

## CS 664 Segmentation (2)

**Daniel Huttenlocher** 

#### Recap

- Last time covered perceptual organization more broadly, focused in on pixel-wise segmentation
- Covered local graph-based methods such as MST and Felzenszwalb-Huttenlocher method
- Today
  - Cut-based methods such as grab cut, normalized cuts
  - Iterative local update methods such as mean shift

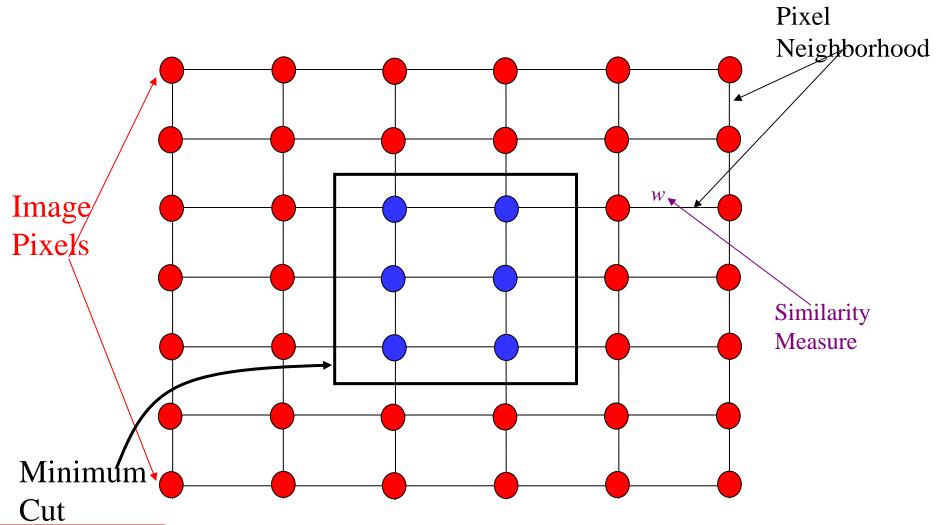


#### **Cut Based Techniques**

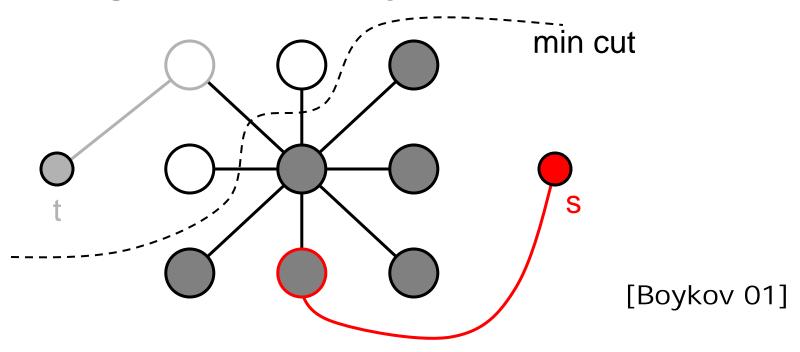
- For costs, natural to consider minimum cost cuts
  - Removing edges with smallest total cost, that cut graph in two parts
  - Graph only has finite-weight edges
- Manually assisted techniques, foreground vs. background
- General segmentation, recursively cut resulting components
  - Question of when to stop



#### **Image Segmentation & Minimum Cut**



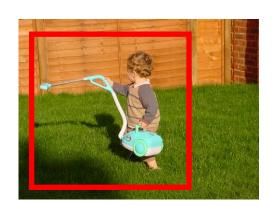
#### Segmentation by Min (s-t) Cut



- Manually select a few fg and bg pixels
  - Infinite cost link from each bg pixel to the "t" node, and each fg pixel to "s" node
  - Compute min cut that separates s from t

#### Grabcut

# [Rother et al., SIGGRAPH 2004]





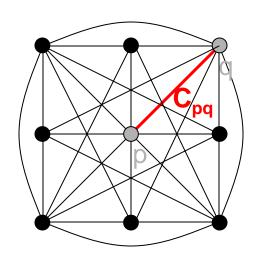


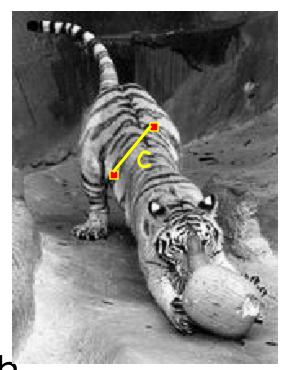






#### **Automatic Cut-Based Segmentation**



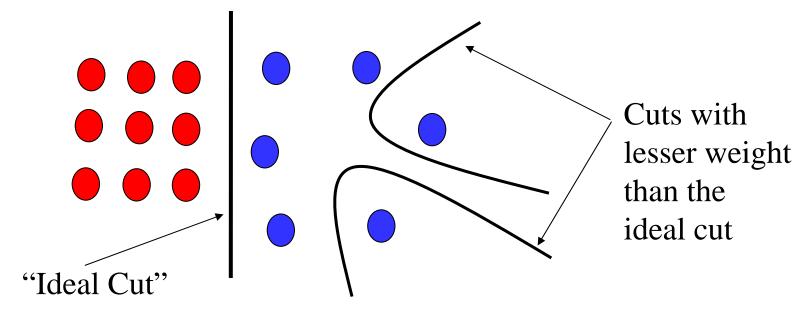


- Fully-connected graph
  - Node for every pixel
  - Link between every pair of pixels, p,q
  - Cost for each link measures similarity



#### **Drawbacks of Minimum Cut**

- Weight of cut proportional to number of edges – preference for small regions
  - Motivation for Shi-Malik normalized cuts



#### **Normalized Cuts**

- A number of normalization criteria have been proposed
- One that is commonly used [Shi&Malik ]

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

Where cut(A,B) is standard definition

$$\sum_{i \in A, j \in B} W_{ij}$$

• And assoc(A,V) =  $\sum_{j} \sum_{i \in A} w_{ij}$ 

## **Computing Normalized Cuts**

 Has been shown this is equivalent to an integer programming problem, minimize

$$\frac{y^{T} (D-W)y}{y^{T} D y}$$

- Subject to the constraint that y<sub>i</sub>∈{1,b} and y<sup>T</sup>D1=0
  - Where 1 vector of all 1's
- W is the affinity matrix
- D is the degree matrix (diagonal)

$$D(i,i) = \sum_{j} w_{ij}$$



## **Approximating Normalized Cuts**

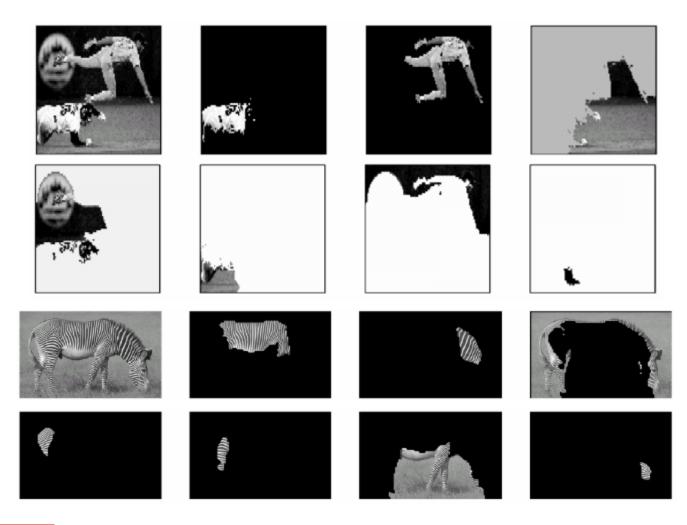
- Integer programming problem NP hard
  - Instead simply solve continuous (real-valued) version
  - This corresponds to finding second smallest eigenvector of

$$(D-W)y_i = \lambda_i Dy_i$$

- Widely used method
  - Works well in practice
    - Large eigenvector problem, but sparse matrices
    - Often resolution reduce images, e.g, 100x100
  - But no longer clearly related to cut problem



## **Normalized Cut Examples**





#### **Another Look [Weiss 99]**

- Consider eigen analysis of affinity matrix
   W = [ w<sub>ii</sub> ]
  - Note W is symmetric; for images w<sub>ij</sub>=w<sub>ji</sub>
  - W also essentially block diagonal
    - With suitable rearrangement of rows/cols so that vertices with higher affinity have nearer indices
    - Entries far from diagonal are small (though not quite zero)
- Eigenvectors of W
  - Recall for real, symmetric matrix forms an orthogonal basis
    - Axes of decreasing "importance"



#### Structure of W

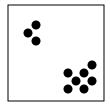
- Eigenvectors of block diagonal matrix consist of eigenvectors of the blocks
  - Padded with zeroes
- Note rearrangement so that clusters lie near diagonal only conceptual
  - Eigenvectors of permuted matrix are permutation of original eigenvectors
- Can think of eigenvectors as being associated with high affinity "clusters"
  - Eigenvectors with large eigenvalues
  - Approximately the case



#### Structure of W

• Consider case of point set where affinities  $w_{ij} = \exp(-(y_i-y_j)^2/\sigma^2)$ 

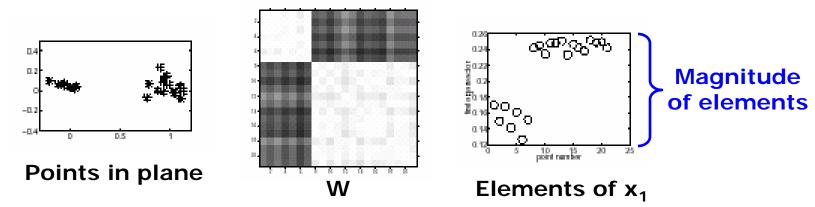
- With two clusters
  - Points indexed to respect clusters for clarity
- Block diagonal form of W
  - Within cluster affinities A, B for clusters
  - Between cluster affinity C



$$M = \begin{pmatrix} C_{L} & B \\ C_{L} & B \end{pmatrix}$$

## First Eigenvector of W

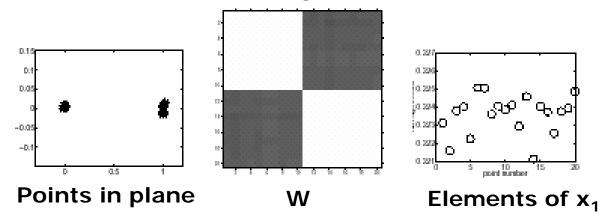
- Recall, vectors x<sub>i</sub> satisfying Wx<sub>i</sub>=λ<sub>i</sub>x<sub>i</sub>
- Consider ordered by eigenvalues λ<sub>i</sub>
  - First eigenvector  $x_1$  has largest eigenvalue  $\lambda_1$
- Elements of first eigenvector serve as "index vector" [Perona, Freeman]
  - Selecting elements of highest affinity cluster





## Clustering

- First eigenvector of W has been suggested as clustering or segmentation criterion
  - For selecting most significant segment
  - Then recursively segment remainder
- Problematic when nonzero non-diagonal blocks (similar affinity clusters)





#### **Understanding Normalized Cuts**

- Intractable discrete graph problem used to motivate continuous (real valued) problem
  - Find second *smallest* "generalized eigenvector"  $(D-W)x_i = \lambda_i Dx_i$
  - Where D is (diagonal) degree matrix  $d_{ii} = \sum_{j} w_{ij}$
- Can be viewed in terms of first two eigenvectors of normalized affinity matrix
  - Let  $N = D^{-1/2}WD^{-1/2}$
  - Note  $n_{ij} = w_{ij} / (\sqrt{d_{ii}} \sqrt{d_{jj}})$ 
    - Affinity normalized by degree of the two nodes

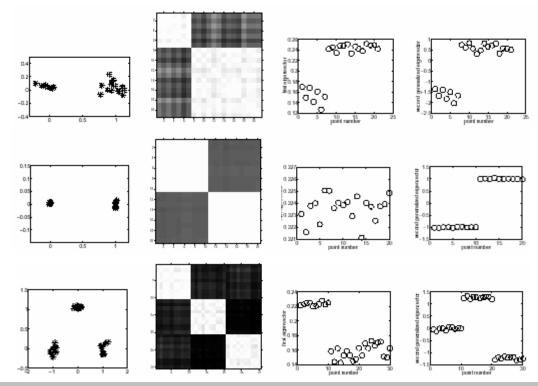
#### **Normalized Affinities**

- Can be shown that
  - If x is an eigenvector of N with eigenvalue  $\lambda$  then D<sup>-1/2</sup>x is a generalized eigenvector of W with eigenvalue 1- $\lambda$
  - The vector D<sup>-1/2</sup>1 is an eigenvector of N with eigenvalue 1
- It follows that
  - Second smallest generalized eigenvector of W is ratio of first two eigenvectors of N
  - So ncut uses normalized affinity matrix N and first two eigenvectors rather than affinity matrix W and first eigenvector



## Contrasting W and N

- Three simple point clustering examples
  - W, first eigenvector of W, ratio of first two eigenvectors of N (generalized eigenvector of W)

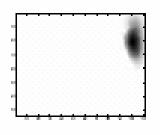


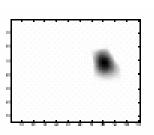


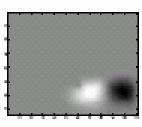
## **Image Segmentation**

- Considering W and N for segmentation
  - Affinity a negative exponential based on distance in x,y,b space
- Eigenvectors of N more correlated with regions

First 4 eigenvectors of W





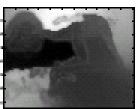


First 4 eigenvectors of N











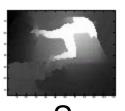
## **Using More Eigenvectors**

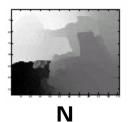
- Based on k largest eigenvectors
  - Construct matrix Q such that (ideally)  $q_{ij}=1$  if i and j in same cluster, 0 otherwise
- Let V be matrix whose columns are first k eigenvectors of W
- Normalize rows of V to have unit Euclidean norm
  - Ideally each node (row) in one cluster (col)
- Let Q=VV<sup>T</sup>
  - Each entry product of two unit vectors

#### Normalization and k Eigenvectors

- Normalized affinities help correct for variations in overall degree of affinity
  - So compute Q for N instead of W
- Contrasting Q with ratio of first two eigenvectors of N (ncut criterion)
  - More clearly selects most significant region
    - Using k=6 eigenvectors
  - Row of Q matrix vs. ratio of eigenvectors of N

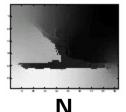














#### **Spectral Methods**

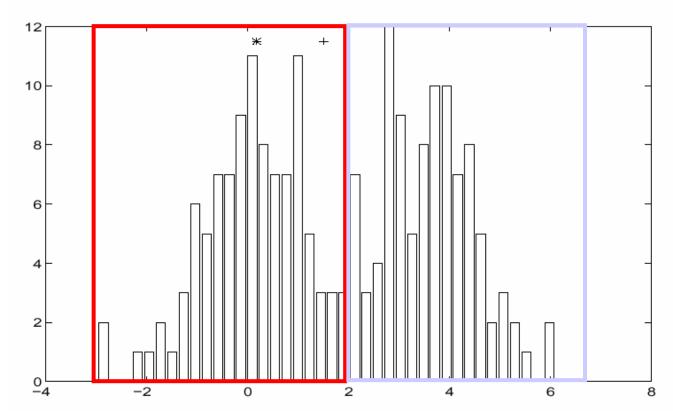
- Eigenvectors of affinity and normalized affinity matrices
- Widely used outside computer vision for graph-based clustering
  - Link structure of web pages, citation structure of scientific papers
  - Often directed rather than undirected graphs

## **Iterative Clustering Methods**

- Techniques such as k-means, but for image segmentation generally have no idea about number of regions
- Mean-shift a nonparametric method

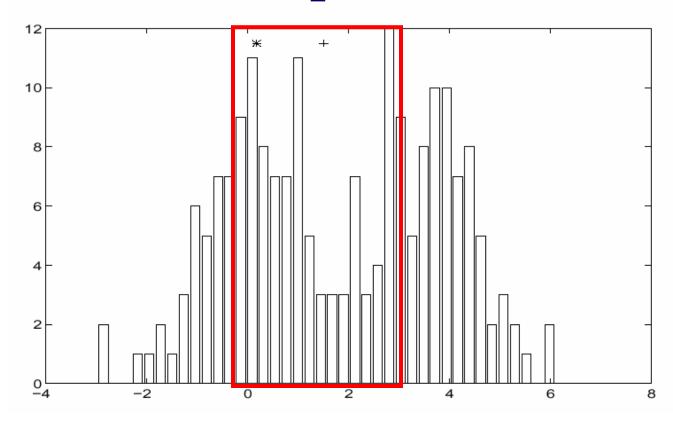


#### Finding Modes in a Histogram



- How Many Modes Are There?
  - Easy to see, less easy to compute

#### Mean Shift [Comaniciu & Meer]



#### Iterative Mode Search

- 1. Initialize random seed, and window W
- 2. Calculate center of gravity ("mean") of W and shift



#### Mean Shift

- Used both for segmentation and for edge preserving filtering
- Operates on collection of points  $X = \{x_1, ..., x_n\}$  in  $R^d$
- Replace each point with value derived from mean shift procedure
  - Searches for a local density maximum by repeatedly shifting a d-dimensional hypersphere of fixed radius h
  - Differs from most clustering, such as k-means in that no fixed number of clusters

#### Mean Shift Procedure

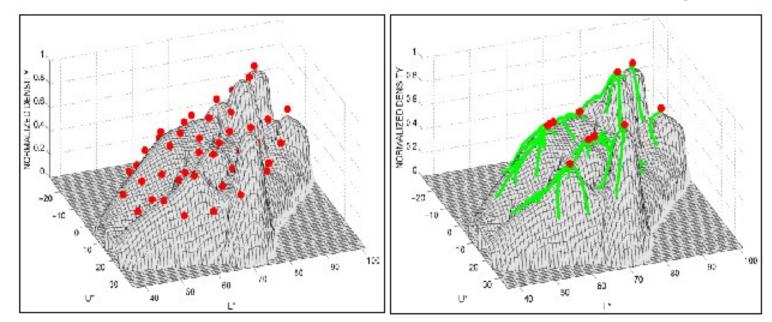
■ For given point  $x \in X$  let  $y_1, ..., y_T$  denote successive locations of that point

$$y_1 = x$$
  
 $y_{k+1} = 1/|S(y_k)| \sum_{x \in S(y^k)} x$ 

- Where  $S(y_k)$  is the subset of X contained in a hyper-sphere of radius h centered at  $y_k$ 
  - The radius h is a fixed parameter of the method
- For a point set X, the mean shift procedure is applied separately to all the points

#### Mean-Shift

- Initialize window around each point
  - Where it shifts determines which region it's in
  - Multiple points will shift to the same region



Mean shift trajectories



## Mean Shift Image Filtering

 Map each image pixel to point in u,v,b space

$$x_i = (u_i, v_i, b_i/\sigma)$$

- Analogous for color images, with three intensity values instead of one
- Scale factor σ normalizes intensity vs. spatial dimensions
- Perform mean shift for each point
  - Let  $Y_i = (U_i, V_i, B_i)$  denote mean shifted value
- Assign result z<sub>i</sub>=(u<sub>i</sub>,v<sub>i</sub>,B<sub>i</sub>)
  - Original spatial coords, mean shifted intensity

# Mean Shift Example





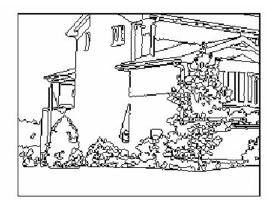


# Mean Shift Example



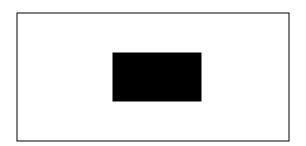
Figure 2: The house image,  $255\times192$  pixels, 9603 colors.





## **Edge Preserving Filtering**

- Mean shift tends to preserve edges
- Edges are where intensity is changing rapidly
- Rapid changes in intensity will result in lower density regions in joint spatialintensity space
- Mean shift finds local density maxima



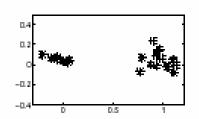
## Mean Shift Clustering

- Run mean shift procedure for each point
- Cluster resulting convergence points that closer than some small constant
- Assign each point label of its cluster
- Analogous to filtering, but with added step of merging cluster that are nearby in the joint spatial-intensity domain



#### **About Mean Shift**

- Convergence to local density maximum
  - Where "local" determined by sphere radius
- Consider simple point set



- Over wide range of sphere radii end up with two clusters
  - Relationship to MST