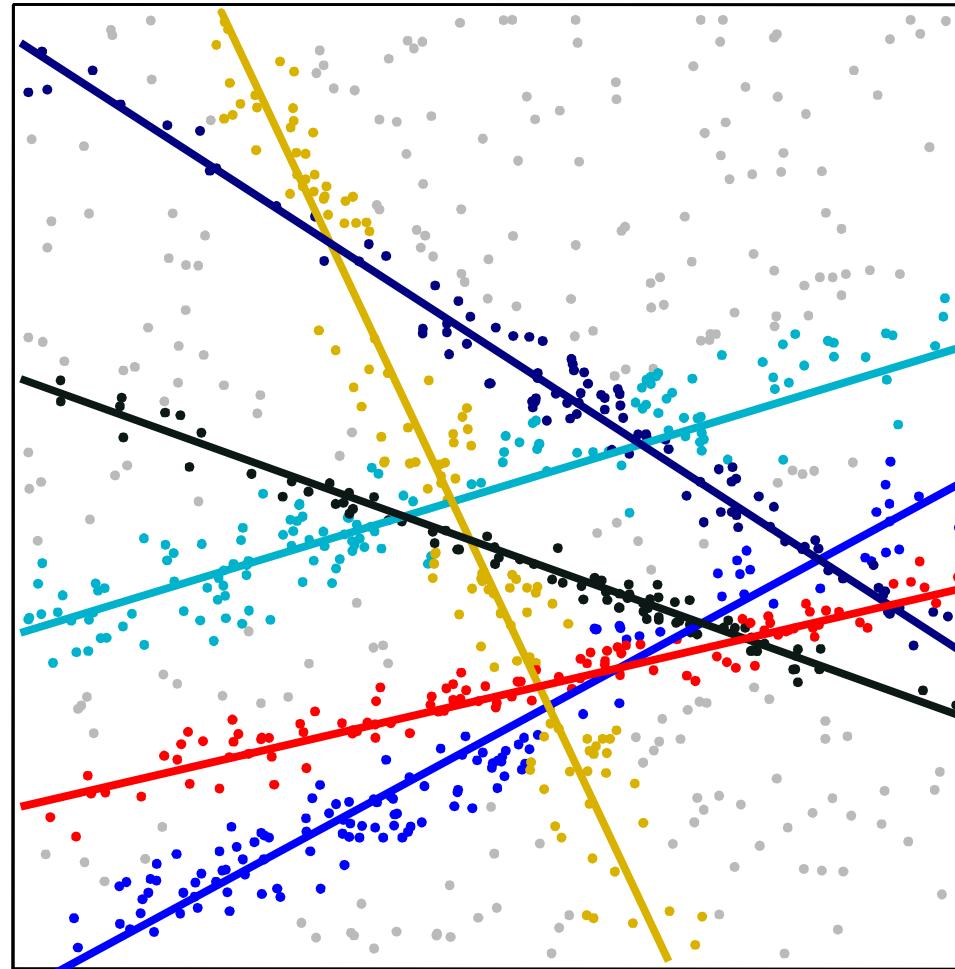
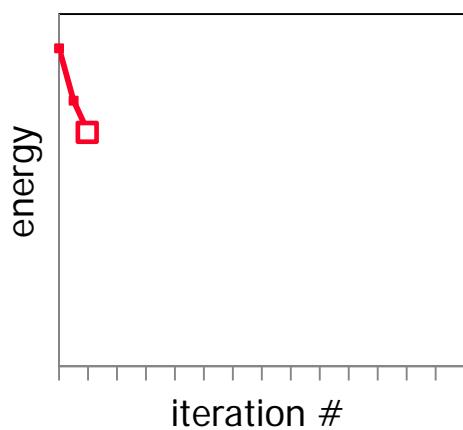
$$E(f) = \sum_p \| p, f_p \| + \sum_{pq \in N} \theta_{pq} \cdot [f_p \neq f_q]$$

$$f_p \in \Lambda$$

# PEARL

---

Propose  
Expand  
And  
Reestimate  
Labels



iteration 2: optimize labeling  $f$

**a-expansion:**  
minimize  $E(f)$   
over a fixed  
set of labels

$$\Lambda$$

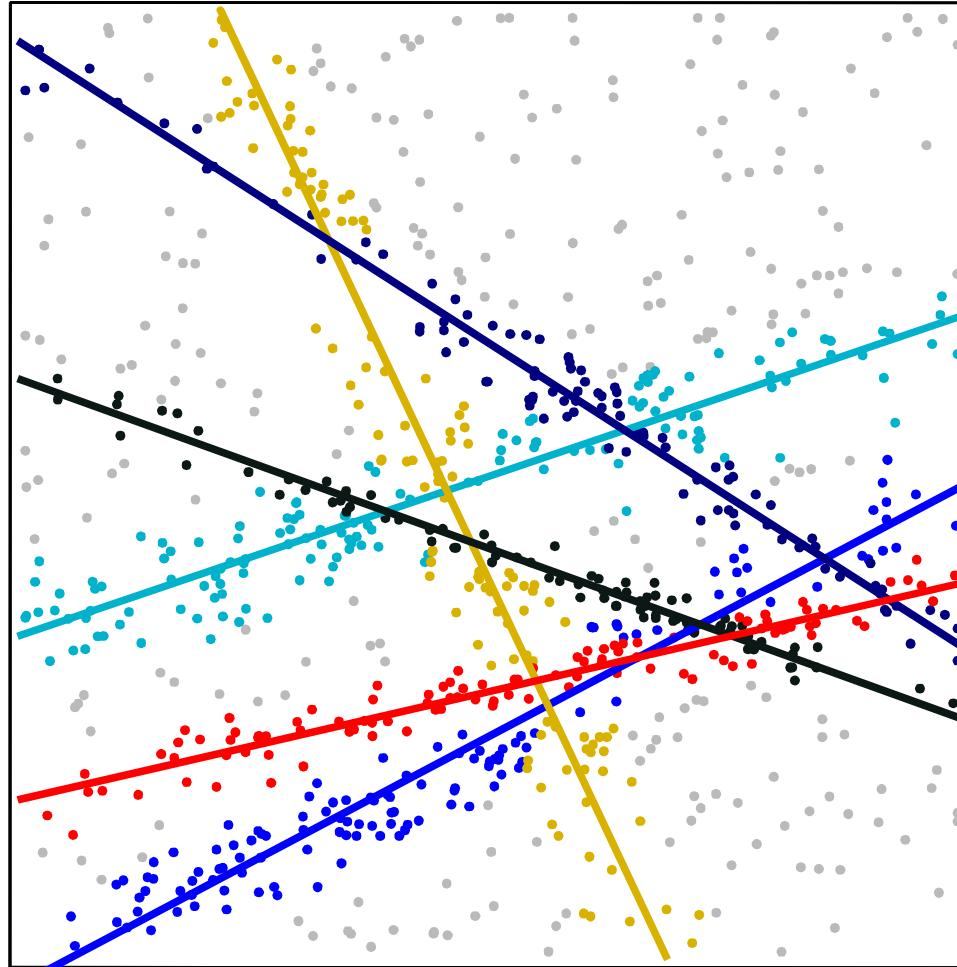
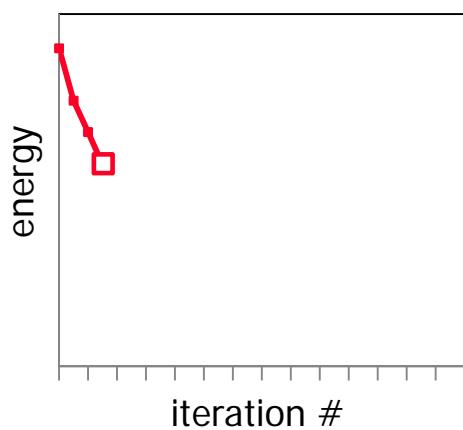
$$E(f) = \sum_p \| p, f_p \| + \sum_{pq \in N} \theta_{pq} \cdot [f_p \neq f_q]$$

fixed

# PEARL

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Propose  
Expand  
And  
**Reestimate**  
Labels

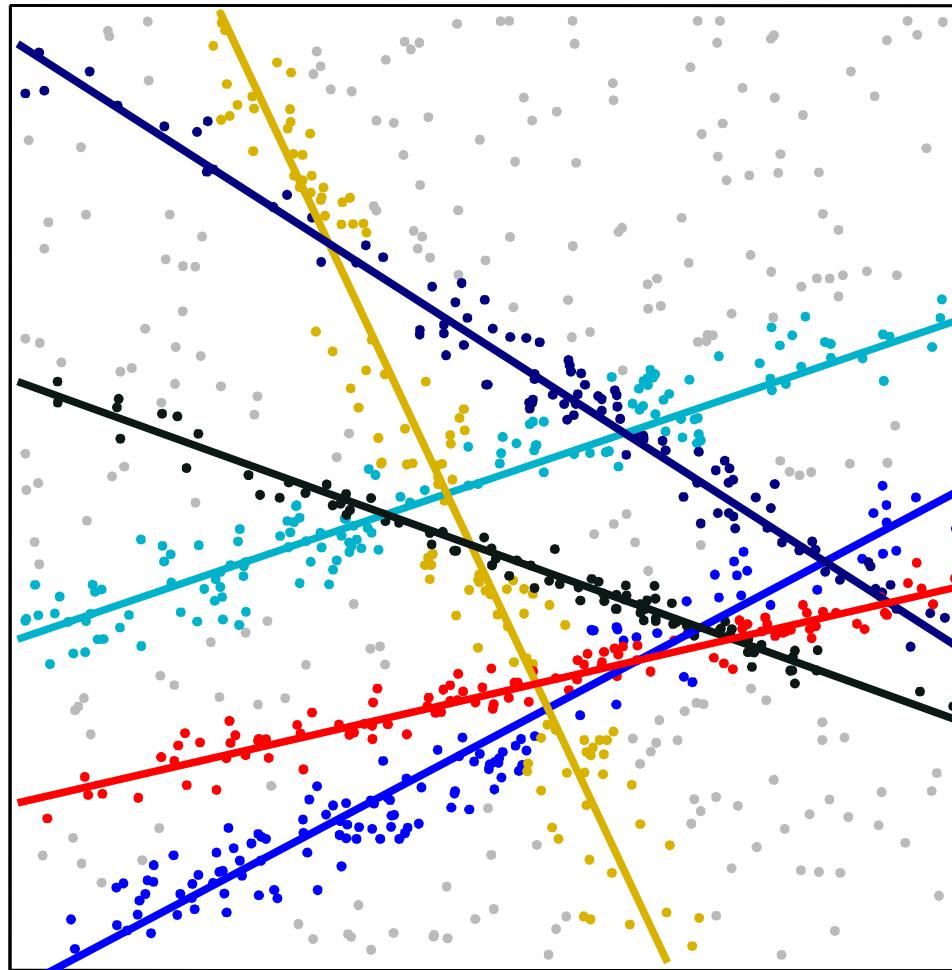
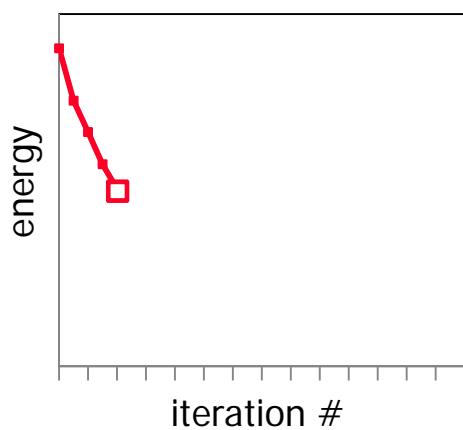


$$E(f) = \sum_p \| p, f_p \| + \sum_{pq \in N} \theta_{pq} \cdot [f_p \neq f_q]$$

# PEARL

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Propose  
Expand  
And  
Reestimate  
Labels

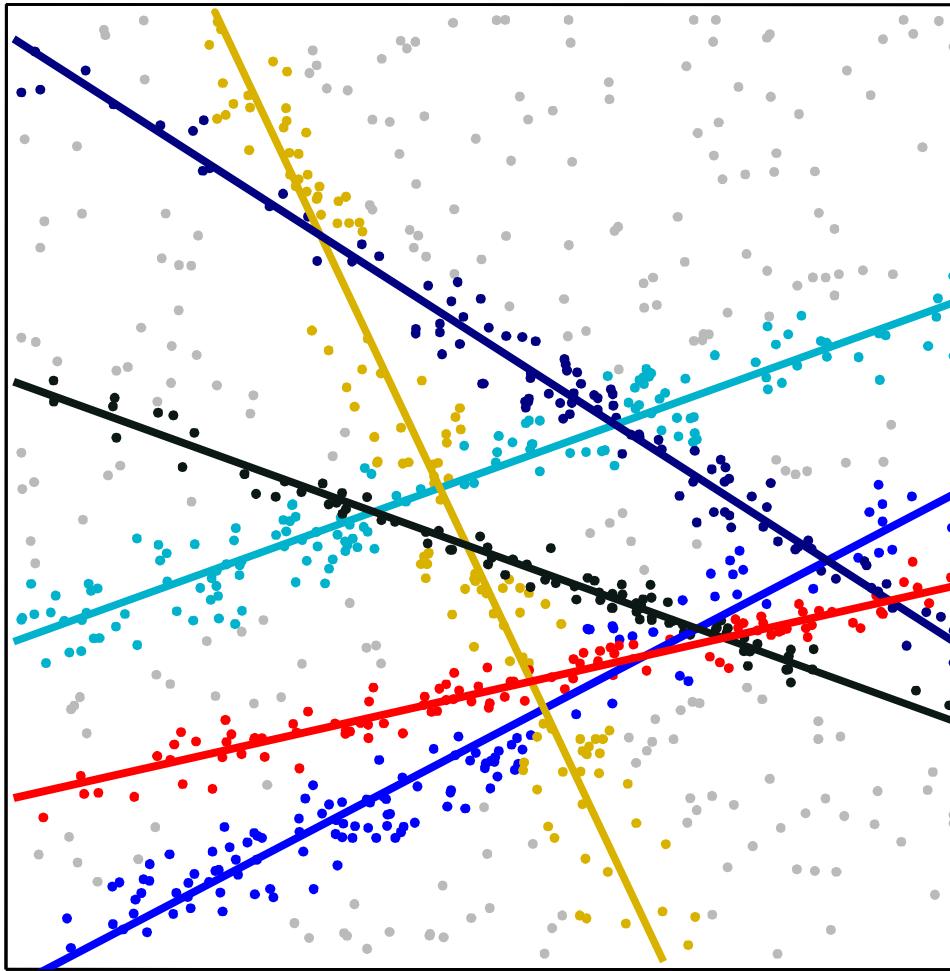
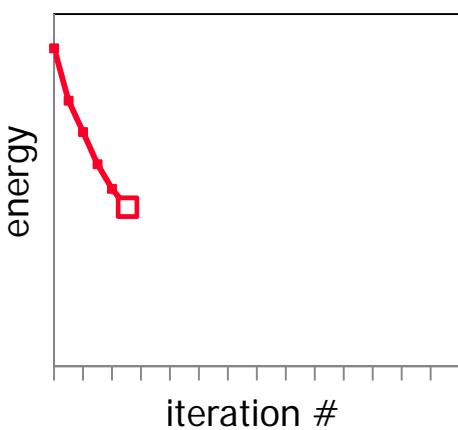


$$E(f) = \sum_p \| p, f_p \| + \sum_{pq \in N} \theta_{pq} \cdot [f_p \neq f_q]$$

# PEARL

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Propose  
Expand  
And  
**Reestimate**  
Labels

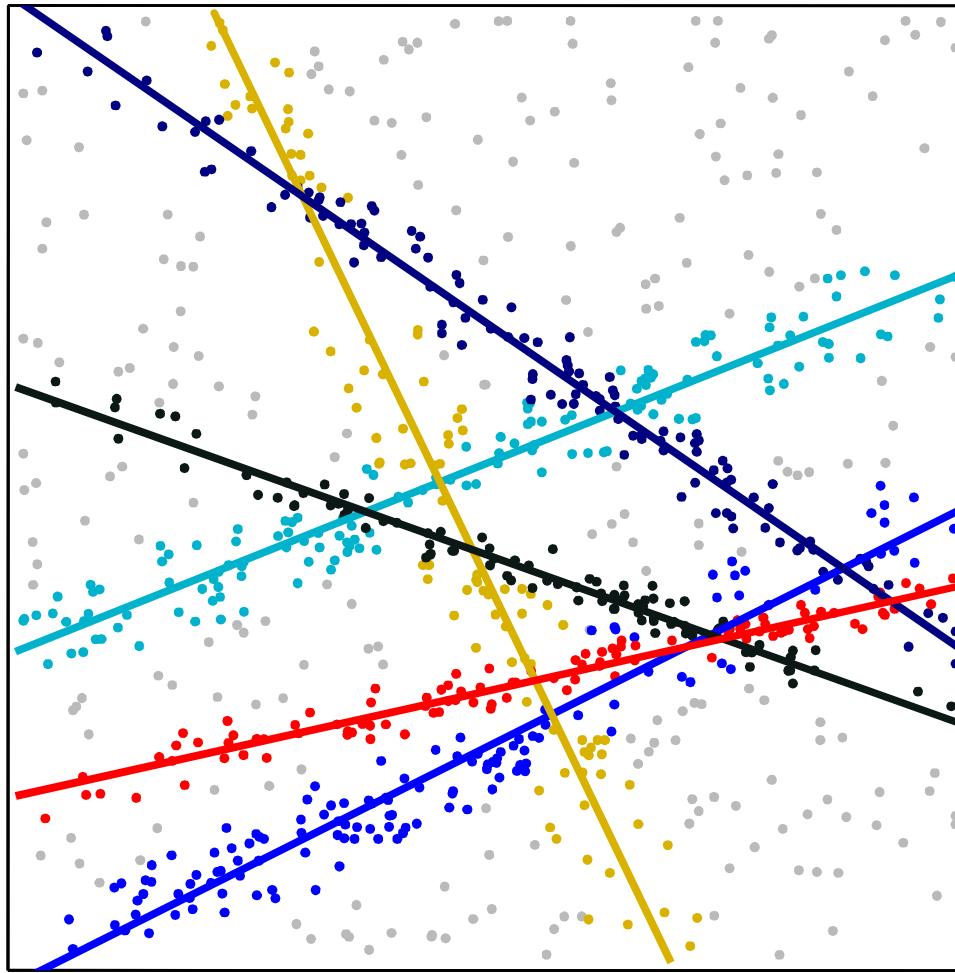
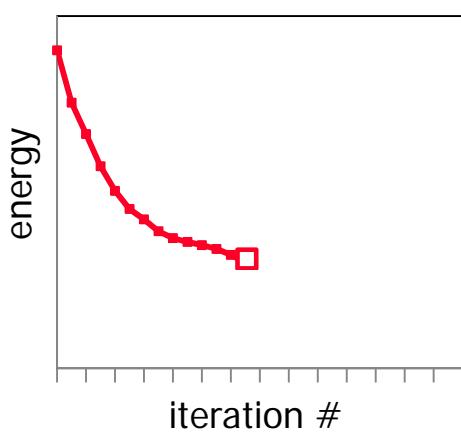


iteration 3: reestimate models

$$E(f) = \sum_p \| p, f_p \| + \sum_{pq \in N} \theta_{pq} \cdot [f_p \neq f_q]$$

# PEARL

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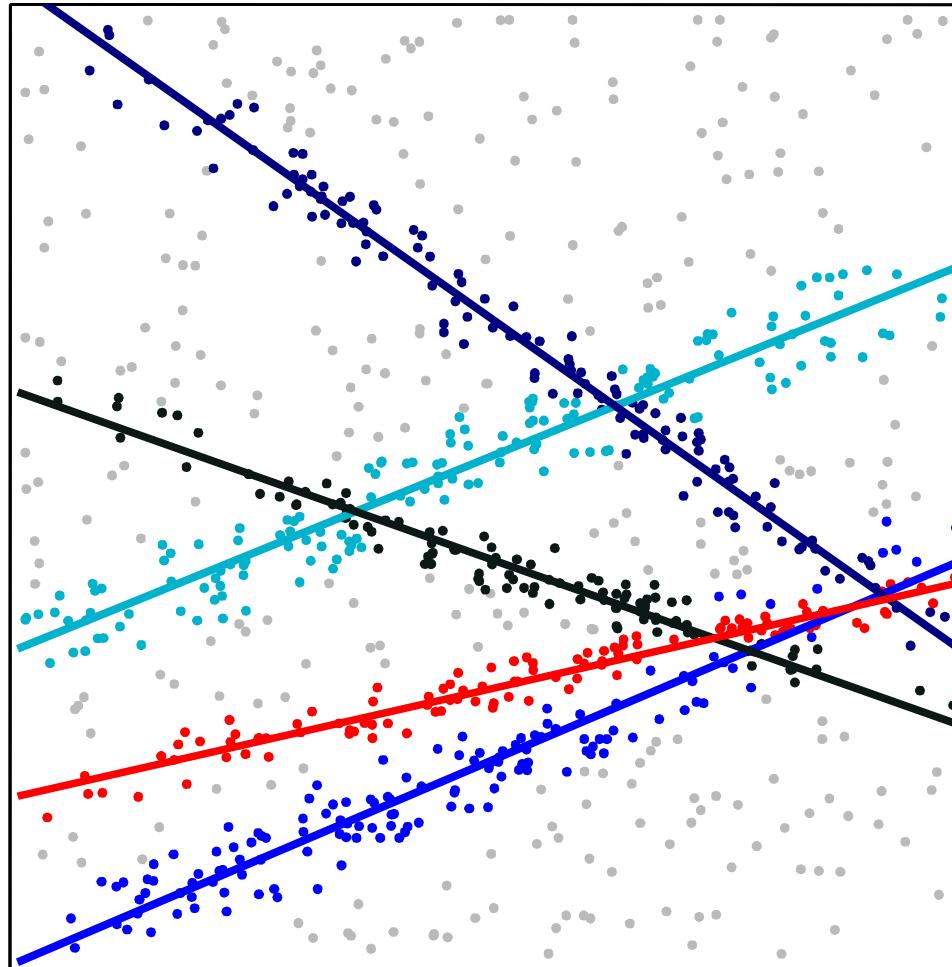
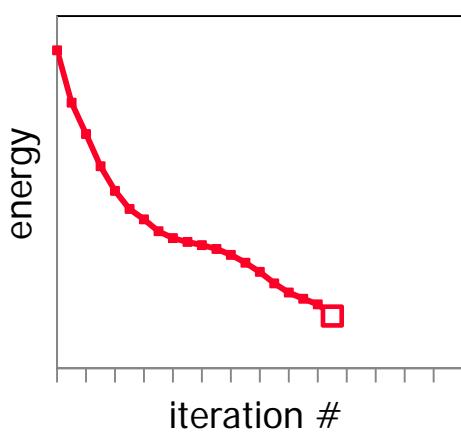


iteration 7...

$$E(f) = \sum_p \| p, f_p \| + \sum_{pq \in N} \theta_{pq} \cdot [f_p \neq f_q]$$

# PEARL

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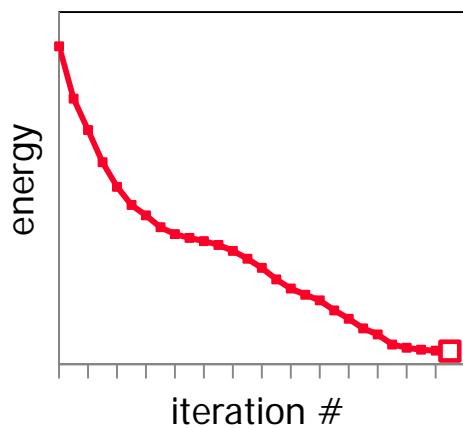
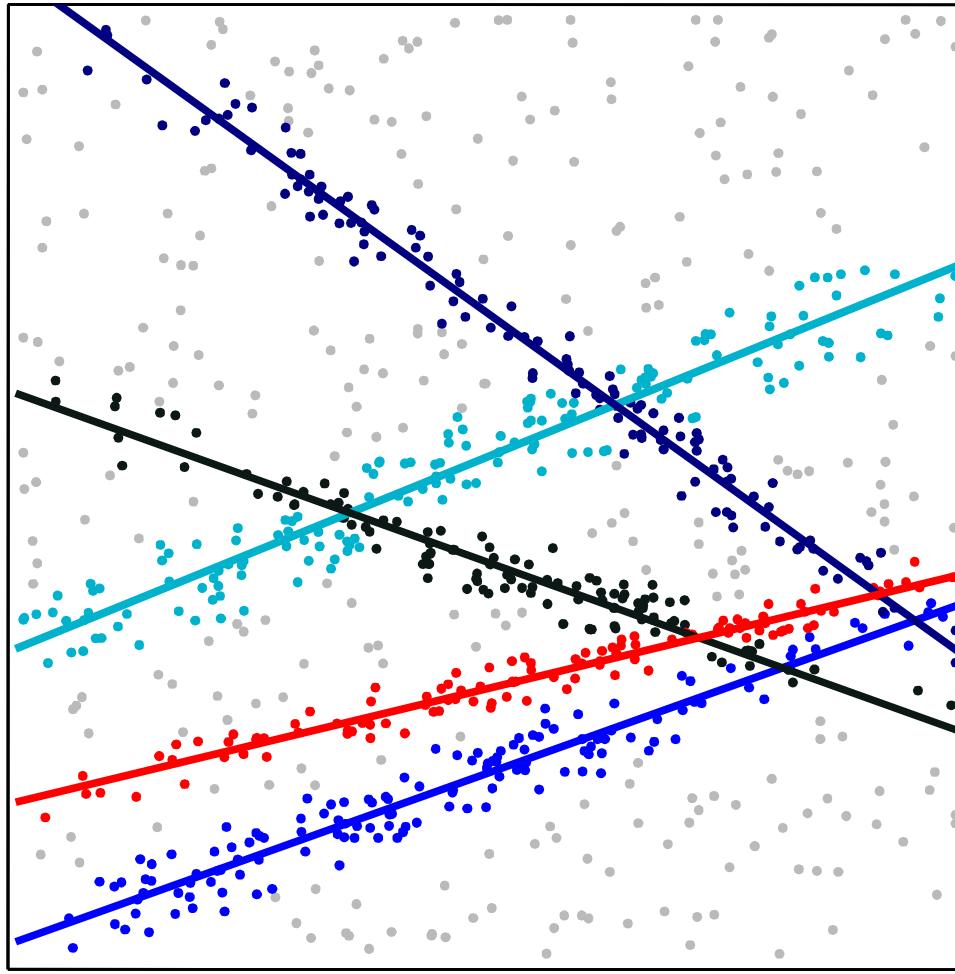


iteration 10...

$$E(f) = \sum_p \| p, f_p \| + \sum_{pq \in N} \theta_{pq} \cdot [f_p \neq f_q]$$

# PEARL

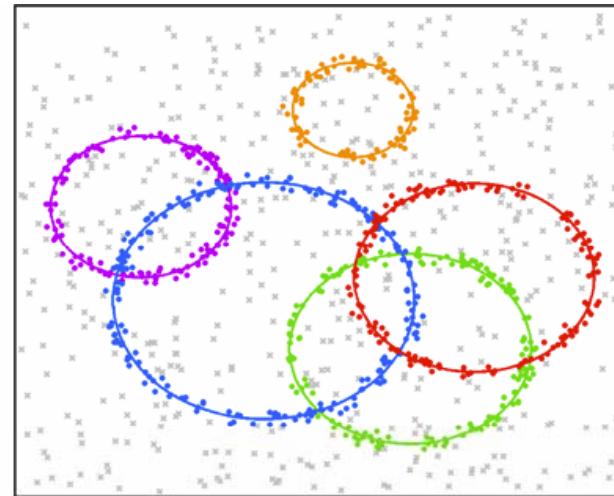
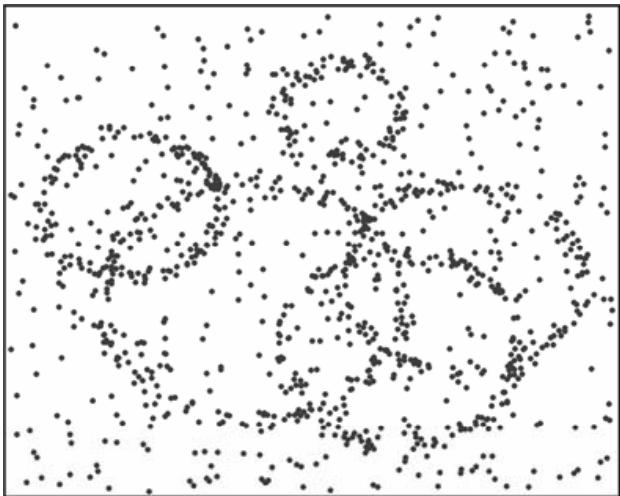
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iteration 15... converged.

# Fitting circles

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regularization with label costs only

Here spatial regularization does not work well

(unsupervised image segmentation)

## Fitting color models

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(unsupervised image segmentation)

## Fitting color models

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(c) Spatial regularity + label costs

Zhu and Yuille 96  
used continuous  
variational formulation  
(gradient descent)

(unsupervised image segmentation)

## Fitting color models

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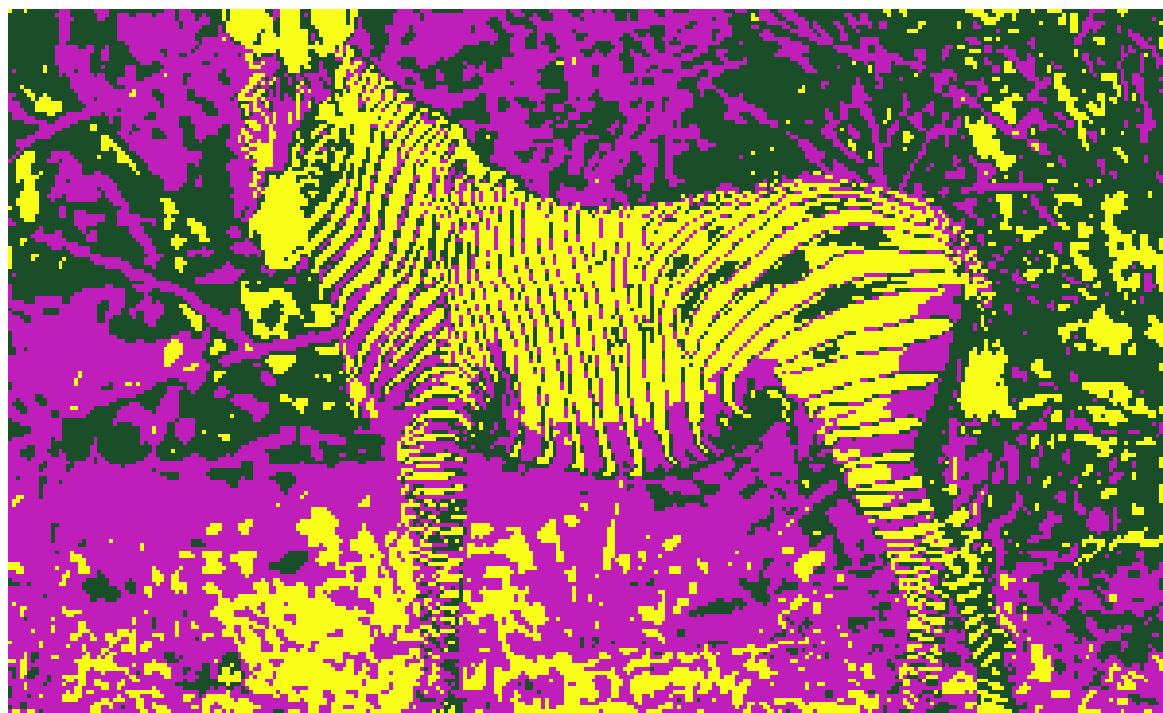


(b) Spatial regularity only [Zabih&Kolmogorov CVPR 04]

(unsupervised image segmentation)

## Fitting color models

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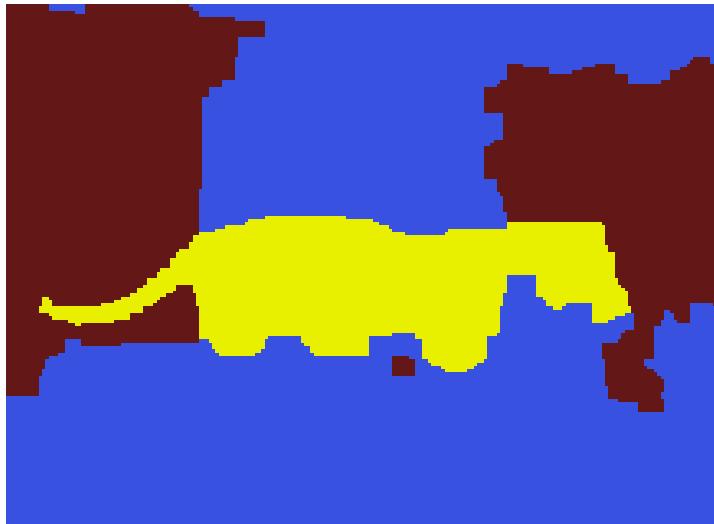


(a) Label costs only [Li, CVPR 2007]

(unsupervised image segmentation)

## Fitting color models

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Spatial regularity + label costs

(unsupervised image segmentation)

## Fitting color models

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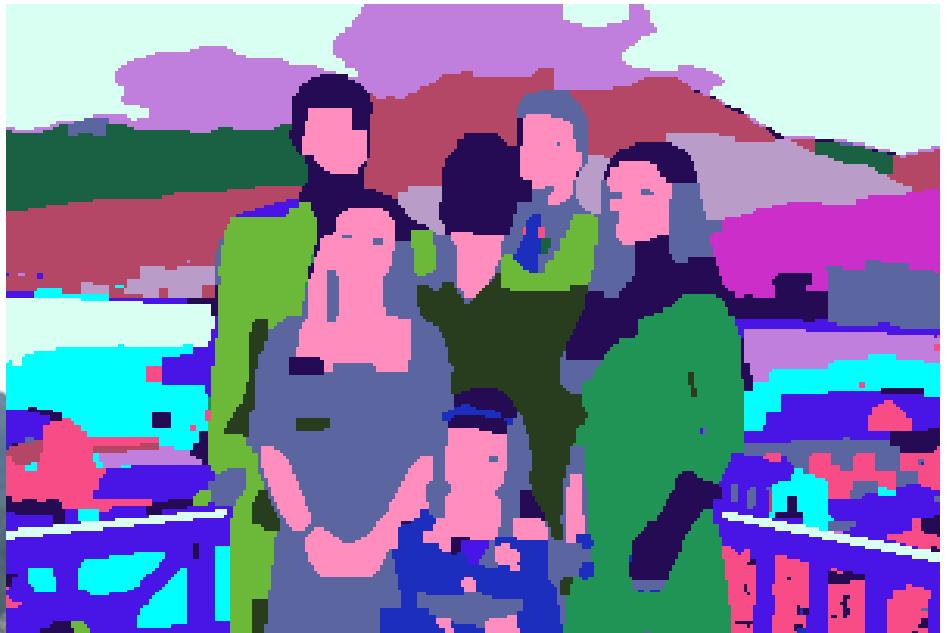


Spatial regularity + label costs

(unsupervised image segmentation)

## Fitting color models

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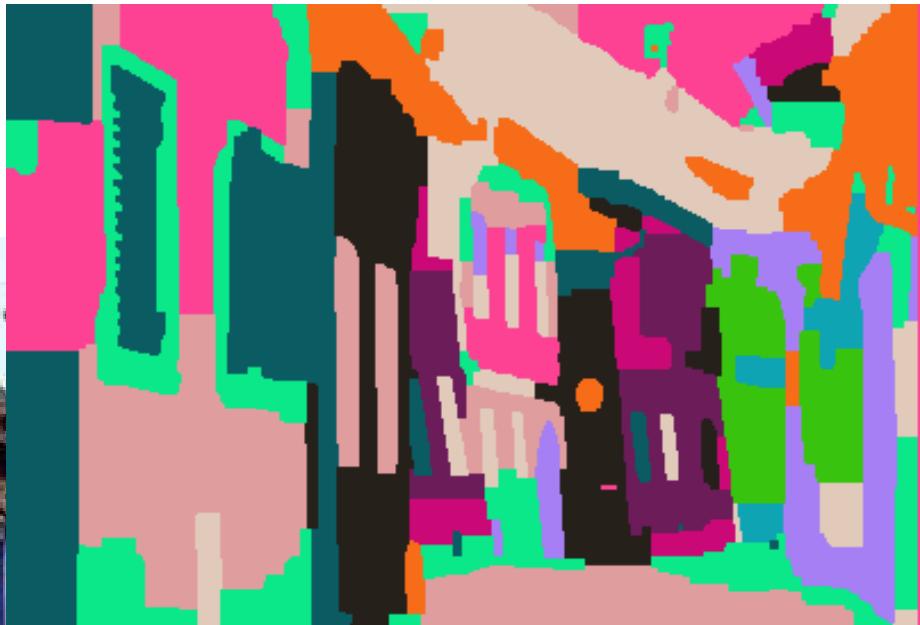


Spatial regularity + label costs

(unsupervised image segmentation)

## Fitting color models

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Spatial regularity + label costs

# Fitting planes (homographies)

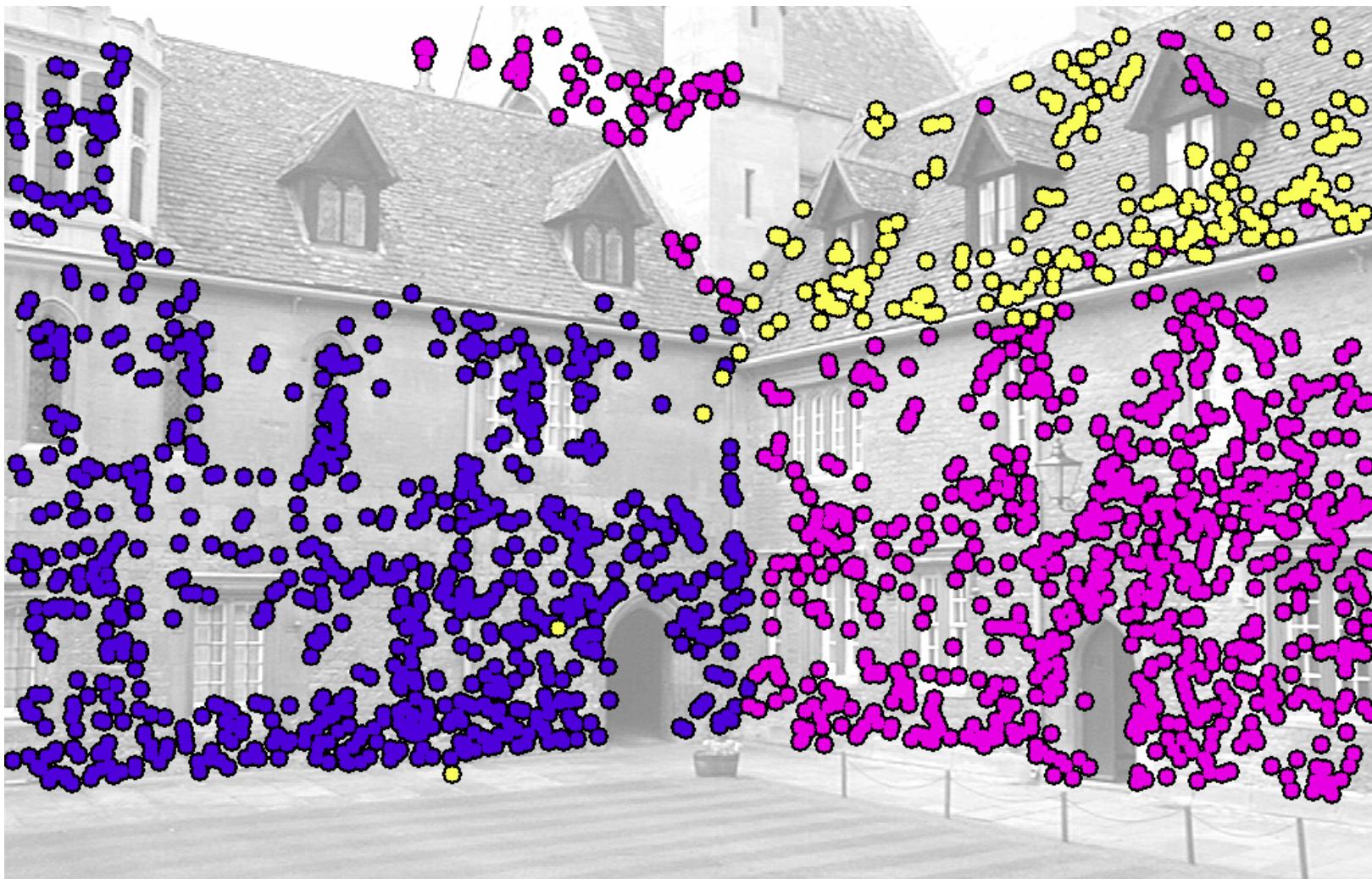
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Original image (one of 2 views)

# Fitting planes (homographies)

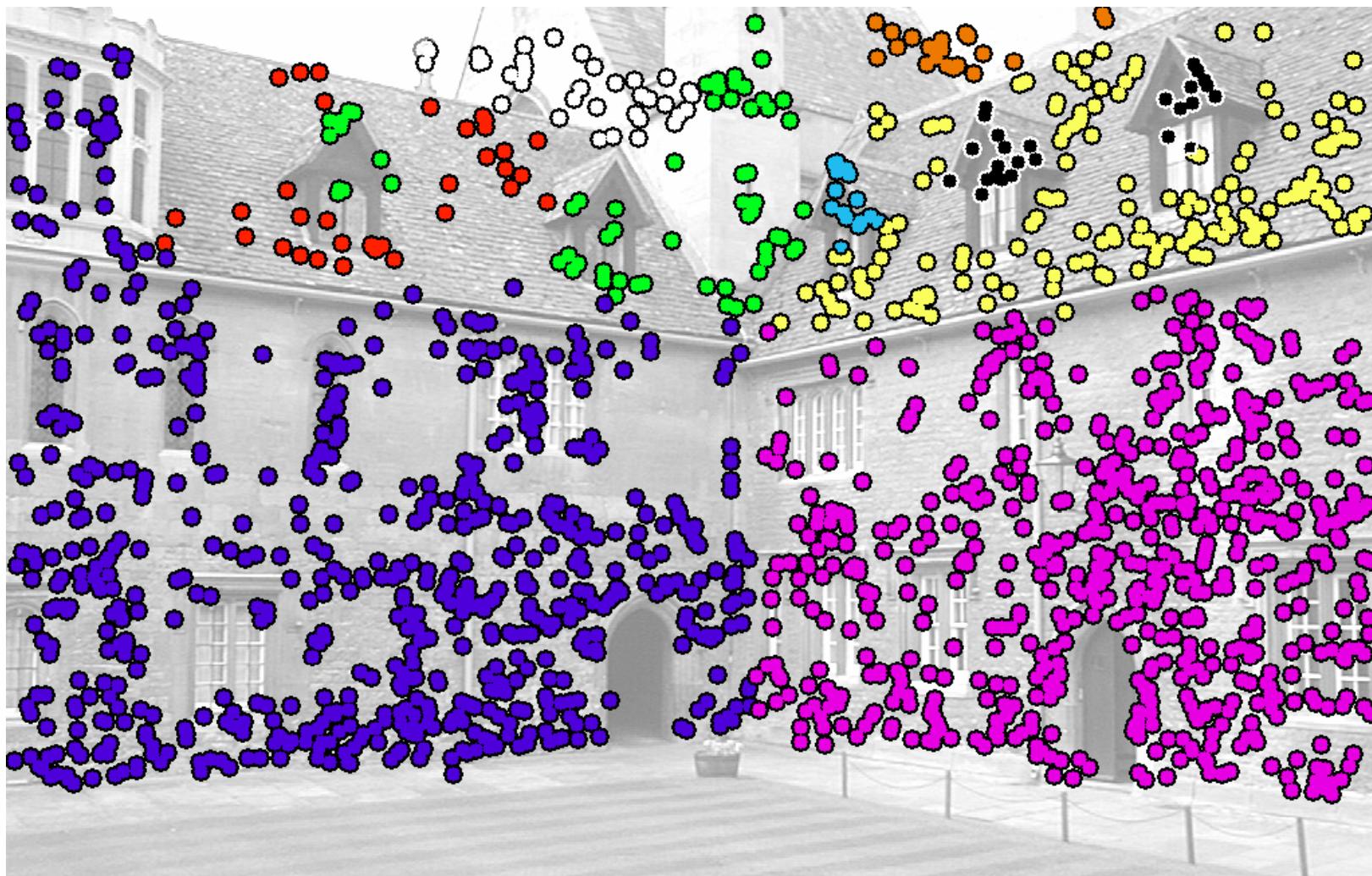
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(a) Label costs only

# Fitting planes (homographies)

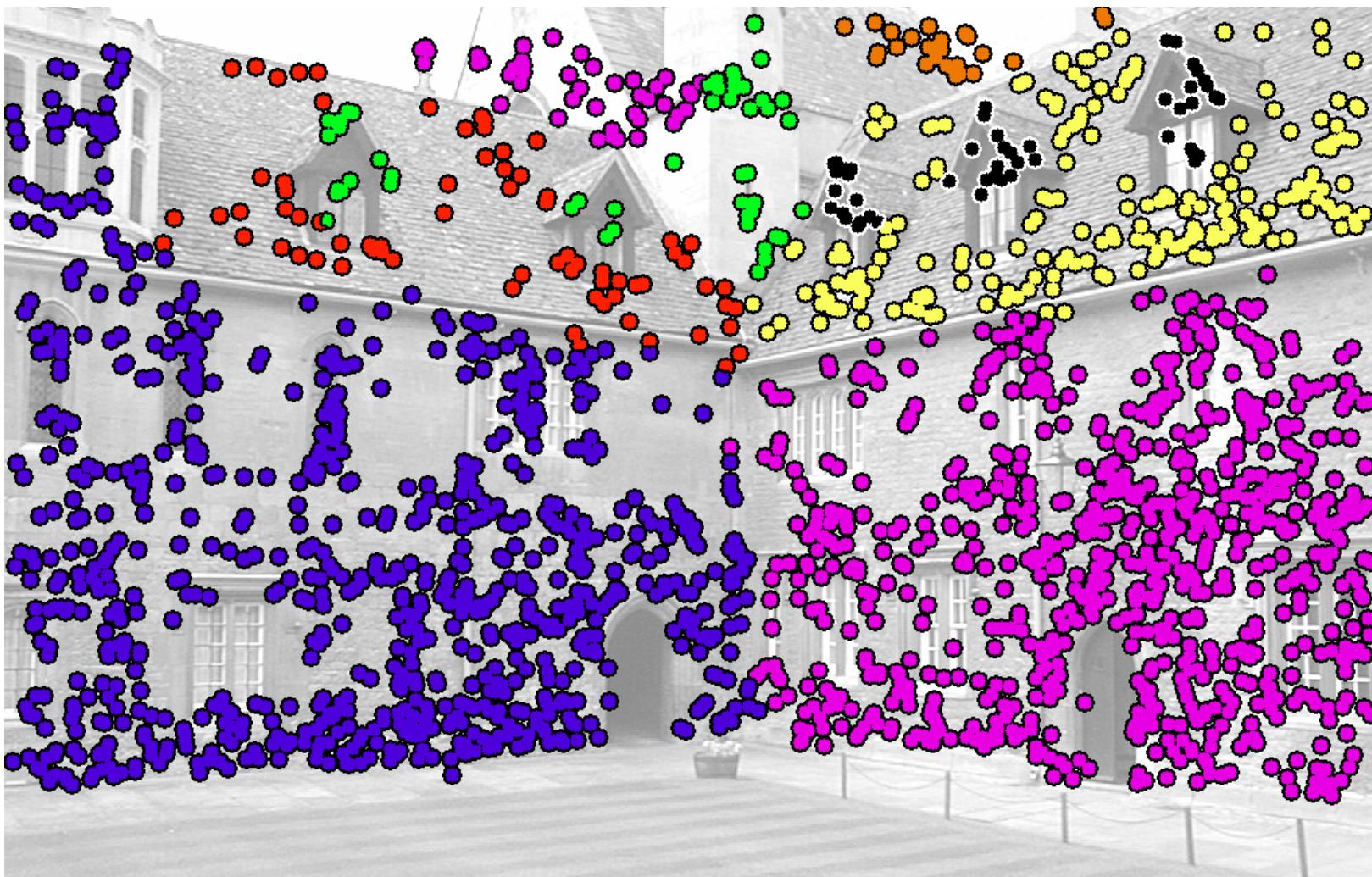
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(b) Spatial regularity only

# Fitting planes (homographies)

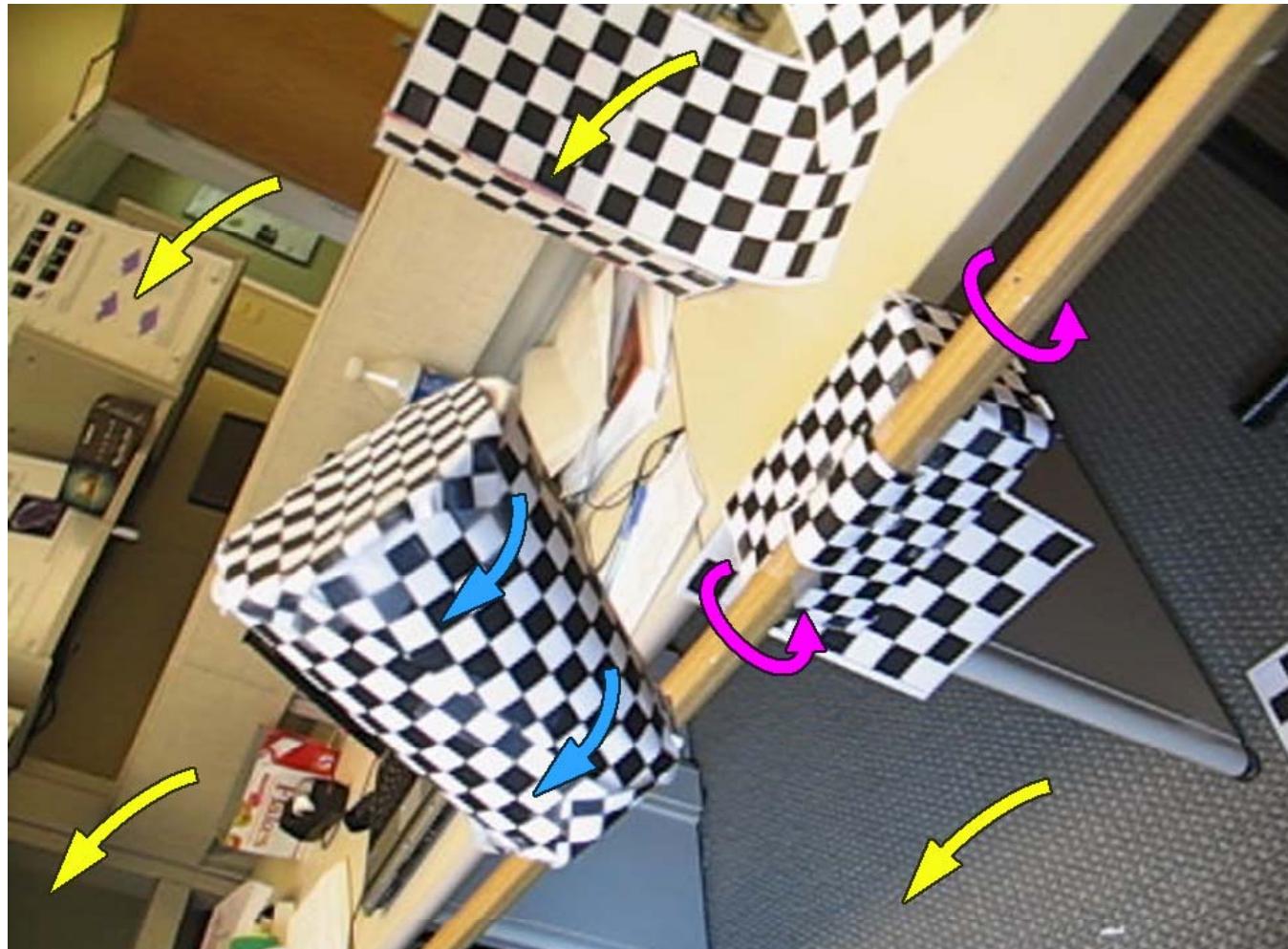
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(c) Spatial regularity + label costs

# (rigid) Motion Estimation

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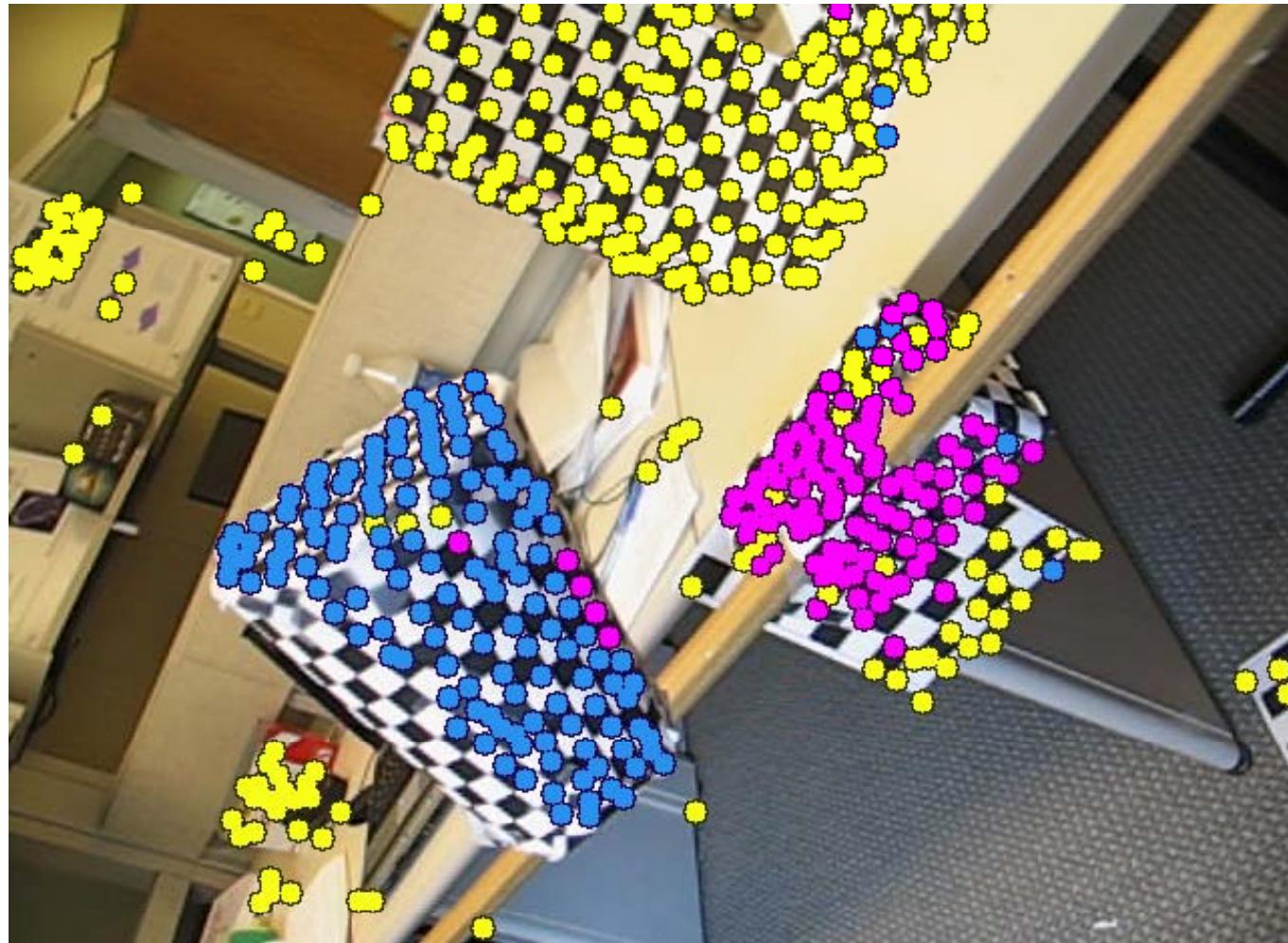


Original image

[Rene Vidal]

# (rigid) Motion Estimation

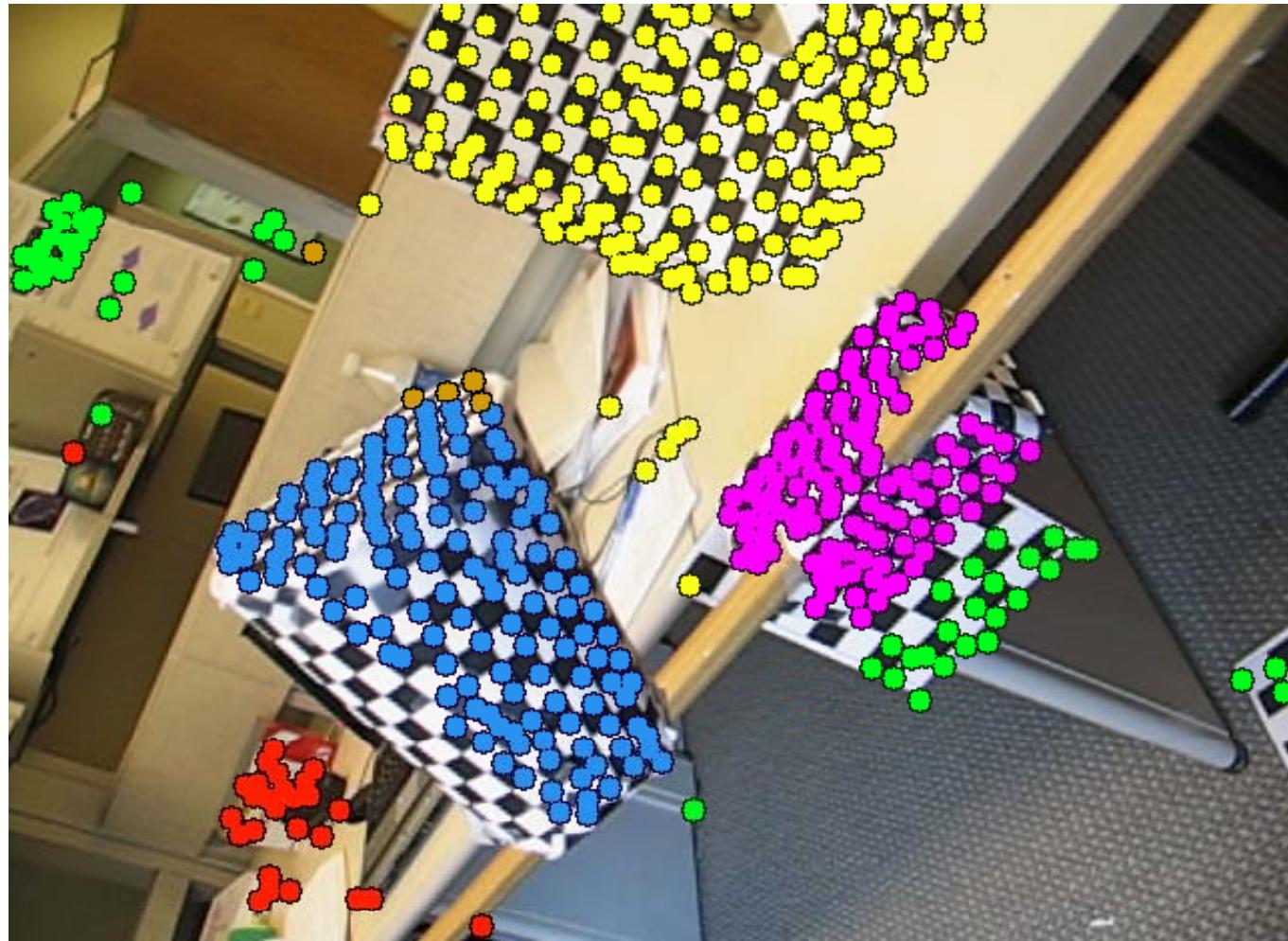
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(a) Label costs only

# (rigid) Motion Estimation

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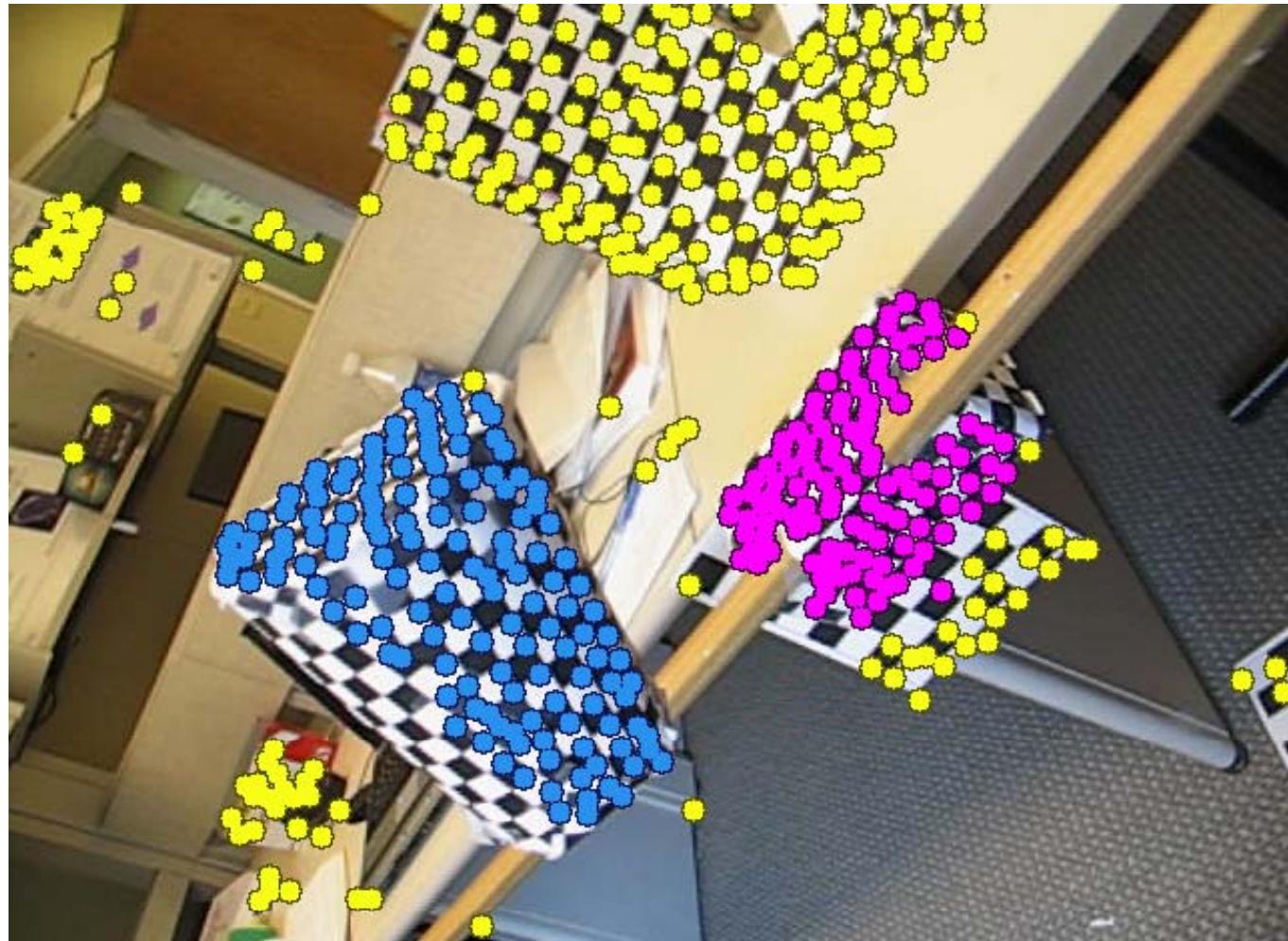


(b) Spatial regularity only

7  
motions

# (rigid) Motion Estimation

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(c) Spatial regularity + label costs

# (rigid) Motion Estimation

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# Multi-Label Energy Formulation

Input: Set of data points  $P$  ← pixels, features,  
matches,...  
Set of candidate labels  $\Lambda$  ← objects, motions,  
homographies,...

Goal: Labeling  $f$  that minimizes energy  $E$

$$\min_{labeling} \left\{ \text{data costs} + \text{smooth costs} + \text{label costs} \right\}$$

$$E(f) = \sum_p D_p(f_p) + \sum_{(p,q)} V_{pq}(f_p, f_q) + \sum_l h_l \delta(\exists p : f_p = l)$$

## $\alpha$ -expansion

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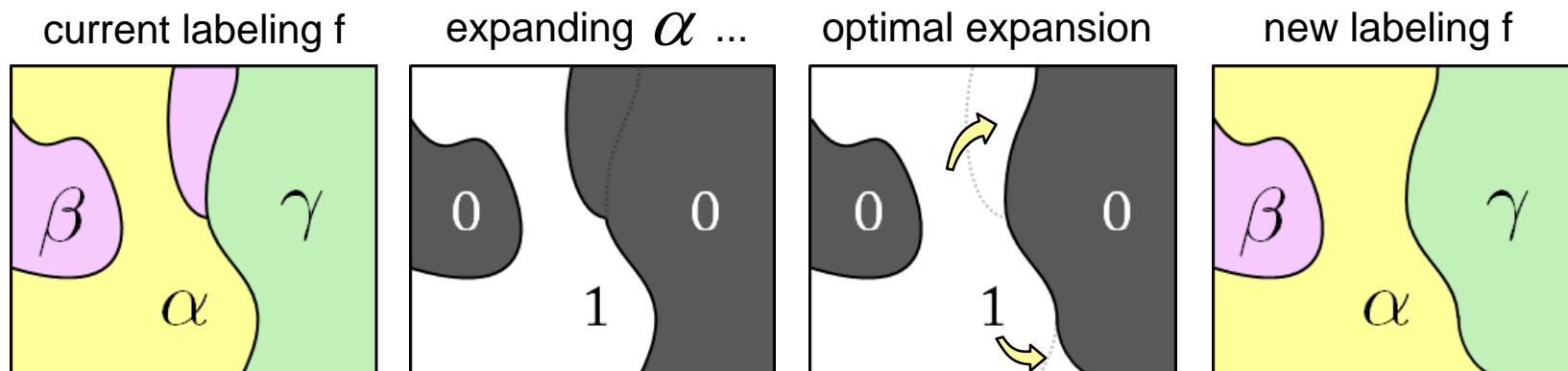
$$E(f) = \sum_p D_p(f_p) + \sum_{(p,q)} V_{pq}(f_p, f_q)$$

- NP-hard in general when  $|\Lambda| \geq 3$
- $\alpha$ -expansion is standard algorithm
  - finds local optimum w.r.t. “expansions”
  - optimality guarantees
  - fast & effective in practice

# $\alpha$ -expansion

## ■ $\alpha$ -expansion main idea:

- convert multi-label problem into sequence of *binary* problems
- choose label  $\alpha$  , and only let it “expand”



## Deriving the Graph Construction

Let  $f$  be current labeling

$\mathbf{x}$  be labeling of binary subproblem

$f^\alpha$  be the labeling induced by  $\mathbf{x}$

$$\begin{array}{lcl} x_p = 0 & \iff & f_p^\alpha = f_p \xleftarrow{\text{keep current label}} \\ x_p = 1 & \iff & f_p^\alpha = \alpha \xleftarrow{\text{switch to } \alpha} \end{array}$$

# Deriving the Graph Construction

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- How to add cost  $h_\beta$  to binary problem?

$$E_h^\alpha(x) = E^\alpha(x) + h_\beta \underbrace{(1 - x_1 x_5 x_6)}$$

indicator function  $\delta_\beta(f^\alpha) = \begin{cases} 1 & \exists p : f_p^\alpha = \beta \\ 0 & \text{otherwise.} \end{cases}$

$f$	$\beta$	$\alpha$	$\gamma$	$\gamma$	$\beta$	$\beta$
$x$	?	1	?	?	?	?
	1	2	3	4	5	6

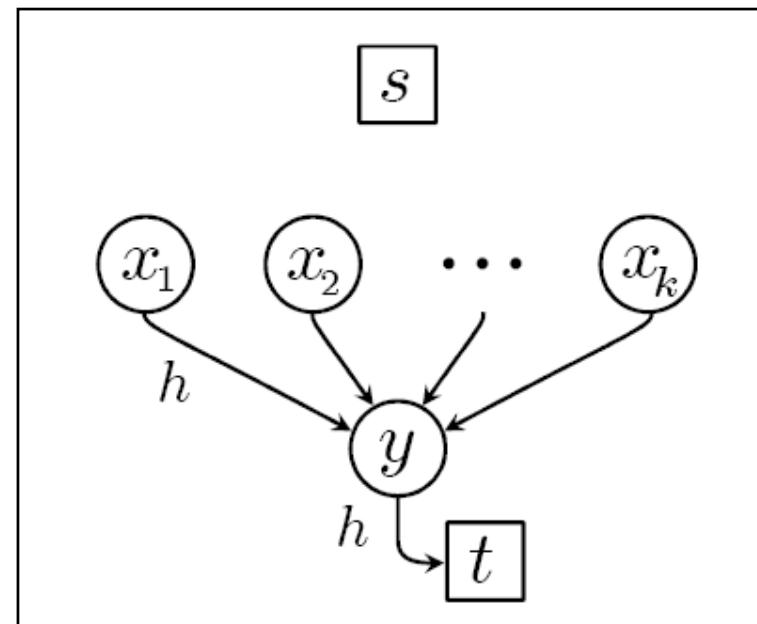
Modified energy  $E_h^\alpha(\mathbf{x})$   
pays  $h_\beta$  iff  $f^\alpha$  contains  
label  $\beta$

# Deriving the Graph Construction

Add one auxiliary variable:

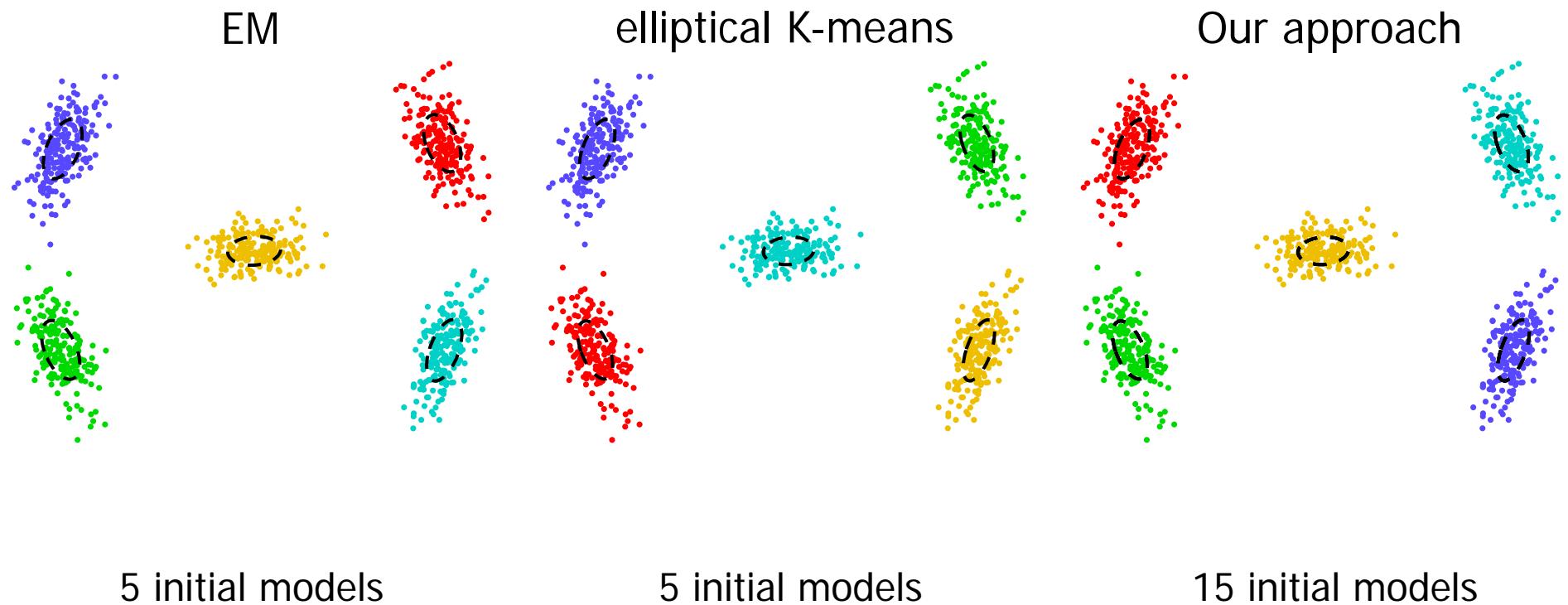
$$h(1 - x_1 x_5 x_6) = h \min_{y \in \{0,1\}} [\bar{y} + \bar{x}_1 y + \bar{x}_5 y + \bar{x}_6 y]$$

Same in terms of graph:



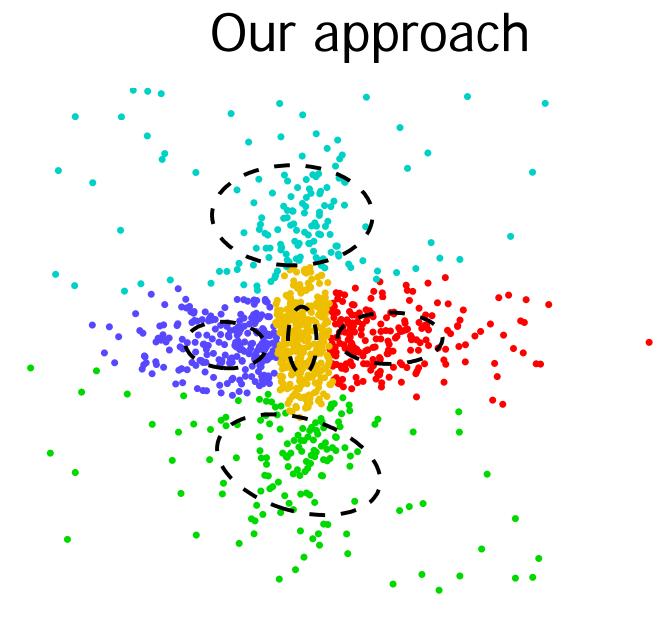
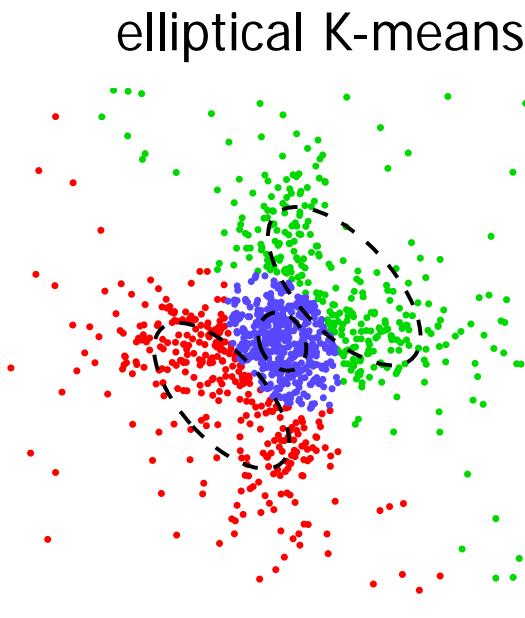
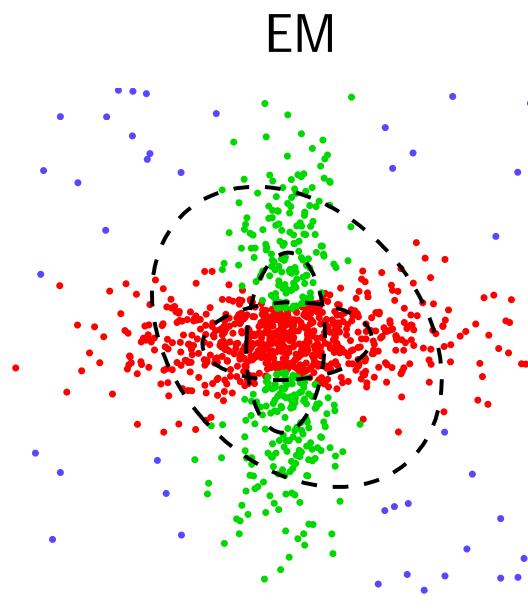
# FMM

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# FMM

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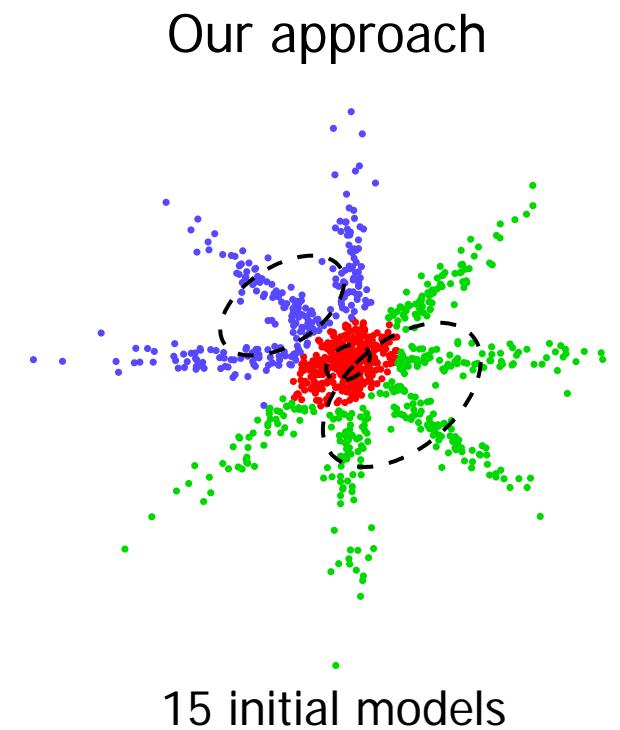
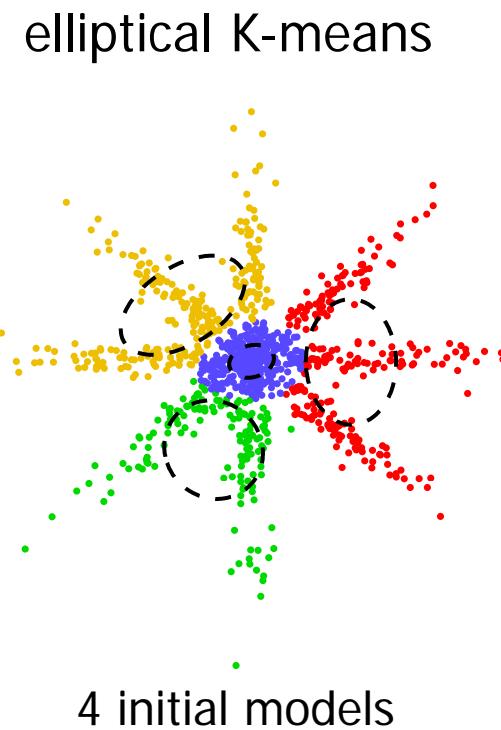
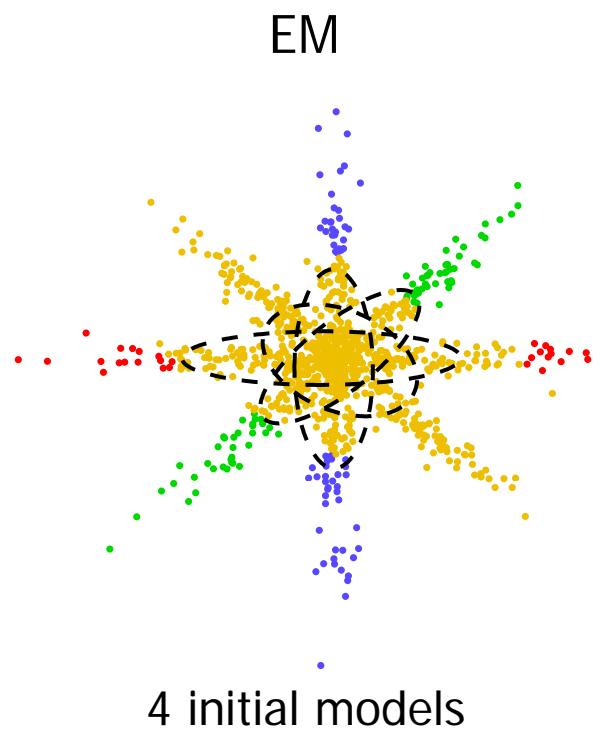


*works well for  
overlapping models  
due to **soft assignments***

***hard assignments*** fail if models overlap spatially

# FMM

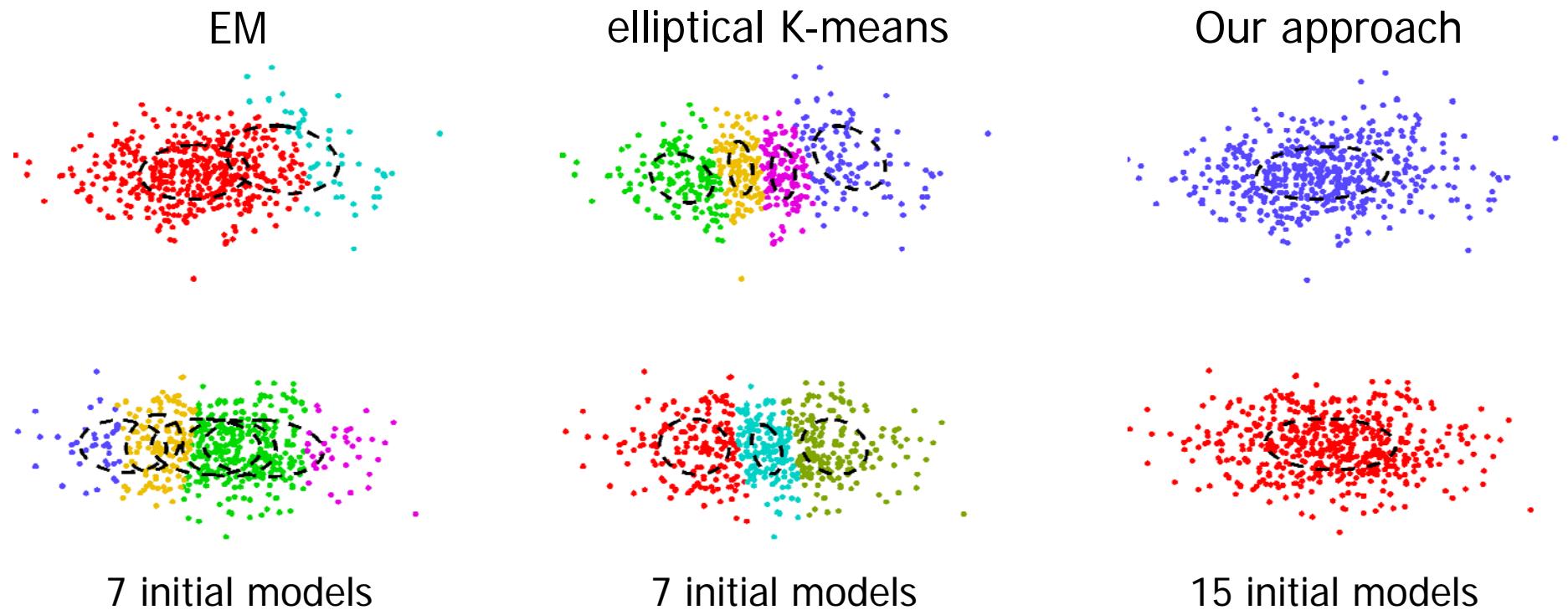
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***soft assignments***  
*is not a panacea*

# FMM

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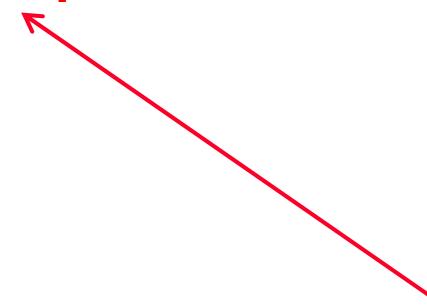


*standard techniques must know the exact number of models*

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*observation:*

our labeling approach makes hard assignments  
which may cause problems if  
**models have spatial overlap**



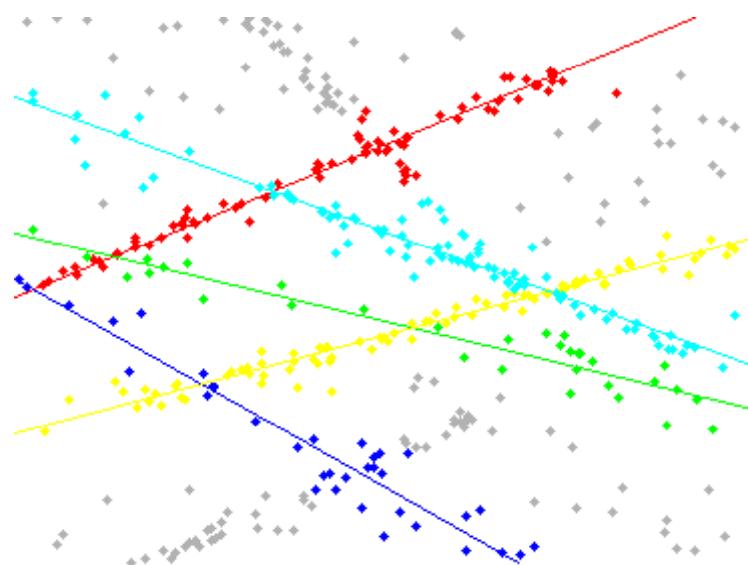
Does not happen in vision

# K-means vs. PEARL

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$$E(f) = \sum_p \| p, f_p \| + \text{hard constraint on number of models}$$

K-means



5 random initial lines + outlier model

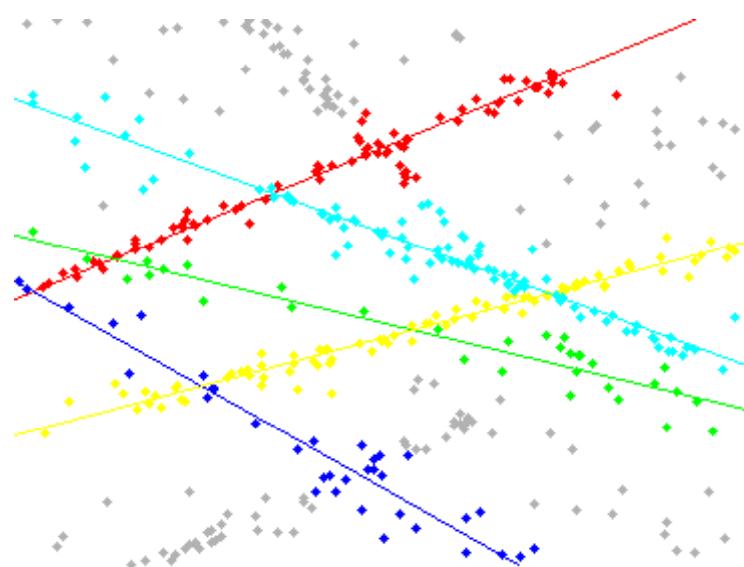
gets stuck in local minima

# K-means vs. PEARL

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$$E(f) = \sum_p \| p, f_p \| + \sum_{f \in \Lambda} h_f \cdot \delta_f(f)$$

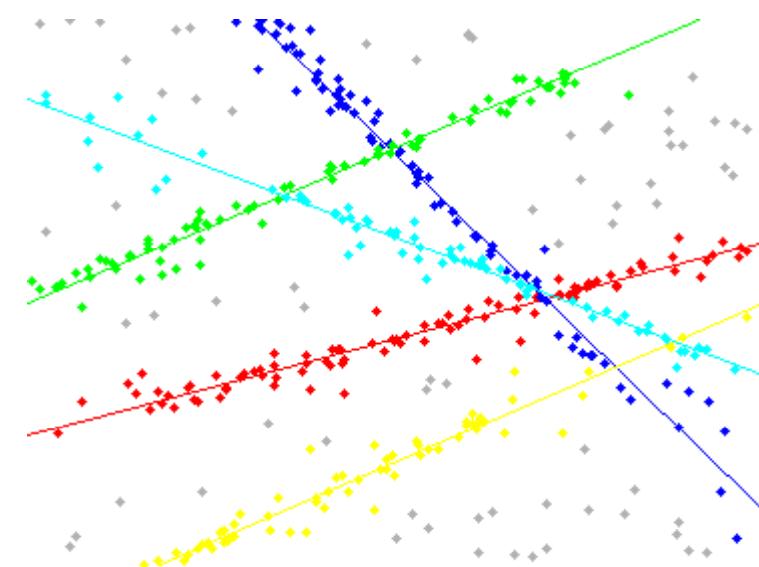
K-means



5 random initial lines + outlier model

gets stuck in local minima

Our approach  $h_f = 1000$



1000 initial lines + outlier model

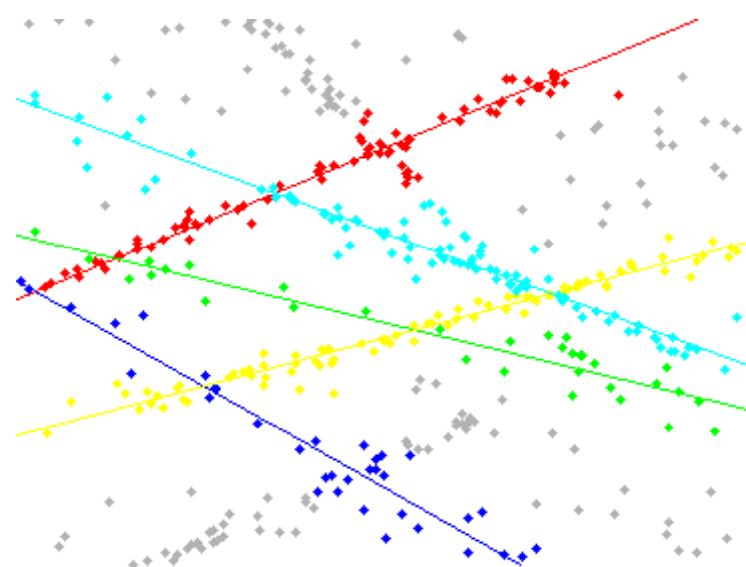
better explores label space

# K-means vs. PEARL

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$$E(f) = \sum_p \| p, f_p \| + \sum_{f \in \Lambda} h_f \cdot \delta_f(f)$$

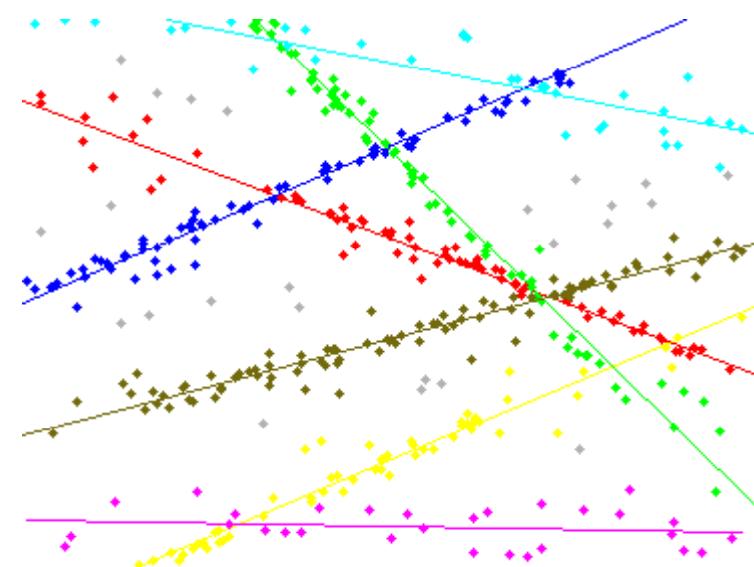
K-means



5 random initial lines + outlier model

gets stuck in local minima

Our approach  $h_f = 500$



1000 initial lines + outlier model

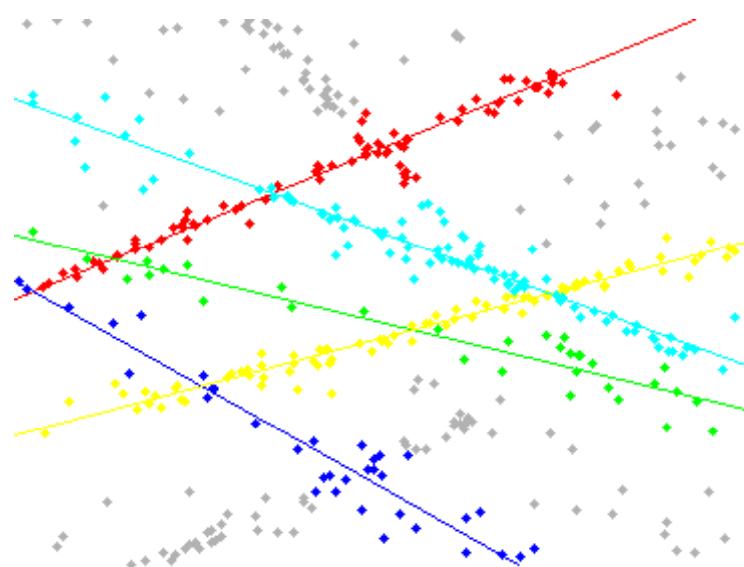
better explores label space

# K-means vs. PEARL

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$$E(f) = \sum_p \| p, f_p \| + \sum_{f \in \Lambda} h_f \cdot \delta_f(f)$$

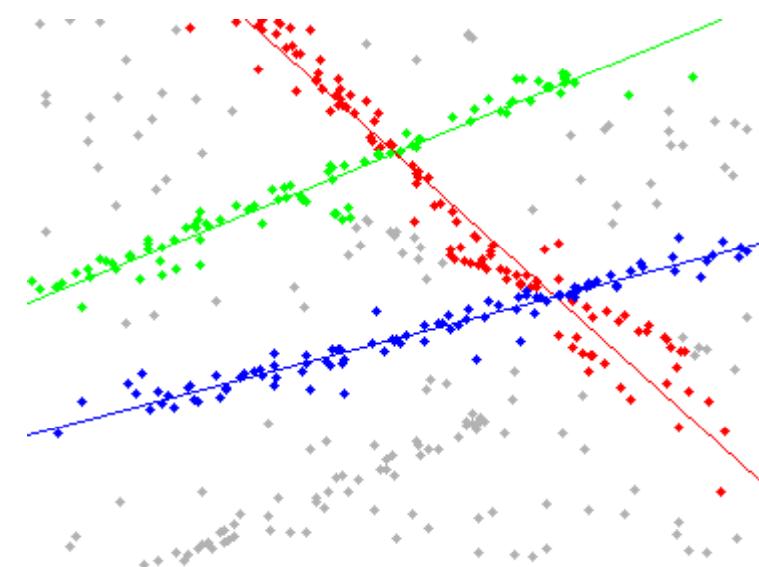
K-means



5 random initial lines + outlier model

**gets stuck in local minima**

Our approach  $h_f = 2000$



1000 initial lines + outlier model

**better explores label space**