Image Segmentation

Philipp Krähenbühl

Stanford University

April 24, 2013

Philipp Krähenbühl (Stanford University)

Segmentation

April 24, 2013 1 / 63

3

< ロ > < 同 > < 三 > < 三

• Goal: identify groups of pixels that go together



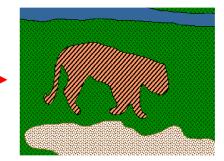


Image: A = 1 = 1

Success Story



Philipp Krähenbühl (Stanford University)

Segmentation

Success Story



Success Story



Philipp Krähenbühl (Stanford University)

Segmentation

Gestalt Theory

- Gestalt: whole or group
 - The whole is greater than the sum of its parts
 - Relationships between parts can yield new properties/features
- Psychologists identified series of factors that predispose set of elements to be grouped (by human visual system)

Max Wertheimer (1880-1943)

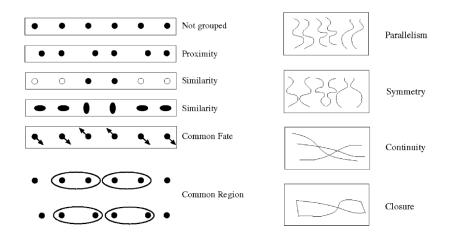
I stand at the window and see a house, trees, sky. Theoretically I might say there were 327 brightnesses and nuances of color. Do I have "327"? No. I have sky, house, and trees.

Untersuchungen zur Lehre von der Gestalt, Psychologische Forschung, Vol. 4, pp. 301-350, 1923 http://psy.ed.asu.edu/~classics/Wertheimer/Forms/ forms.htm





Gestalt Theory



 These factors make intuitive sense, but are very difficult to translate into algorithms.

Philipp Krähenbühl (Stanford University)

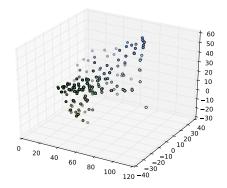
Segmentation

3

< ロ > < 同 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ >

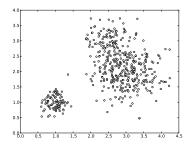
Segmentation as clustering





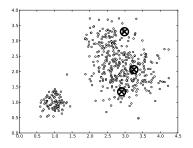
• Pixels are points in a high dimensional space

- color: 3d
- color+location:5d
- Cluster pixels into segment



- Randomly initialize K cluster centers, c₁,..., c_k
- ② Given cluster centers, determine points in each cluster
 - For each point p, find the closest c_i. Put p into cluster i.
- **(3)** Given points in each cluster, solve for c_i
 - Set c_i to be the mean of points in cluster i
- If c_i have changed, repeat Step 2

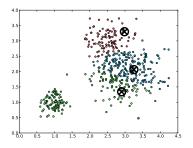
4 A N



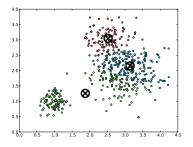
Randomly initialize K cluster centers, c₁,..., c_k

② Given cluster centers, determine points in each cluster

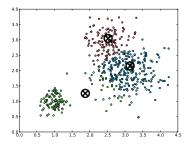
- For each point p, find the closest c_i. Put p into cluster i.
- Given points in each cluster, solve for c_i
 - Set c_i to be the mean of points in cluster i
- If c_i have changed, repeat Step 2



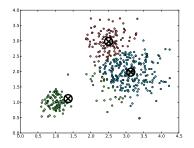
- Randomly initialize K cluster centers, c₁,..., c_k
- Q Given cluster centers, determine points in each cluster
 - ▶ For each point *p*, find the closest *c_i*. Put *p* into cluster *i*.
- Given points in each cluster, solve for c_i
 - Set c_i to be the mean of points in cluster i
- If c_i have changed, repeat Step 2



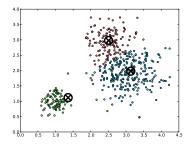
- Randomly initialize K cluster centers, c₁,..., c_k
- Q Given cluster centers, determine points in each cluster
 - ▶ For each point *p*, find the closest *c_i*. Put *p* into cluster *i*.
- Given points in each cluster, solve for c_i
 - Set c_i to be the mean of points in cluster i
- If c_i have changed, repeat Step 2



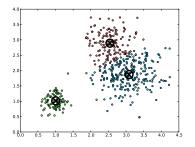
- Randomly initialize K cluster centers, c₁,..., c_k
- Q Given cluster centers, determine points in each cluster
 - ▶ For each point *p*, find the closest *c_i*. Put *p* into cluster *i*.
- Given points in each cluster, solve for c_i
 - Set c_i to be the mean of points in cluster i
- If c_i have changed, repeat Step 2



- Randomly initialize K cluster centers, c₁,..., c_k
- Q Given cluster centers, determine points in each cluster
 - ▶ For each point *p*, find the closest *c_i*. Put *p* into cluster *i*.
- Given points in each cluster, solve for c_i
 - Set c_i to be the mean of points in cluster i
- If c_i have changed, repeat Step 2



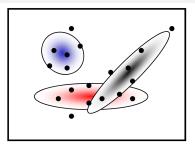
- Randomly initialize K cluster centers, c₁,..., c_k
- Q Given cluster centers, determine points in each cluster
 - ▶ For each point *p*, find the closest *c_i*. Put *p* into cluster *i*.
- Given points in each cluster, solve for c_i
 - Set c_i to be the mean of points in cluster i
- If c_i have changed, repeat Step 2



- Randomly initialize K cluster centers, c₁,..., c_k
- Q Given cluster centers, determine points in each cluster
 - ▶ For each point *p*, find the closest *c_i*. Put *p* into cluster *i*.
- Given points in each cluster, solve for c_i
 - Set c_i to be the mean of points in cluster i
- If c_i have changed, repeat Step 2



Expectation Maximization (EM)



Goal

Find blob parameters θ that maximize the likelihood function:

$$\mathsf{P}(\mathsf{data}| heta) = \prod_{\mathsf{x}} \mathsf{P}(\mathsf{x}| heta)$$

• Approach:

- E-step: given current guess of blobs, compute ownership of each point
- Omega M-step: given ownership probabilities, update blobs to maximize likelihood function
- 8 Repeat until convergence

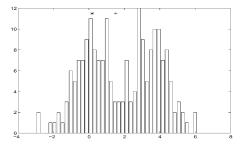
Expectation Maximization (EM)



∃ →

Image: A match a ma

Mean-Shift Algorithm



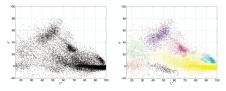
Iterative Mode search

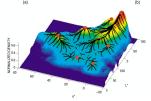
- **1** Initialize random seed, and window W
- **2** Calculate center of gravity (the "mean") of $W: \sum_{x \in W} xH(x)$
- Shift the search window to the mean
- Repeat Step 2 until convergence

Mean-Shift Segmentation

Iterative Mode search

- Find features (color, gradients, texture, etc)
- Initialize windows at individual pixel locations
- Perform mean shift for each window until convergence
- Merge windows that end up near the same "peak" or mode





Expectation Maximization (EM)



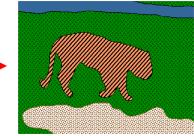
- ∢ ≣ →

Image: A math a math

Back to Image Segmentation

• Goal: identify groups of pixels that go together



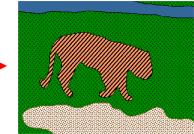


- Up to now, we have focused on ways to group pixels into image segments based on their appearance...
 - Segmentation as clustering.
- We also want to enforce region constraints.
 - Spatial consistency
 - Smooth borders

Back to Image Segmentation

• Goal: identify groups of pixels that go together



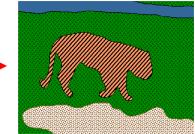


- Up to now, we have focused on ways to group pixels into image segments based on their appearance...
 - Segmentation as clustering.
- We also want to enforce region constraints.
 - Spatial consistency
 - Smooth borders

Back to Image Segmentation

• Goal: identify groups of pixels that go together





- Up to now, we have focused on ways to group pixels into image segments based on their appearance...
 - Segmentation as clustering.
- We also want to enforce region constraints.
 - Spatial consistency
 - Smooth borders

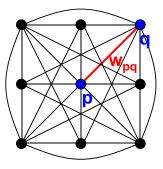
- Graph theoretic segmentation
 - Normalized Cuts
 - Using texture features
- Segmentation as Energy Minimization
 - Markov Random Fields (MRF) / Conditional Random Fields (CRF)
 - Graph cuts for image segmentation
 - Applications

What we will learn today?

Graph theoretic segmentation

- Normalized Cuts
- Using texture features
- Segmentation as Energy Minimization
 - Markov Random Fields (MRF) / Conditional Random Fields (CRF)
 - Graph cuts for image segmentation
 - Applications

Images as Graphs





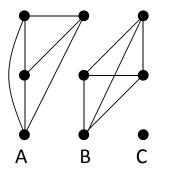
- (Fully-Connected) Graph
 - Node (vertex) for every pixel
 - Link between (every) pair of pixels, (p,q)
 - Affinity weight w_{pq} for each link (edge)
 - \star w_{pq} measures similarity
 - * Inverse proportional to distance (difference in color and position)

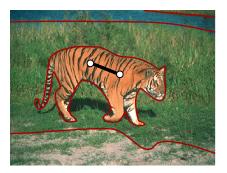
Slide Credit: Steve Seitz

Philipp Krähenbühl (Stanford University)

April 24, 2013 16 / 63

Segmentation by Graph Cuts





A = A = A = A = A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A
 A

• Break Graph into Segments (cliques)

- Delete links that cross between segments
- Easiest to break links that low similarity (low affinity weight)
 - * Similar pixels should be in the same segment
 - ★ Dissimilar pixels should be if different segments

Slide Credit: Steve Seitz

• Distance $\exp(-\frac{1}{2\sigma^2}\|x-y\|^2)$

Intensity

$$\exp(-\frac{1}{2\sigma^2} \|I(x) - I(y)\|^2)$$

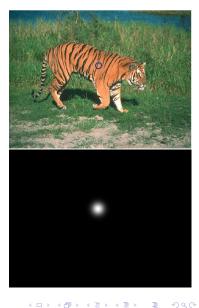
• Color

$$\exp(-\frac{1}{2\sigma^2}\underbrace{dist(c(x),c(y))^2}_{\text{suitable color distance}})$$

Texture

$$\exp(-\frac{1}{2\sigma^2} \| \underbrace{f(x) - f(y)}_{\text{Filter output}} \|^2)$$

Source: Forsyth & Ponce



• Distance
$$\exp(-\frac{1}{2\sigma^2}\|x-y\|^2)$$

Intensity

$$\exp(-\frac{1}{2\sigma^2}\|I(x) - I(y)\|^2)$$

Color

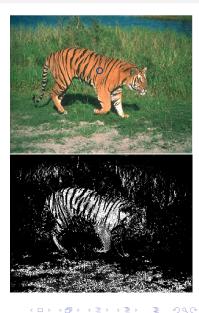
$$\exp(-\frac{1}{2\sigma^2}\underbrace{dist(c(x),c(y))^2}_{\text{suitable color distance}})$$

suitable color distance

Texture

$$\exp(-\frac{1}{2\sigma^2} \| \underbrace{f(x) - f(y)}_{\text{Filter output}} \|^2)$$

Source: Forsyth & Ponce



• Distance
$$\exp(-\frac{1}{2\sigma^2}\|x-y\|^2)$$

Intensity

۵

$$\exp(-\frac{1}{2\sigma^2}\|I(x) - I(y)\|^2)$$

Color

$$exp(-\frac{1}{2\sigma^2} \underbrace{dist(c(x), c(y))^2}_{\text{suitable color distance}}$$

$$\exp(-\frac{1}{2\sigma^2} \|\underbrace{f(x) - f(y)}_{\text{Filter output}}\|^2)$$

Source: Forsyth & Ponce

イロト イポト イヨト イヨト

3

• Distance
$$\exp(-\frac{1}{2\sigma^2}\|x-y\|^2)$$

Intensity

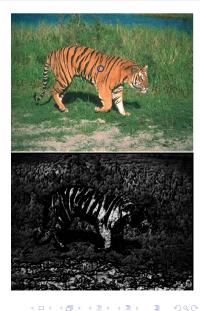
$$\exp(-\frac{1}{2\sigma^2}\|I(x) - I(y)\|^2)$$

$$\exp(-\frac{1}{2\sigma^2}\underbrace{dist(c(x), c(y))^2}_{\text{suitable color distance}})$$

Texture

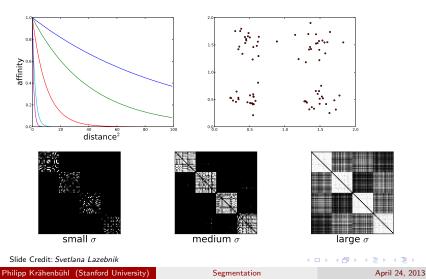
$$\exp(-\frac{1}{2\sigma^2} \|\underbrace{f(x) - f(y)}_{\text{Filter output}}\|^2)$$

Source: Forsyth & Ponce



Scale Affects Affinity

- Small σ : group only nearby points
- Large σ : group far-away points



19 / 63

Graph Cut: Using Eigenvalues

• Affinity matrix W

• Extract a single good cluster (v_n)

 \triangleright $v_n(i)$: probability of point *i* belonging to the cluster

Elements have high affinity with each other

$$v_n^{\perp} W v_n$$

Constraint v_n^T v_n = 1
 Prevents v_n → constraints

• Constraint objective

$$v_n^{\top} W v_n - \lambda (1 - v_n^{\top} v_n)$$



• Reduces to Eigenvalue problem

$$v_n^\top W = \lambda v_n$$

April 24, 2013 20 / 63

イロト イヨト イヨト イヨト

Graph Cut: Using Eigenvalues

- Affinity matrix W
- Extract a single good cluster (v_n)
 - $v_n(i)$: probability of point *i* belonging to the cluster
 - Elements have high affinity with each other

$$v_n^\top W v_n$$

- Constraint $v_n^\top v_n = 1$ • Prevents $v_n \to \infty$
- Constraint objective

$$v_n^\top W v_n - \lambda (1 - v_n^\top v_n)$$



• Reduces to Eigenvalue problem

$$v_n^\top W = \lambda v_n$$

- Affinity matrix W
- Extract a single good cluster (v_n)
 - $v_n(i)$: probability of point *i* belonging to the cluster
 - Elements have high affinity with each other

$$v_n^\top W v_n$$

• Constraint
$$v_n^\top v_n = 1$$

• Prevents $v_n \to \infty$

• Constraint objective

$$v_n^{\top} W v_n - \lambda (1 - v_n^{\top} v_n)$$



• Reduces to Eigenvalue problem

$$v_n^\top W = \lambda v_n$$

April 24, 2013 20 / 63

• • • • • • • • • • • •

- Affinity matrix W
- Extract a single good cluster (v_n)
 - $v_n(i)$: probability of point *i* belonging to the cluster
 - Elements have high affinity with each other

$$v_n^+ W v_n$$

• Constraint objective

$$v_n^\top W v_n - \lambda (1 - v_n^\top v_n)$$



• Reduces to Eigenvalue problem

$$v_n^\top W = \lambda v_n$$

April 24, 2013 20 / 63

< ロ > < 同 > < 三 > < 三

- Affinity matrix W
- Extract a single good cluster (v_n)
 - v_n(i): probability of point i belonging to the cluster
 - Elements have high affinity with each other

$$v_n^\top W v_n$$

• Constraint
$$v_n^\top v_n = 1$$

★ Prevents $v_n \rightarrow \infty$

• Constraint objective

$$v_n^{\top} W v_n - \lambda (1 - v_n^{\top} v_n)$$



• Reduces to Eigenvalue problem

$$v_n^\top W = \lambda v_n$$

April 24, 2013 20 / 63

(日) (同) (三) (三)

- Affinity matrix W
- Extract a single good cluster (v_n)
 - v_n(i): probability of point i belonging to the cluster
 - Elements have high affinity with each other

$$v_n^\top W v_n$$

• Constraint objective

$$v_n^{\top} W v_n - \lambda (1 - v_n^{\top} v_n)$$



• Reduces to Eigenvalue problem

$$v_n^\top W = \lambda v_n$$

April 24, 2013 20 / 63

(日) (同) (三) (三)

- Affinity matrix W
- Extract a single good cluster (v_n)
 - v_n(i): probability of point i belonging to the cluster
 - Elements have high affinity with each other

$$v_n^\top W v_n$$

• Constraint
$$v_n^\top v_n = 1$$

★ Prevents
$$v_n \rightarrow \infty$$

Constraint objective

$$v_n^\top W v_n - \lambda (1 - v_n^\top v_n)$$



• Reduces to Eigenvalue problem

$$v_n^\top W = \lambda v_n$$

April 24, 2013 20 / 63

< ロ > < 同 > < 三 > < 三

- Affinity matrix W
- Extract a single good cluster (v_n)
 - $v_n(i)$: probability of point *i* belonging to the cluster
 - Elements have high affinity with each other

$$v_n^\top W v_n$$

• Constraint
$$v_n^\top v_n = 1$$

★ Prevents
$$v_n \to \infty$$

Constraint objective

$$v_n^\top W v_n - \lambda (1 - v_n^\top v_n)$$

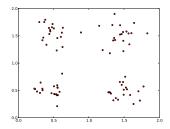


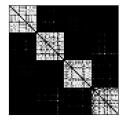
• Reduces to Eigenvalue problem

$$v_n^\top W = \lambda v_n$$

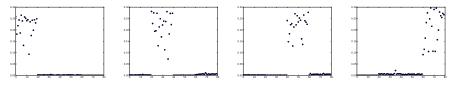
April 24, 2013 20 / 63

・ロト ・ 同ト ・ ヨト ・ ヨ





4 largest eigenvalues



Segmentation

April 24, 2013 21 / 63

Construct an affinity matrix

- Ompute the eigenvalues and eigenvectors of the affinity matrix
- Ontil there are sufficient clusters
 - Take the eigenvector corresponding to the largest unprocessed eigenvalue
 - zero all components corresponding to elements that have already been clustered
 - threshold the remaining components to determine which element belongs to this cluster,
 - A choose a threshold by clustering the components, or using a threshold fixed in advance.
 - If all elements have been accounted for, there are sufficient clusters: end

< ロ > < 同 > < 三 > < 三

Construct an affinity matrix

- Output the eigenvalues and eigenvectors of the affinity matrix
- Ontil there are sufficient clusters
 - Take the eigenvector corresponding to the largest unprocessed eigenvalue
 - zero all components corresponding to elements that have already been clustered
 - threshold the remaining components to determine which element belongs to this cluster,
 - choose a threshold by clustering the components, or using a threshold freed in advance.
 - If all elements have been accounted for, there are sufficient clusters: end

< ロ > < 同 > < 三 > < 三

- Construct an affinity matrix
- Ompute the eigenvalues and eigenvectors of the affinity matrix
- Ontil there are sufficient clusters
 - ▶ Take the eigenvector corresponding to the largest unprocessed eigenvalue
 - zero all components corresponding to elements that have already been clustered
 - threshold the remaining components to determine which element belongs to this cluster,
 - * choose a threshold by clustering the components, or using a threshold fixed in advance.
 - ▶ If all elements have been accounted for, there are sufficient clusters: end

- Construct an affinity matrix
- Ompute the eigenvalues and eigenvectors of the affinity matrix
- Ontil there are sufficient clusters
 - ► Take the eigenvector corresponding to the largest unprocessed eigenvalue
 - zero all components corresponding to elements that have already been clustered
 - threshold the remaining components to determine which element belongs to this cluster,
 - * choose a threshold by clustering the components, or using a threshold fixed in advance.
 - ▶ If all elements have been accounted for, there are sufficient clusters: end

- Construct an affinity matrix
- Ompute the eigenvalues and eigenvectors of the affinity matrix
- Ontil there are sufficient clusters
 - ► Take the eigenvector corresponding to the largest unprocessed eigenvalue
 - zero all components corresponding to elements that have already been clustered
 - threshold the remaining components to determine which element belongs to this cluster,
 - choose a threshold by clustering the components, or using a threshold fixed in advance.
 - ▶ If all elements have been accounted for, there are sufficient clusters: end

- Construct an affinity matrix
- Ompute the eigenvalues and eigenvectors of the affinity matrix
- Ontil there are sufficient clusters
 - ► Take the eigenvector corresponding to the largest unprocessed eigenvalue
 - zero all components corresponding to elements that have already been clustered
 - threshold the remaining components to determine which element belongs to this cluster,
 - * choose a threshold by clustering the components, or using a threshold fixed in advance.
 - ▶ If all elements have been accounted for, there are sufficient clusters: end

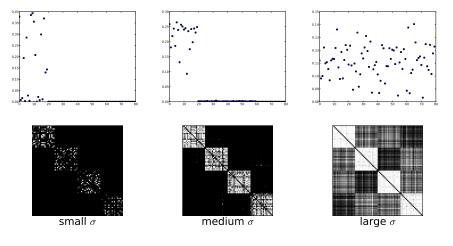
- Construct an affinity matrix
- Ompute the eigenvalues and eigenvectors of the affinity matrix
- Ontil there are sufficient clusters
 - Take the eigenvector corresponding to the largest unprocessed eigenvalue
 - zero all components corresponding to elements that have already been clustered
 - threshold the remaining components to determine which element belongs to this cluster,
 - choose a threshold by clustering the components, or using a threshold fixed in advance.
 - ▶ If all elements have been accounted for, there are sufficient clusters: end

(日) (周) (三) (三)

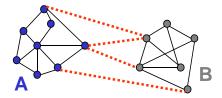
- Construct an affinity matrix
- Ompute the eigenvalues and eigenvectors of the affinity matrix
- Ontil there are sufficient clusters
 - Take the eigenvector corresponding to the largest unprocessed eigenvalue
 - zero all components corresponding to elements that have already been clustered
 - threshold the remaining components to determine which element belongs to this cluster,
 - choose a threshold by clustering the components, or using a threshold fixed in advance.
 - ▶ If all elements have been accounted for, there are sufficient clusters: end

(日) (周) (三) (三)

Effects of the scaling



Graph Cut

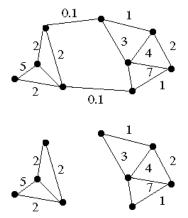


- Find set of edges whose removal makes graph disconnected
- Cost of a cut
 - Sum of weights of cut edges: $cut(A, B) = \sum_{p \in A, q \in B} w_{pq}$
- Graph cut gives us a segmentation
 - What is a "good" graph cut and how do we find one?

Slide Credit: Steve Seitz

イロト 不得下 イヨト イヨト

Graph Cut





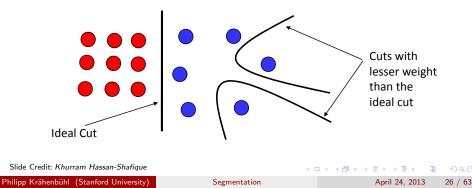
Here, the cut is nicely defined by the block-diagonal structure of the affinity matrix.

⇒ How can this be generalized?

Image Source: Forsyth & Ponce

Minimum Cut

- We can do segmentation by finding the minimum cut in a graph
 - a minimum cut of a graph is a cut whose cutset has the smallest affinity.
 - Efficient algorithms exist for doing this (max-flow)
- Drawback
 - Weight of cut proportional to number of edges in the cut
 - Minimum cut tends to cut off very small, isolated components



Normalized Cut (NCut)

- A minimum cut penalizes large segments
- This can be fixed by normalizing for size of segments
- The normalized cut cost is:

$$Ncut(A, B) = \frac{cut(A, B)}{assoc(A, V)} + \frac{cut(A, B)}{assoc(B, V)}$$
$$= cut(A, B) \left[\frac{1}{\sum_{\rho \in A, q} w_{\rho, q}} + \frac{1}{\sum_{q \in B, \rho} w_{\rho, q}} \right]$$

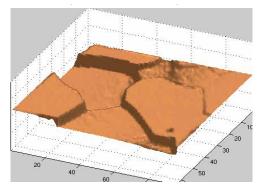
• assoc(A, V) = sum of weights of all edges in V that touch A

- The exact solution is NP-hard but an approximation can be computed by solving a generalized eigenvalue problem.
- J. Shi and J. Malik. Normalized cuts and image segmentation. PAMI 2000

▲□▶ ▲□▶ ▲□▶ ▲□▶ = ののの

Interpretation as a Dynamical System





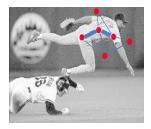
• Treat the links as springs and shake the system

- Elasticity proportional to cost
- Vibration "modes" correspond to segments
 - $\star\,$ Can compute these by solving a generalized eigenvector problem

Slide Credit: Steve Seitz

- Definitions
 - ► W: the affinity matrix
 - D: diagonal matrix, $D_{ii} = \sum_{j} W_{ij}$
 - x: a vector in $\{-1, 1\}^N$,
- Rewriting the Normalized Cut in matrix form

$$Ncut(A, B) = \frac{cut(A, B)}{assoc(A, V)} + \frac{cut(A, B)}{assoc(B, V)}$$
$$= \dots$$



Slide Credit: Jitentra Malik

イロト イヨト イヨト イヨト

Some more math...

We see again this is an unbiased measure, which reflects how tightly on average nodes within the group are connected to each other.

Another important property of this definition of association and disassociation of a partition is that they are naturally related:

$$\begin{split} \operatorname{Noak}(A,B) &= \frac{\operatorname{out}(A,B)}{\operatorname{out}(A,V)} + \frac{\operatorname{out}(A,B)}{\operatorname{out}(A,V)} \\ &= \frac{\operatorname{aux}(A,V) - \operatorname{aux}(A,A)}{\operatorname{out}(A,V) - \operatorname{aux}(A,A)} \\ &+ \frac{\operatorname{sum}(A,V) - \operatorname{aux}(A,B)}{\operatorname{out}(A,K)} \\ &= 2 - (\frac{\operatorname{aux}(A,A)}{\operatorname{out}(A,A)} + \frac{\operatorname{aux}(A,B)}{\operatorname{out}(A,A)} + \frac{\operatorname{aux}(A,B)}{\operatorname{out}(A,A)} \end{split}$$

= 2 - Namo(A, B)

Hence the two partition criteria that we seek in our grouping algorithm, minimizing the disassectiation between the groups and maximizing the association within the group, are in fact identical, and can be satished annukaneously. In our algorithm, we will use this associated out as the partition criterion.

Having defined the graph partition oriterion that we want to optimize, we will show how such an optimal partition can be computed efficiently.

2.1. Computing the optimal partition Given a particle of nodes of a graph, V, into two sets A and B, let a be an N = |V| dimensional indicator vector, $z_i = 1$ if node is k in A, and -1 otherwise. Let $d(I) = \sum_{i=1}^{N} u(i,j)$, he the total connection from nodes it call other nodes. With the definitions a and are can service N cuis(A, B) as:

$$\begin{split} \mathsf{N}\mathsf{cut}(A,B) &= \frac{\mathsf{cut}(A,B)}{\mathsf{cus}(A,V)} + \frac{\mathsf{cut}(B,A)}{\mathsf{aux}(B,V)} \\ &= \frac{\sum_{\{\underline{a}>0,\underline{a}_i<\underline{a}\}} - \mathsf{su}_{ij}|\underline{a}_i|\underline{a}_i}{\sum_{\underline{a}>0} \cdot \underline{a}_i} \\ &+ \frac{\sum_{\{\underline{a}>0,\underline{a}_i<\underline{a}\}} - \mathsf{su}_{ij}|\underline{a}_i|\underline{a}_i}{\sum_{\underline{a}>0} \cdot \underline{a}_i} d_i \end{split}$$

Let D be an $N \times N$ diagonal matrix with d to its diagonal, W be an $N \times N$ dynametrical matrix with $W(i_3) = w_{i_3} k = \frac{\sum_{i > 1} d_i}{\sum_{i < 1} d_i} a_i \text{ at be an } N \times 1$ vector of all cose. Using the fact $\frac{1}{2} \frac{d_i}{d_i}$ and $\frac{1}{2} \frac{d_i}{d_i}$ mindes are write $(N \otimes (a_i) \text{ at } k = 1, 2)$ of any exciton i_2 , $j \in \mathbb{N}$ and $i_2 < 0$ supertively, we can rewrite $(N \otimes (a_i) \text{ at } k)$

$$= \frac{(\mathbf{z}+\mathbf{z})^{\mathbf{v}}(\mathbf{D}-\mathbf{W})(\mathbf{z}+\mathbf{z})}{\mathbf{z}^{\mathbf{v}}\mathbf{D}\mathbf{z}} + \frac{(\mathbf{z}-\mathbf{z})^{\mathbf{v}}(\mathbf{D}-\mathbf{W})(\mathbf{z}-\mathbf{z})}{(1-\mathbf{z})^{\mathbf{v}}\mathbf{D}\mathbf{z}}$$

$$= \frac{(\mathbf{z}^{\mathbf{v}}(\mathbf{D}-\mathbf{W})\mathbf{z}_{\mathbf{z}}+\mathbf{z}^{\mathbf{v}}(\mathbf{D}-\mathbf{W})\mathbf{z})}{\mathbf{z}(1-\mathbf{z})^{\mathbf{z}}(\mathbf{D}-\mathbf{W})\mathbf{z}} + \frac{\mathbf{z}(1-\mathbf{z})\mathbf{z}^{\mathbf{v}}(\mathbf{D}-\mathbf{W})\mathbf{z}}{\mathbf{z}(1-\mathbf{z})^{\mathbf{z}}\mathbf{T}}\mathbf{D}\mathbf{z}}$$

Let
$$\alpha(\mathbf{a}) = \mathbf{a}^T (\mathbf{D} - \mathbf{W}) \mathbf{a}$$
, $\beta(\mathbf{a}) = \mathbf{x}^T (\mathbf{D} - \mathbf{W}) \mathbf{a}$, $\gamma = \mathbf{x}^T (\mathbf{D} - \mathbf{W}) \mathbf{x}$, and $M = \mathbf{x}^T \mathbf{D} \mathbf{x}$, we can then further expand the above equation as:

$$\frac{(\alpha(\mathbf{a}) + \gamma) + 2(1 - 2k)\beta(\mathbf{a})}{k(1 - k)M}$$

$$= \frac{(\alpha(\mathbf{a}) + \gamma) + 2(1 - 2k)\beta(\mathbf{a})}{k(1 - k)M} - \frac{2(\alpha(\mathbf{a}) + \gamma)}{M}$$

$$= \frac{\gamma(\alpha(\mathbf{a}))}{M} + \frac{\gamma(\mathbf{a})}{M} + \frac{\gamma(\mathbf{a})}{M}$$

dropping the last constant term, which in this case equals 0, we get

$$= \frac{\frac{(1-2k+2k^2)(\alpha(\mathbf{a})+\gamma)+2(1-2k)\beta(\mathbf{a})}{k(1-k)M} + \frac{2\alpha(\mathbf{a})}{M}}{\sum_{\substack{(1-2k+2k^2)\\(1-k)^2}}(\alpha(\mathbf{a})+\gamma) + \frac{2(-2k)}{(1-k)^2}\beta(\mathbf{a})} + \frac{2\alpha(\mathbf{a})}{M}$$

Letting $\delta = \frac{\delta}{1-\delta}$, and since $\gamma = 0$, it becomes,

$$\begin{array}{l} \displaystyle = \frac{(1+\delta^2)(z(u)+\gamma+2(1-\delta^2)(u)}{\delta M} + \frac{2\partial (z(u)}{\delta M} \\ \displaystyle = \frac{(1+\delta^2)(z(u)+\gamma)}{\delta M} + \frac{2\partial (z(u)}{\delta M} - \frac{2\partial (z(u)$$

Setting $y = (x + a) - \delta(x - a)$, it is easy to see that

$${}^{T}D\mathbf{x} = \sum_{\mathbf{x},j=0} \mathbf{d}_{i} - b \sum_{\mathbf{x},i\neq 0} \mathbf{d}_{i} = 0 \quad (4)$$

$$\begin{aligned} & \operatorname{ce} \, b = \frac{s}{1-\kappa} = \frac{\sum_{i=\infty}^{1-\mu} d_i^{i}}{\sum_{i=\infty}^{1} d_i^{i}}, \, \operatorname{and} \\ & y^T Dy = \sum_{m \geq 0} d_i + b^T \sum_{m < 0} d_i \\ & = \quad b \sum_{n < 0} d_i + b \sum_{m < 0} d_i \\ & = \quad b (\sum_{m < 0} d_i + b \sum_{m < 0} d_i) \\ & = \quad b n^T D_i. \end{aligned}$$

Slide Credit: Jitentra Malik

Philipp Krähenbühl (Stanford University)

Segmentation

410

April 24, 2013 30 / 63

3

< ロ > < 同 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ >

• After simplifications, we get

$$Ncut(A,B) = \frac{y^{\top}(D-W)y}{y^{\top}Dy}$$

with $y_i \in \{-1, b\}$ and $y^{\top} D1 = 0$

- This is the Rayleigh Quotient
 - Solution given by the generalized eigenvalue problem

 $(D - W)y = \lambda Dy$

♥
Continuous

approximation

Subtleties

- Optimal solution is second smallest eigenvector
- Gives continuous result—must convert into discrete values of y

Slide Credit: Jitentra Malik Philipp Krähenbühl (Stanford University)

April 24, 2013 31 / 63

(日) (同) (三) (三)

• After simplifications, we get

$$Ncut(A, B) = \frac{y^{\top}(D - W)y}{y^{\top}Dy}$$

with
$$y_i \in \{-1, b\}$$
 and $y^{ op} D1 = 0$

• This is the Rayleigh Quotient

 Solution given by the generalized eigenvalue problem

$$(D - W)y = \lambda Dy$$

Hard as a discrete problem

∜

Continuous approximation

Subtleties

- Optimal solution is second smallest eigenvector
- Gives continuous result—must convert into discrete values of y

Slide Credit: Jitentra Malik Philipp Krähenbühl (Stanford University)

April 24, 2013 31 / 63

- **(())) (())) ())**

After simplifications, we get

$$Ncut(A, B) = \frac{y^{\top}(D - W)y}{y^{\top}Dy}$$

Hard as a discrete problem

 \downarrow

with
$$y_i \in \{-1, b\}$$
 and $y^{\top}D1 = 0$

- This is the Rayleigh Quotient
 - Solution given by the generalized eigenvalue problem

$$(D - W)y = \lambda Dy$$

Continuous approximation

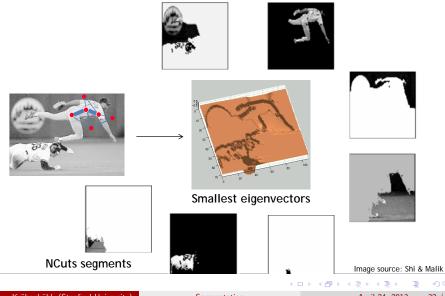
Subtleties

- Optimal solution is second smallest eigenvector
- Gives continuous result—must convert into discrete values of y

Slide Credit: litentra Malik Philipp Krähenbühl (Stanford University)

April 24, 2013 31 / 63

NCuts Example



Philipp Krähenbühl (Stanford University)

Segmentation

April 24, 2013 32 / 63

NCuts Example

- Problem: eigenvectors take on continuous values
 - How to choose the splitting point to binarize the image?



- Possible procedures
 - Pick a constant value (0, or 0.5).
 - Pick the median value as splitting point.
 - Look for the splitting point that has the minimum NCut value:
 - Choose n possible splitting points.
 - 2 Compute NCut value.
 - Ick minimum.

• Construct a weighted graph G = (V, E) from an image.

- Onnect each pair of pixels, and assign graph edge weights W_{ij} = Prob. that i and j belong to the same region.
- 3 Solve $(D W)y = \lambda Dy$ for the smallest few eigenvectors. This yields a continuous solution.
- Threshold eigenvectors to get a discrete cut
 - This is where the approximation is made (we're not solving NP).
- In the security of the secu

NCuts Matlab code available at http://www.cis.upenn.edu/~jshi/software/

- Construct a weighted graph G = (V, E) from an image.
- ② Connect each pair of pixels, and assign graph edge weights $W_{ij} =$ Prob. that i and j belong to the same region.
- 3 Solve $(D W)y = \lambda Dy$ for the smallest few eigenvectors. This yields a continuous solution.
- Threshold eigenvectors to get a discrete cut
 This is where the approximation is made (we're not solving NP)
- In the security of the secu

NCuts Matlab code available at http://www.cis.upenn.edu/~jshi/software/

- Construct a weighted graph G = (V, E) from an image.
- Onnect each pair of pixels, and assign graph edge weights W_{ij} = Prob. that i and j belong to the same region.
- Solve $(D W)y = \lambda Dy$ for the smallest few eigenvectors. This yields a continuous solution.
- Threshold eigenvectors to get a discrete cut
 This is where the approximation is made (we're not solving NP).
- In the second second

NCuts Matlab code available at http://www.cis.upenn.edu/~jshi/software/

- Construct a weighted graph G = (V, E) from an image.
- Onnect each pair of pixels, and assign graph edge weights W_{ij} = Prob. that i and j belong to the same region.
- Solve $(D W)y = \lambda Dy$ for the smallest few eigenvectors. This yields a continuous solution.
- Threshold eigenvectors to get a discrete cut
 - This is where the approximation is made (we're not solving NP).
- Secursively subdivide if NCut value is below a pre-specified value.

NCuts Matlab code available at http://www.cis.upenn.edu/~jshi/software/

- Construct a weighted graph G = (V, E) from an image.
- ② Connect each pair of pixels, and assign graph edge weights $W_{ij} =$ Prob. that i and j belong to the same region.
- Solve $(D W)y = \lambda Dy$ for the smallest few eigenvectors. This yields a continuous solution.
- Threshold eigenvectors to get a discrete cut
 - This is where the approximation is made (we're not solving NP).

Recursively subdivide if NCut value is below a pre-specified value.

NCuts Matlab code available at http://www.cis.upenn.edu/~jshi/software/

- Construct a weighted graph G = (V, E) from an image.
- Onnect each pair of pixels, and assign graph edge weights W_{ij} = Prob. that i and j belong to the same region.
- Solve $(D W)y = \lambda Dy$ for the smallest few eigenvectors. This yields a continuous solution.
- Threshold eigenvectors to get a discrete cut
 - This is where the approximation is made (we're not solving NP).
- Secursively subdivide if NCut value is below a pre-specified value.

NCuts Matlab code available at http://www.cis.upenn.edu/~jshi/software/

・ロト ・四ト ・ヨト ・ヨト ・ヨ

NCuts Results







Image Source: Shi & Malik

Philipp Krähenbühl (Stanford University)

Segmentation

April 24, 2013 35 / 63

3

イロト イヨト イヨト イヨト

Using Texture Features for Segmentation

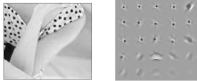
• Texture descriptor is vector of filter bank outputs



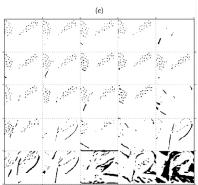
J. Malik, S. Belongie, T. Leung and J. Shi.

"Contour and Texture Analysis for Image Segmentation". IJCV 43(1),7-27,2001

Using Texture Features for Segmentation



- Texture descriptor is vector of filter bank outputs.
- Textons are found by clustering.
 - Bag of words



Slide Credit: Svetlana Lazebnik

Using Texture Features for Segmentation

- Texture descriptor is vector of filter bank outputs.
- Textons are found by clustering.
 - Bag of words
- Affinities are given by similarities of texton histograms over windows given by the "local scale" of the texture.





- **(())) (())) ())**

Results with Color and Texture



Segmentation

Summary: Normalized Cuts

- Pros:
 - Generic framework, flexible to choice of function that computes weights ("affinities") between nodes
 - Does not require any model of the data distribution

Cons:

- Time and memory complexity can be high

 - $\star\,$ Solving eigenvalue problem for each cut
- Preference for balanced partitions
 - If a region is uniform, NCuts will find the modes of vibration of the image dimensions



Slide Credit: Kristen Grauman

What we will learn today?

Graph theoretic segmentation

- Normalized Cuts
- Using texture features
- Segmentation as Energy Minimization
 - Markov Random Fields (MRF) / Conditional Random Fields (CRF)
 - Graph cuts for image segmentation
 - Applications

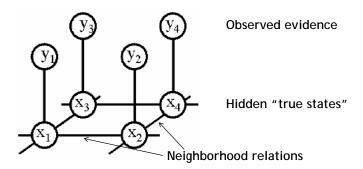
• Graph theoretic segmentation

- Normalized Cuts
- Using texture features
- Segmentation as Energy Minimization
 - Markov Random Fields (MRF) / Conditional Random Fields (CRF)
 - Graph cuts for image segmentation
 - Applications

< ロ > < 同 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ >

Markov Random Fields

- Allow rich probabilistic models for images
- But built in a local, modular way
 - Learn/model local effects, get global effects out



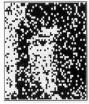
Slide Credit: William Freeman

Philipp Krähenbühl (Stanford University)

MRF Nodes as Pixels



Original image



Degraded image





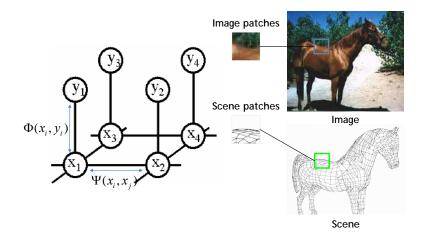
Reconstruction from MRF modeling pixel neighborhood statistics

(日) (同) (三) (三)

Slide Credit: Bastian Leibe

3

MRF Nodes as Patches

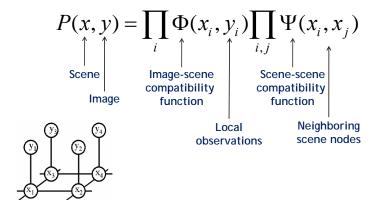


Slide Credit: William Freeman

3

イロト イヨト イヨト イヨト

Network Joint Probability



Slide Credit: William Freeman

Energy Formulation

Joint probability

$$P(x,y) = rac{1}{Z} \prod_i \Phi(x_i,y_i) \prod_{ij} \Psi(x_i,x_j)$$

Taking the log turns this into an Energy optimization

$$E(x,y) = \sum_{i} \varphi(x_i, y_i) + \sum_{ij} \psi(x_i, x_j)$$

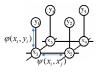
- This is similar to free-energy problems in statistical mechanics (spin glass theory). We therefore draw the analogy and call E an energy function.
- φ and ψ are called potentials.

Energy Formulation

Energy function

$$E(x,y) = \sum_{i} \underbrace{\varphi(x_{i}, y_{i})}_{\text{unary term}} + \sum_{ij} \underbrace{\psi(x_{i}, x_{j})}_{\text{pairwise term}}$$

- Unary potential φ
 - Encode local information about the given pixel/patch
 - How likely is a pixel/patch to belong to a certain class (e.g. foreground/background)?
- Pairwise potential ψ
 - Encode neighborhood information
 - How different is a pixel/patch's label from that of its neighbor? (e.g. based on intensity/color/texture difference, edges)



Slide Credit: Bastian Leibe

April 24, 2013 47 / 63

Segmentation using MRFs/CRFs

• Boykov and Jolly (2001)

$$E(x,y) = \sum_{i} \varphi(x_i, y_i) + \sum_{ij} \psi(x_i, x_j)$$

- Variables
 - x_i: Binary variable
 - \star foreground/background
 - y_i: Annotation
 - ★ foreground/background/empty
- Unary term
 - $\varphi(x_i, y_i) = K[x_i \neq y_i]$
 - Pay a penalty for disregarding the annotation
- Pairwise term
 - $\psi(x_i, x_j) = [x_i \neq x_j] w_{ij}$
 - Encourage smooth annotations
 - *w_{ij}* affinity between pixels *i* and *j*





Efficient solutions

- Grid structured random fields
 - Efficient solution using Maxflow/Mincut
 - Optimal solution for binary labeling
 - Boykov & Kolmogorov, "An Experimental Comparison of Min-Cut/Max-Flow Algorithms for Energy Minimization in Vision", PAMI 26(9): 1124-1137 (2004)
- Fully connected models
 - Efficient solution using convolution mean-field
 - Krähenbühl and Koltun, "Efficient Inference in Fully-Connected CRFs with Gaussian edge potentials", NIPS 2011



→ 3 → 4 3

GrabCut: Interactive Foreground Extraction





Slides credit: Carsten Rother

-

Image: A matrix

Segmentation

April 24, 2013 50 / 63

What GrabCut Does







Intelligent Scissors Mortensen and Barrett (1995)



GrabCut



Result



Regions



Boundary



Regions & Boundary

.∃ >

< 1[™] >

Philipp Krähenbühl (Stanford University)

Segmentation

April 24, 2013 51 / 63

GrabCut

Energy function

$$\mathsf{E}(\mathbf{x},\mathbf{k},\boldsymbol{\theta}|\mathbf{I}) = \sum_{i} \varphi(x_{i},k_{i},\boldsymbol{\theta}|z_{i}) + \sum_{ij} \psi(x_{i},x_{j}|z_{i},z_{j})$$

- Variables
 - $x_i \in \{0,1\}$: Foreground/background label
 - ▶ $k_i \in \{0, ..., K\}$: Gaussian mixture component
 - θ: Model parameters (GMM parameters)
 - $I = \{z_1, \ldots, z_N\}$: RGB Image
- Unary term $\varphi(x_i, k_i, \theta | z_i)$
 - Gaussian mixture model (log of a GMM)
- Pairwise term

$$\psi(x_i, x_j | z_i, z_j) = [x_i \neq x_j] \exp(-\beta \|z_i - z_j\|^2)$$

- 3

・ロン ・聞と ・ ほと ・ ほと

GrabCut



Philipp Krähenbühl (Stanford University)

Segmentation

April 24, 2013 53 / 63

GrabCut



Philipp Krähenbühl (Stanford University)

Segmentation

April 24, 2013 53 / 63

GrabCut - Unary term

• Gaussian Mixture Model

$$P(z_i|x_i, \theta) = \sum_k \pi(x_i, k) p(z_k|k, \theta)$$
• Hard to optimize (\sum_k)

Tractable solution

Assign each variable x_i a single mixture component k_i

$$P(z_i|x_i, k_i, \theta) = \pi(x_i, k_i)p(z_k|k_i, \theta)$$

- Optimize over k_i
- Unary term

$$\begin{split} \varphi(x_i, k_i, \boldsymbol{\theta} | z_i) &= -\log \pi(x_i, k_i) - \log p(z_k | k_i, \boldsymbol{\theta}) \\ &= -\log \pi(x_i, k_i) + \frac{1}{2} \log |\Sigma(k_i)| \\ &+ \frac{1}{2} (z_i - \mu(k_i))^\top \Sigma(k_i)^{-1} (z_i - \mu(k_i)) \end{split}$$

3

- 4 同 6 4 日 6 4 日 6

GrabCut - Unary term

• Unary term

$$egin{aligned} arphi(x_i,k_i,oldsymbol{ heta}|z_i) &= -\log\pi(x_i,k_i) + rac{1}{2}\log|\Sigma(k_i)| \ &+ rac{1}{2}(z_i-\mu(k_i))^{ op}\Sigma(k_i)^{-1}(z_i-\mu(k_i)) \end{aligned}$$

• Model parameters

$$\boldsymbol{\theta} = \{ \underbrace{\pi(x_i, k_i)}_{i \in \mathcal{N}}, \underbrace{\mu(k_i), \Sigma(k_i)}_{i \in \mathcal{N}} \}$$

mixture weight mean and variance

æ

< ロ > < 同 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < 回 > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ > < □ >

Initialize Mixture Models

Assign GMM components

$$k_i = \arg\min_k \varphi(x_i, k_i, \theta | z_i)$$

Icarn GMM parameters

$$\boldsymbol{\theta} = \arg\min\sum_{i} \varphi(x_i, k_i, \boldsymbol{\theta} | z_i)$$

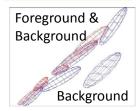
Estimate segmentation using mincut

 $\mathbf{x} = \arg\min E(\mathbf{x}, \mathbf{k}, \boldsymbol{\theta} | \mathbf{I})$

Repeat from 2 until convergence



Initialization



I ∃ ≥

April 24, 2013 56 / 63

- Initialize Mixture Models
- Assign GMM components

 $k_i = \arg\min_k \varphi(x_i, k_i, \theta|z_i)$

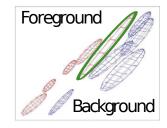
Icarn GMM parameters

$$\boldsymbol{\theta} = \arg\min\sum_{i} \varphi(x_i, k_i, \boldsymbol{\theta} | z_i)$$

Estimate segmentation using mincut

 $\mathbf{x} = \arg\min E(\mathbf{x}, \mathbf{k}, \boldsymbol{\theta} | \mathbf{I})$

Seperat from 2 until convergence



• • • • • • • • • • • •

- Initialize Mixture Models
- Assign GMM components

$$k_i = \arg\min_k \varphi(x_i, k_i, \theta|z_i)$$

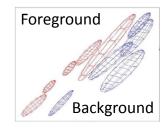
Searn GMM parameters

$$\boldsymbol{\theta} = \arg\min\sum_{i} \varphi(x_i, k_i, \boldsymbol{\theta} | z_i)$$

Istimate segmentation using mincut

 $\mathbf{x} = \arg\min E(\mathbf{x}, \mathbf{k}, \boldsymbol{\theta} | \mathbf{I})$

Seperat from 2 until convergence



- Initialize Mixture Models
- Assign GMM components

$$k_i = \arg\min_k \varphi(x_i, k_i, \theta|z_i)$$

Searn GMM parameters

$$\boldsymbol{ heta} = rgmin\sum_i arphi(x_i,k_i,oldsymbol{ heta}|z_i)$$

Estimate segmentation using mincut

 $\mathbf{x} = \arg\min E(\mathbf{x}, \mathbf{k}, \boldsymbol{\theta} | \mathbf{I})$



Seperat from 2 until convergence

- Initialize Mixture Models
- Assign GMM components

$$k_i = \arg\min_k \varphi(x_i, k_i, \theta|z_i)$$

Searn GMM parameters

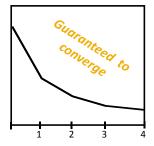
$$oldsymbol{ heta} = rgmin\sum_i arphi(x_i,k_i,oldsymbol{ heta}|z_i)$$

Estimate segmentation using mincut

 $\mathbf{x} = \arg\min E(\mathbf{x}, \mathbf{k}, \boldsymbol{\theta} | \mathbf{I})$

Repeat from 2 until convergence





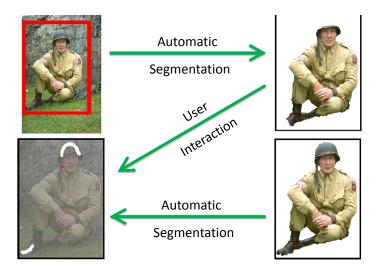
Result

Energy after each Iteration

(日) (同) (三) (三)

3

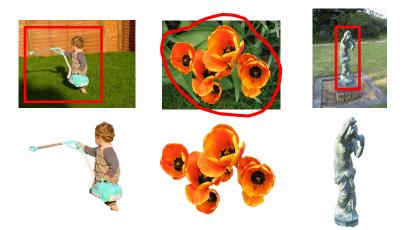
GrabCut - Further editing



April 24, 2013 58 / 63

3

GrabCut - More results



... GrabCut completes automatically

Philipp Krähenbühl (Stanford University)

Segmentation

April 24, 2013 59 / 63

Image: A match a ma

GrabCut - Live demo

• Included in MS Office 2010

3

• Problem: Images contain many pixels

 Even with efficient graph cuts, an MRF formulation has too many nodes for interactive results.

• Efficiency trick: Superpixels

- Group together similar-looking pixels for efficiency of further processing.
- Cheap, local oversegmentation
- Important to ensure that superpixels do not cross boundaries

• Several different approaches possible

- Superpixel code available here
- http:

//www.cs.sfu.ca/~mori/research/superpixels/



< ロ > < 同 > < 三 > < 三

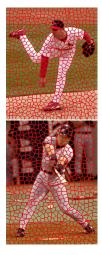
• Problem: Images contain many pixels

 Even with efficient graph cuts, an MRF formulation has too many nodes for interactive results.

• Efficiency trick: Superpixels

- Group together similar-looking pixels for efficiency of further processing.
- Cheap, local oversegmentation
- Important to ensure that superpixels do not cross boundaries
- Several different approaches possible
 - Superpixel code available here
 - http:

//www.cs.sfu.ca/~mori/research/superpixels/



< ロ > < 同 > < 三 > < 三

• Problem: Images contain many pixels

 Even with efficient graph cuts, an MRF formulation has too many nodes for interactive results.

• Efficiency trick: Superpixels

- Group together similar-looking pixels for efficiency of further processing.
- Cheap, local oversegmentation
- Important to ensure that superpixels do not cross boundaries
- Several different approaches possible
 - Superpixel code available here
 - http:

//www.cs.sfu.ca/~mori/research/superpixels/



• Problem: Images contain many pixels

- Even with efficient graph cuts, an MRF formulation has too many nodes for interactive results.
- Efficiency trick: Superpixels
 - Group together similar-looking pixels for efficiency of further processing.
 - Cheap, local oversegmentation
 - Important to ensure that superpixels do not cross boundaries
- Several different approaches possible
 - Superpixel code available here
 - http:

//www.cs.sfu.ca/~mori/research/superpixels/



• Problem: Images contain many pixels

- Even with efficient graph cuts, an MRF formulation has too many nodes for interactive results.
- Efficiency trick: Superpixels
 - Group together similar-looking pixels for efficiency of further processing.
 - Cheap, local oversegmentation
 - Important to ensure that superpixels do not cross boundaries
- Several different approaches possible
 - Superpixel code available here
 - http:

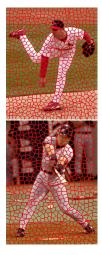
//www.cs.sfu.ca/~mori/research/superpixels/



• Problem: Images contain many pixels

- Even with efficient graph cuts, an MRF formulation has too many nodes for interactive results.
- Efficiency trick: Superpixels
 - Group together similar-looking pixels for efficiency of further processing.
 - Cheap, local oversegmentation
 - Important to ensure that superpixels do not cross boundaries
- Several different approaches possible
 - Superpixel code available here
 - http:

//www.cs.sfu.ca/~mori/research/superpixels/



Improving Efficiency of Segmentation

Problem: Images contain many pixels

- Even with efficient graph cuts, an MRF formulation has too many nodes for interactive results.
- Efficiency trick: Superpixels
 - Group together similar-looking pixels for efficiency of further processing.
 - Cheap, local oversegmentation
 - Important to ensure that superpixels do not cross boundaries
- Several different approaches possible
 - Superpixel code available here
 - http:

//www.cs.sfu.ca/~mori/research/superpixels/



Improving Efficiency of Segmentation

• Problem: Images contain many pixels

- Even with efficient graph cuts, an MRF formulation has too many nodes for interactive results.
- Efficiency trick: Superpixels
 - Group together similar-looking pixels for efficiency of further processing.
 - Cheap, local oversegmentation
 - Important to ensure that superpixels do not cross boundaries
- Several different approaches possible
 - Superpixel code available here
 - http:

//www.cs.sfu.ca/~mori/research/superpixels/



イロト イヨト イヨト イヨト

Improving Efficiency of Segmentation

• Problem: Images contain many pixels

- Even with efficient graph cuts, an MRF formulation has too many nodes for interactive results.
- Efficiency trick: Superpixels
 - Group together similar-looking pixels for efficiency of further processing.
 - Cheap, local oversegmentation
 - Important to ensure that superpixels do not cross boundaries
- Several different approaches possible
 - Superpixel code available here
 - http:

//www.cs.sfu.ca/~mori/research/superpixels/



Pros

- Powerful technique, based on probabilistic model (MRF).
- Applicable for a wide range of problems.
- Very efficient algorithms available for vision problems.
- Becoming a de-facto standard for many segmentation tasks.

• Cons/Issues

- Graph cuts can only solve a limited class of models
 - Submodular energy functions
 - Can capture only part of the expressiveness of MRFs
- Only approximate algorithms available for multi-label case

< ロ > < 同 > < 回 > < 回 > < 回 > < 回

Pros

- Powerful technique, based on probabilistic model (MRF).
- Applicable for a wide range of problems.
- Very efficient algorithms available for vision problems.
- Becoming a de-facto standard for many segmentation tasks.
- Cons/Issues
 - Graph cuts can only solve a limited class of models
 - Submodular energy functions
 - Can capture only part of the expressiveness of MRFs
 - Only approximate algorithms available for multi-label case

Pros

- Powerful technique, based on probabilistic model (MRF).
- Applicable for a wide range of problems.
- Very efficient algorithms available for vision problems.
- Becoming a de-facto standard for many segmentation tasks.
- Cons/Issues
 - Graph cuts can only solve a limited class of models
 - Submodular energy functions
 - Can capture only part of the expressiveness of MRFs.
 - Only approximate algorithms available for multi-label case

Pros

- Powerful technique, based on probabilistic model (MRF).
- Applicable for a wide range of problems.
- Very efficient algorithms available for vision problems.
- Becoming a de-facto standard for many segmentation tasks.
- Cons/Issues
 - Graph cuts can only solve a limited class of models
 - Submodular energy functions
 - \star Can capture only part of the expressiveness of MRFs
 - Only approximate algorithms available for multi-label case

Pros

- Powerful technique, based on probabilistic model (MRF).
- Applicable for a wide range of problems.
- Very efficient algorithms available for vision problems.
- Becoming a de-facto standard for many segmentation tasks.

• Cons/Issues

- Graph cuts can only solve a limited class of models
 - Submodular energy functions
 - \star Can capture only part of the expressiveness of MRFs
- Only approximate algorithms available for multi-label case

Pros

- Powerful technique, based on probabilistic model (MRF).
- Applicable for a wide range of problems.
- Very efficient algorithms available for vision problems.
- Becoming a de-facto standard for many segmentation tasks.

Cons/Issues

- Graph cuts can only solve a limited class of models
 - * Submodular energy functions
 - Can capture only part of the expressiveness of MRFs
- Only approximate algorithms available for multi-label case

・ロン ・四 ・ ・ ヨン ・ ヨン

Pros

- Powerful technique, based on probabilistic model (MRF).
- Applicable for a wide range of problems.
- Very efficient algorithms available for vision problems.
- Becoming a de-facto standard for many segmentation tasks.
- Cons/Issues
 - Graph cuts can only solve a limited class of models
 - * Submodular energy functions
 - Can capture only part of the expressiveness of MRFs
 - Only approximate algorithms available for multi-label case

Pros

- Powerful technique, based on probabilistic model (MRF).
- Applicable for a wide range of problems.
- Very efficient algorithms available for vision problems.
- Becoming a de-facto standard for many segmentation tasks.
- Cons/Issues
 - Graph cuts can only solve a limited class of models
 - ★ Submodular energy functions
 - Can capture only part of the expressiveness of MRFs
 - Only approximate algorithms available for multi-label case

Pros

- Powerful technique, based on probabilistic model (MRF).
- Applicable for a wide range of problems.
- Very efficient algorithms available for vision problems.
- Becoming a de-facto standard for many segmentation tasks.
- Cons/Issues
 - Graph cuts can only solve a limited class of models
 - ★ Submodular energy functions
 - ★ Can capture only part of the expressiveness of MRFs
 - Only approximate algorithms available for multi-label case

Pros

- Powerful technique, based on probabilistic model (MRF).
- Applicable for a wide range of problems.
- Very efficient algorithms available for vision problems.
- Becoming a de-facto standard for many segmentation tasks.
- Cons/Issues
 - Graph cuts can only solve a limited class of models
 - ★ Submodular energy functions
 - ★ Can capture only part of the expressiveness of MRFs
 - Only approximate algorithms available for multi-label case

- Graph theoretic segmentation
 - Normalized Cuts
 - Using texture features
- Segmentation as Energy Minimization
 - Markov Random Fields (MRF) / Conditional Random Fields (CRF)
 - Graph cuts for image segmentation
 - Applications