

# CS 664 Slides #11

## Image Segmentation



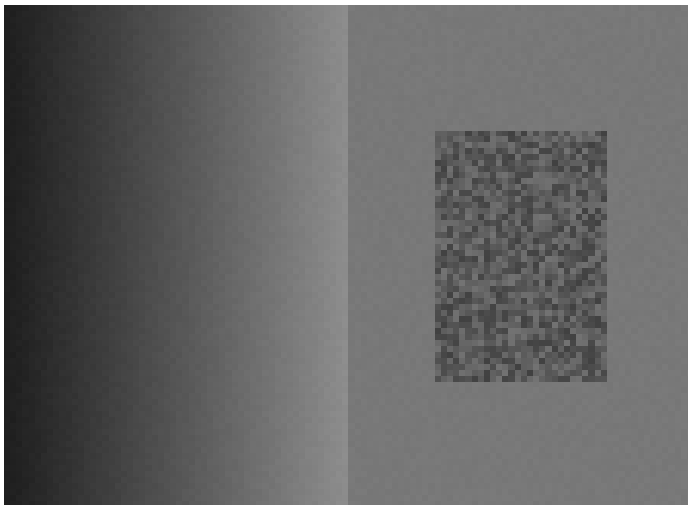
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Fall 2003

# Image Segmentation

- Find regions of image that are “coherent”
- “Dual” of edge detection
  - Regions vs. boundaries
- Related to clustering problems
  - Early work in image processing and clustering
- Many approaches
  - Graph-based
    - Cuts, spanning trees, MRF methods
  - Feature space clustering
  - Mean shift

# A Motivating Example

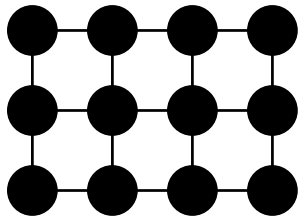
- Image segmentation plays a powerful role in human visual perception
  - Independent of particular objects or recognition



This image has three perceptually distinct regions

# Graph Based Formulation

- $G=(V,E)$  with vertices corresponding to pixels and edges connecting neighboring pixels



4-connected or 8-connected

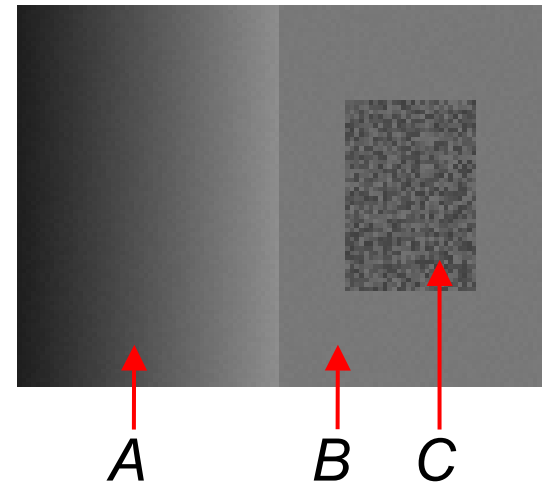
- Weight of edge is magnitude of intensity difference between connected pixels
- A *segmentation*,  $S$ , is a partition of  $V$  such that each  $C \in S$  is connected

# Important Characteristics

- Efficiency
  - Run in time essentially linear in the number of image pixels
    - With low constant factors
    - E.g., compared to edge detection
- Understandable output
  - Way to describe what algorithm does
    - E.g., Canny edge operator and step edge plus noise
- Not purely local
  - Perceptually important

# Motivating Example

- Purely local criteria are inadequate
  - Difference along border between A and B is less than differences within C
- Criteria based on piecewise constant regions are inadequate (e.g., Potts MRF)
  - Will arbitrarily split A into subparts



# MST Based Approaches

- Graph-based representation
  - Nodes corresponding to pixels, edge weights are intensity difference between connected pixels
- Compute minimum spanning tree (MST)
  - Cheapest way to connect all pixels into single component or “region”
- Selection criterion
  - Remove certain MST edges to form components
    - Fixed threshold
    - Threshold based on neighborhood
      - How to find neighborhood

# Component Measure

- Instead of constructing MST based on just the edge weights
  - Consider properties of two components being merged when adding an edge
- Recall Kruskal's MST algorithm adds edges from lowest to highest weight
  - Only when connect distinct components
- Apply criterion based on components to further filter added edges
  - Form of criterion limited by considering edges weight ordered



# Measuring Component Difference

- Let *internal difference* of a component be maximum edge weight in its MST

$$Int(C) = \max_{e \in MST(C,E)} w(e)$$

- Smallest weight such that all pixels of  $C$  are connected by edges of at most that weight

- Let *difference* between two components be minimum edge weight connecting them

$$Dif(C_1, C_2) = \min_{v_i \in C_1, v_j \in C_2} w((v_i, v_j))$$

- Note: infinite if there is no such edge

# Region Comparison Function

- Two components judged to be distinct when  $Dif(C_1, C_2)$  large relative to  $Int(C_1)$  or  $Int(C_2)$ 
  - Require that it be *sufficiently* larger
  - Controlled by (non-negative) threshold function  $\tau$

- Region comparison function  $g(C_1, C_2)$  is true when regions should be distinct, i.e., when

$$Dif(C_1, C_2) > MInt(C_1, C_2)$$

where  $MInt(C_1, C_2)$

$$= \min(Int(C_1) + \tau(C_1), Int(C_2) + \tau(C_2))$$

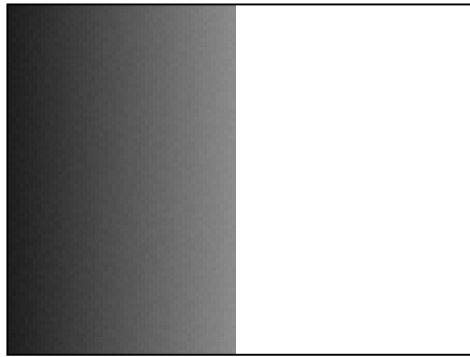
# About the Threshold Function $\tau$

- Intuitively  $Int(C)$  estimates local differences over component
  - Small components give underestimate of local difference – neighboring pixels tend to be similar
    - Thus  $\tau$  should be large in this case
- Use a function inversely proportional to component size  $\tau(C) = k / |C|$ 
  - $k$  is a parameter of the method that captures “scale of observation”
    - Larger  $k$  means prefer larger components
  - Other functions possible, e.g., based on shape

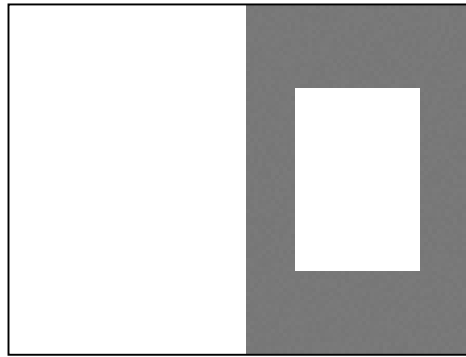
# The Algorithm

0. Sort edges of  $E$  into  $(e_1, \dots, e_n)$ , in order of non-decreasing edge weight
1. Initialize  $S$  with one component per pixel
2. For each  $e_q$  in  $(e_1, \dots, e_n)$  do step 3
3. If weight of  $e_q$  small relative to internal difference of components it connects then merge components, otherwise do nothing  
I.e., if  $w(e_q) \leq MInt(C_i, C_j)$ , where  $C_i, C_j \in S$  are distinct components connected by  $e_q$ , then update  $S$  by merging  $C_i$  and  $C_j$

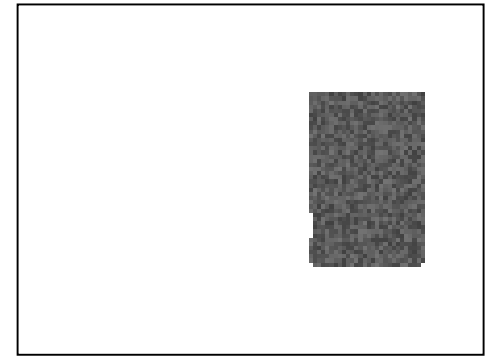
# Regions Found by the Algorithm



*A*



*B*



*C*

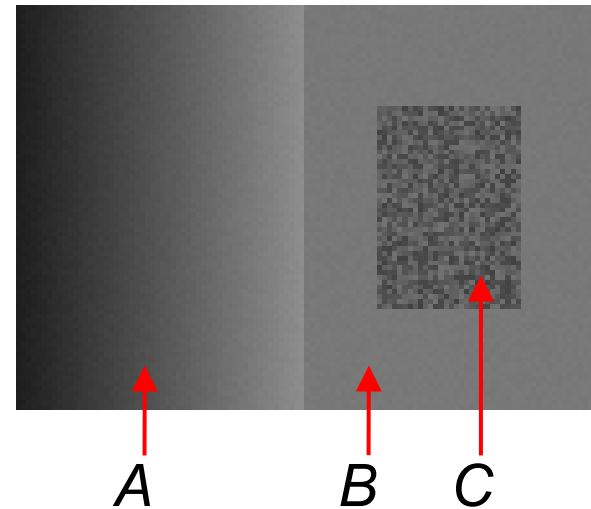
- Three main regions plus a few small ones
- Why the algorithm stops growing these
  - Weight of edges between A and B large wrt max weight MST edges of A and of B
  - Weight of edges between B and C large wrt max weight MST edge of B (but not of C)

# Criteria for a Good Segmentation

- Some predicate for comparing two regions
  - Intuitively, evaluates whether there is evidence for a boundary between two regions
- A segmentation is *too fine* when predicate says no evidence for a boundary
  - Some pair of neighboring regions where predicate false
- A segmentation is *too coarse* when there is some refinement that is not too fine
  - A *refinement* is obtained by splitting one or more regions of a segmentation

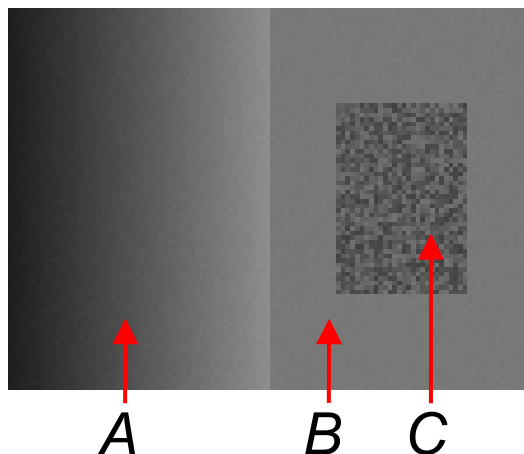
# Good Segmentations and the Example

- Splitting A, B or C would be too fine
- Not splitting A from B or B from C would be too coarse



# Other Algorithms and the Criteria

- Piecewise constant regions (or compact clusters in a color-based feature space)
  - Too fine: arbitrarily split ramp in A into pieces
- Breaking high cost edges in the MST of a graph corresponding to the image
  - Both: merge A with B or split C into multiple pieces





# Properties of the Algorithm

- It is fast,  $O(n \log n)$  for sorting in step 0 and  $O(n\alpha(n))$  for the remaining steps
  - Using union-find with path compression to represent the partition,  $S$
- It produces good segmentations
  - Neither too coarse nor too fine according to the above definitions
    - Despite being a greedy algorithm
- It yields the same results regardless of the order that equal-weight edges are considered
  - Proof a bit involved, won't discuss here

# Components “Freeze”

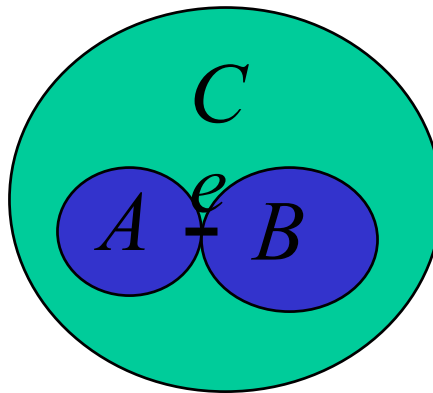
- When two components do not merge, one will be a component of the final segmentation
  - A merge decision is made for an edge  $e_q$  and the two components that it connects  $C_i, C_j$
  - Say the merge does not occur because  $w(e_q) > Int(C_i) + \tau(C_i)$ 
    - Then any subsequent merge involving  $C_i$  will also not occur, because edges are considered in non-decreasing weight order
  - Analogous for  $C_j$ , so when a merge fails one or both of the components involved “freeze”

# Segmentation Not Too Fine

- Follows readily from fact that components “freeze”
  - An edge between two components in final segmentation implies the algorithm decided not to merge when considering this edge
    - Component that caused this decision is frozen, so appears in the final segmentation
- Thus the decision that was true when the edge was considered remains true for the final segmentation

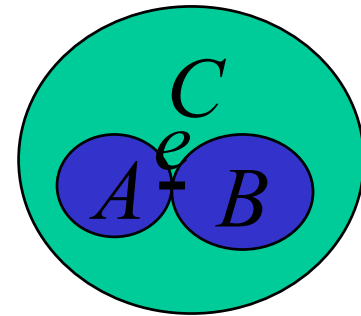
# Segmentation Not Too Coarse

- Means any proper refinement is too fine
- Suppose there was a proper refinement,  $T$ , of the final segmentation,  $S$ , that is not too fine
  - Consider the minimum weight edge,  $e$ , that is between two components  $A, B$  of  $T$  but is within a single component  $C$  of  $S$



# Sketch Continued

- All edges in MST of either  $A$  or  $B$  have weights smaller than  $w(e)$ , say it is  $A$ 
  - Definition of not too fine, and predicate
- Thus algorithm creates  $A$  before considering  $e$ 
  - Because all edges on boundary of  $A$ , but internal to  $C$ , have weight larger than  $w(e)$
- Since  $T$  not too fine, the decision criterion implies the algorithm would freeze  $A$  when considering  $e$



# Closely Related Problems Hard

- What appears to be a slight change
  - Make *Dif* be quantile instead of min
$$k\text{-th}_{v_i \in C_1, v_j \in C_2} w((v_i, v_j))$$
  - Desirable for addressing “cheap path” problem of merging based on one low cost edge
- Makes problem NP hard
  - Reduction from min ratio cut
    - Ratio of “capacity” to “demand” between nodes
- Other methods that we will see are also NP hard and approximated in various ways

# Some Implementation Issues

- Smooth images slightly before processing
  - Remove high variation due to digitization artifacts
- Sorting is dominant time in processing
  - For known edge distribution can in principle do better by binning
- Treat color images as three separate images
  - Components of segmentation are “intersection” of components from each of the three color planes
    - Motivation: significant change in any color channel should result in a region boundary

# Some Example Segmentations



k=300  
320 components  
larger than 10

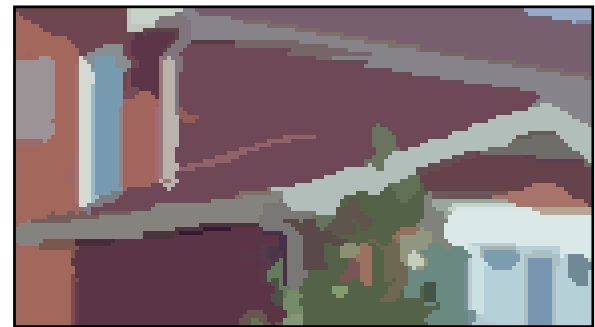


k=200  
323 components  
larger than 10



# Some Shortcomings

- Smoothing can introduce problems
  - “Extra regions” at boundaries
  - Creates “ramps” between regions, thus merge

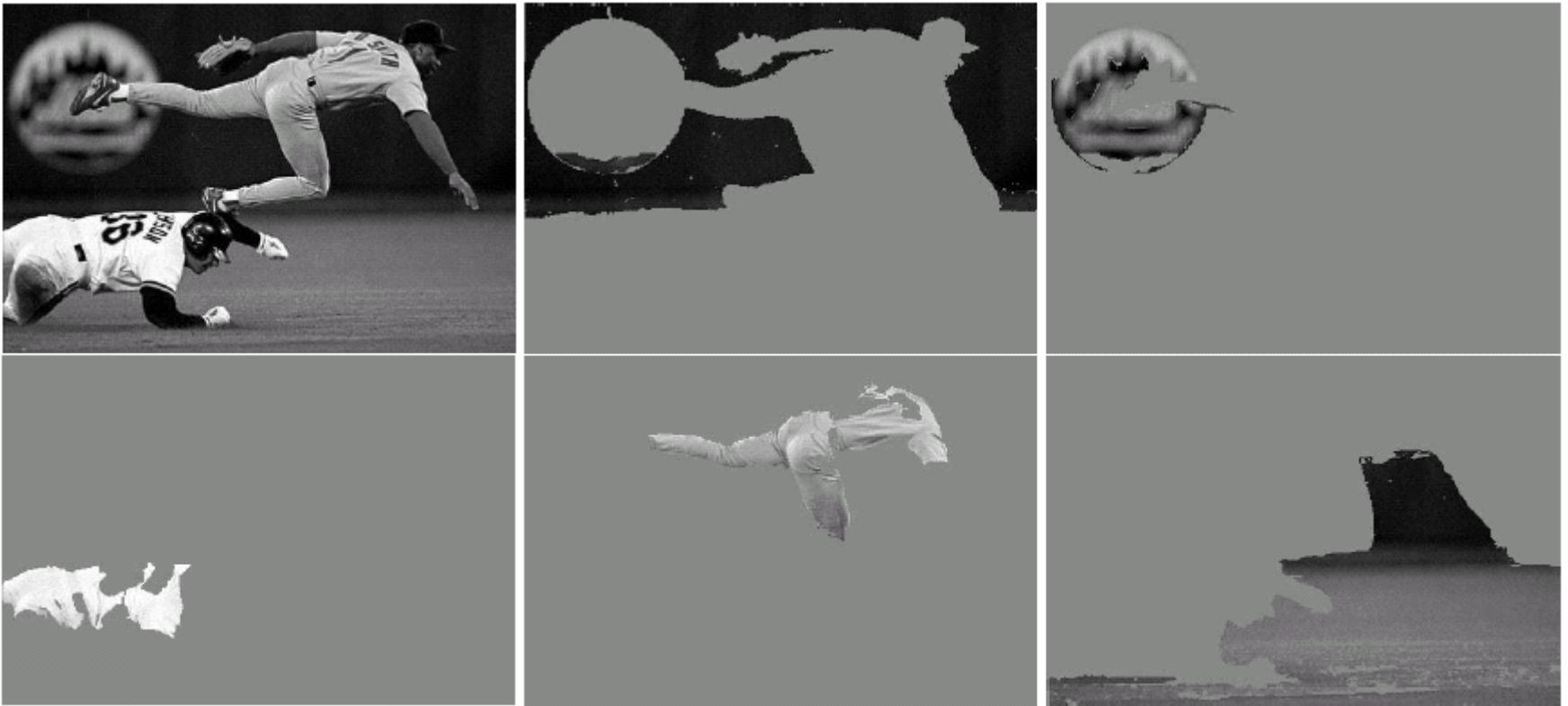


# Simple Object Examples



# Monochrome Example

- Components locally connected (grid graph)
  - Sometimes not desirable

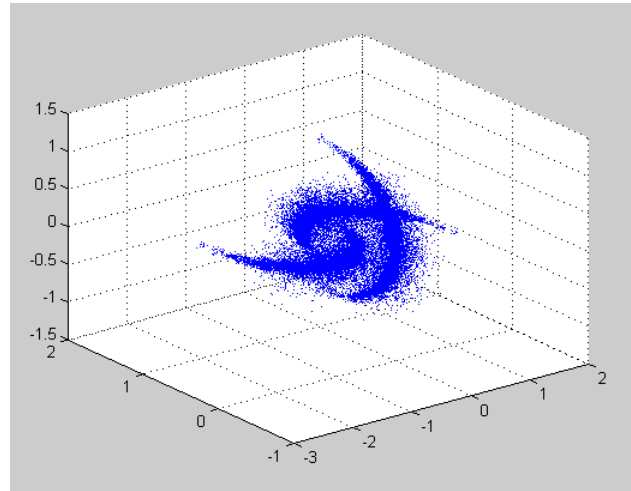


# Clustering: Non-Local Components

- Points in  $d$ -dimensional space
  - Vertex for each point, edge weights based on distance in this space
- Intuitively,  $Int$  measures “density” of clusters
  - Smallest dilation radius such that all points in the cluster are connected
  - When clusters separated by nearly same distance as their “densities” then segmentation is too fine
- For efficiency use a graph with  $O(|V|)$  edges
  - Use Mount’s approximate nearest neighbor algorithm to find nearest neighbors

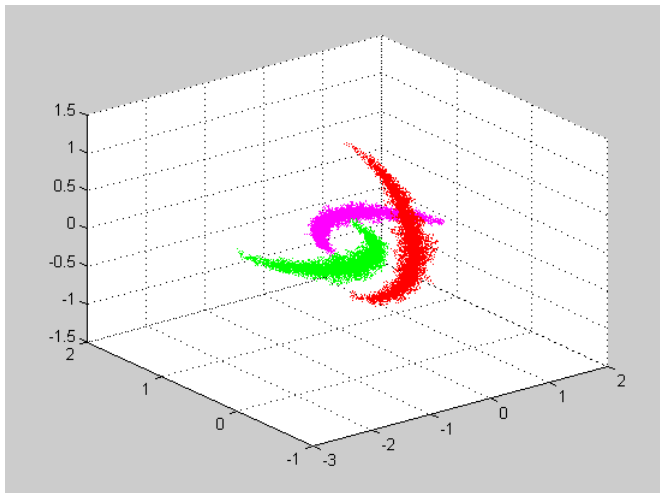
# Clustering Gaussian Point Data

Note: Gaussian not constant density

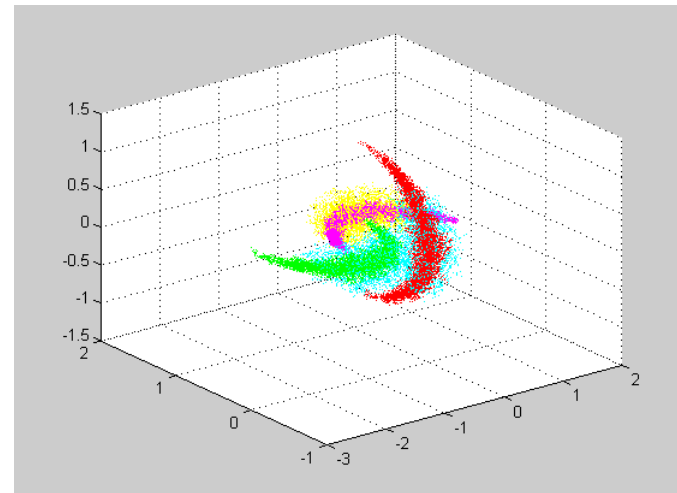


Graph connecting four nearest neighbors to each vertex

$$k = 1$$



*3 largest clusters, 75% classified*



*5 largest clusters, 95% classified*

# Clustering for Image Segmentation

- Treat each pixel as a point in a feature space
  - More than just local intensity or color, incorporate spatial, texture, motion or other differences
- Now regions of segmentation need not be connected in image
- Practical issue, relatively expensive to find nearest neighbors for graph
  - Can use neighbors in some fixed distance, but restricts regions that can be found
  - In examples here use 4 nearest neighbors

# Example Clustering of Image Data

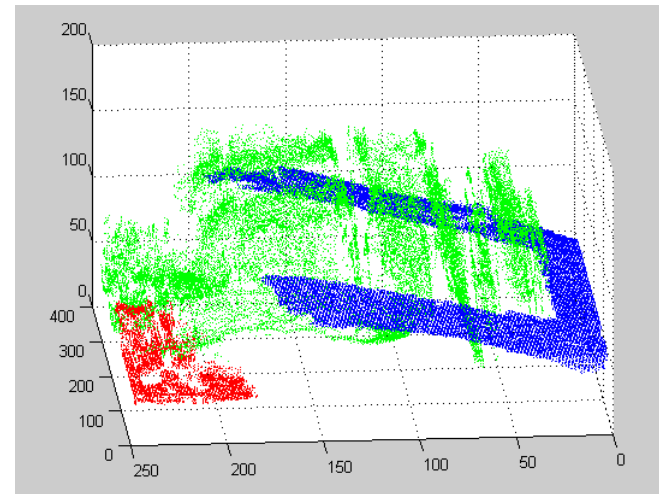
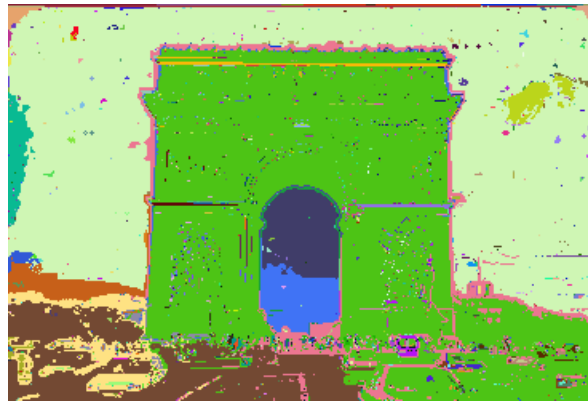
- Segmentation using difference in R,G,B values and in position
  - Distance of 5 pixels same as 1 intensity unit

Non-Local  
Component



# About Clustering for Image Data

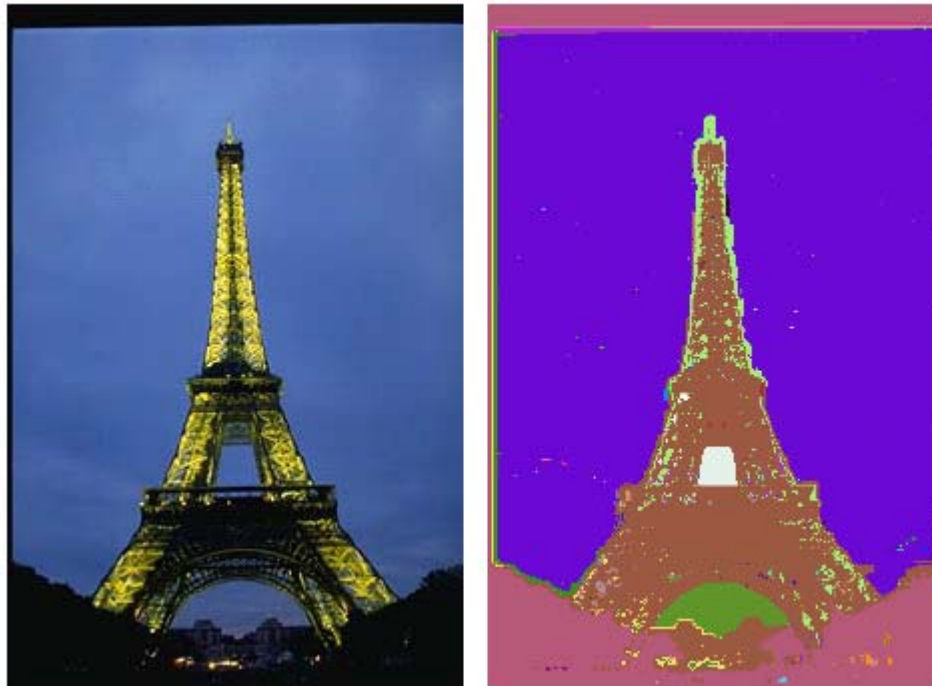
- Meaningful regions in image are not necessarily compact in feature space
- Cheap path in feature space not always apparent in image





# Additional Example

- High variability in illuminated tower pixels



# Beyond Grid Graphs

- Image segmentation methods using affinity (or cost) matrices
  - For each pair of vertices  $v_i, v_j$  an associated weight  $w_{ij}$ 
    - Affinity if larger when vertices more related
    - Cost if larger when vertices less related
  - Matrix  $W = [w_{ij}]$  of affinities or costs
    - $W$  is large, avoid constructing explicitly
    - For images affinities tend to be near zero except for pixels that are nearby
      - E.g., decrease exponentially with distance
    - $W$  is sparse

# 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 non-infinite-weight edges
- For segmentation, recursively cut resulting components
  - Question of when to stop
- Problem is that cuts tend to split off small components
  - Few edges

# Normalized Cuts

- A number of normalization criteria have been proposed
- One that is commonly used

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

- 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

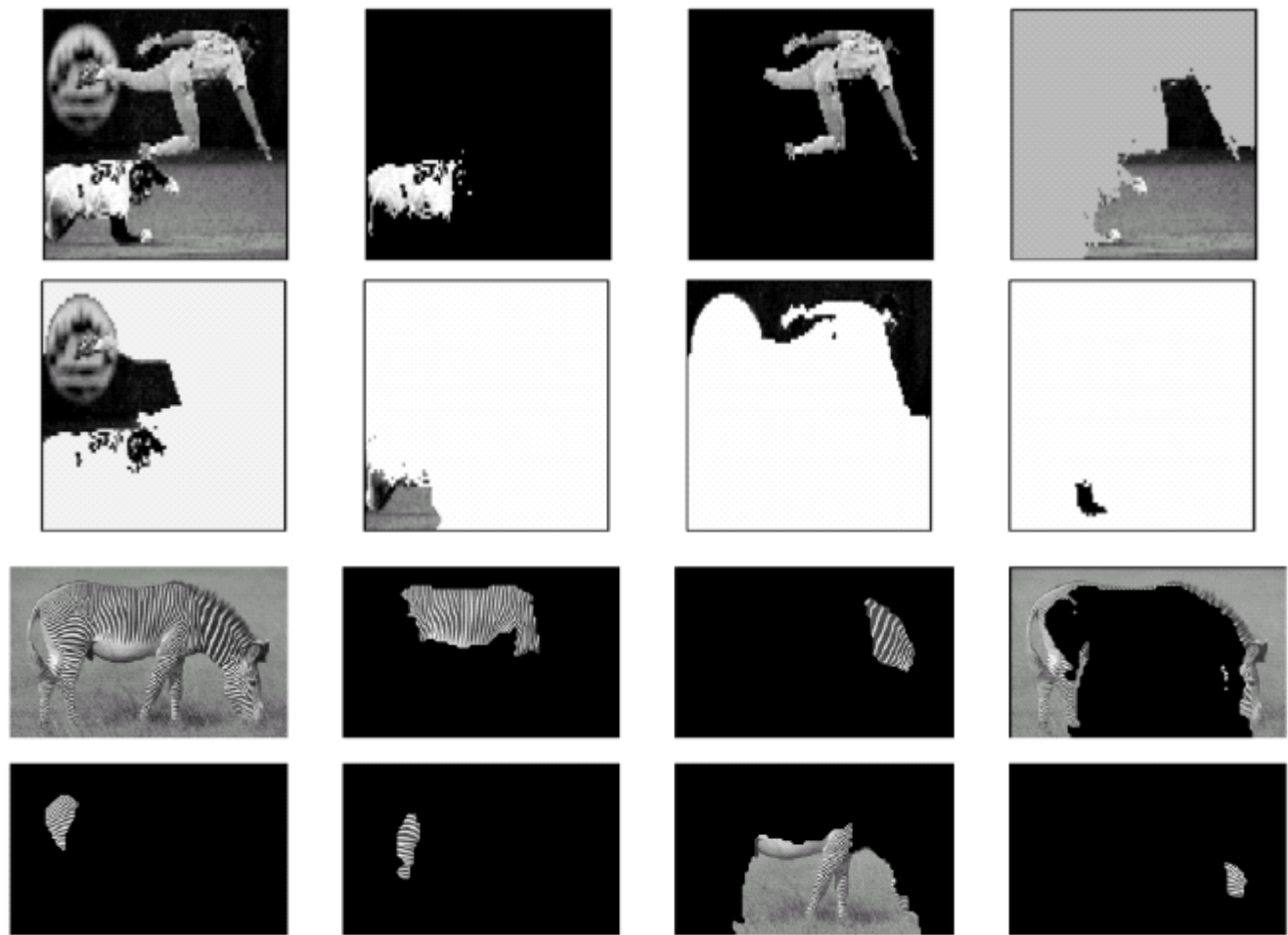
$$\frac{y^T (D-W)y}{y^T D y}$$

- Subject to the constraint that  $y_i \in \{1, b\}$  and  $y^T D \mathbf{1} = 0$ 
  - Where  $\mathbf{1}$  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
$$(D-W)y_i = \lambda_i D y_i$$
- 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 at the Problem

- Consider eigen analysis of affinity matrix
$$W = [ w_{ij} ]$$
  - Note  $W$  is symmetric; for images  $w_{ij}=w_{ji}$
  - $W$  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”

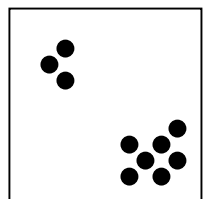


# Structure of $W$

- 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 $W$

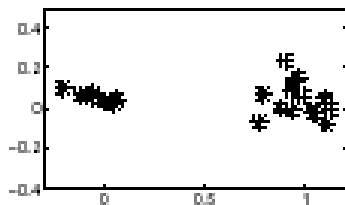
- Consider case of point set where affinities
$$w_{ij} = \exp(-(y_i - y_j)^2 / \sigma^2)$$
- With two clusters
  - Points indexed to respect clusters for clarity
- Block diagonal form of  $W$ 
  - Within cluster affinities  $A, B$  for clusters
  - Between cluster affinity  $C$



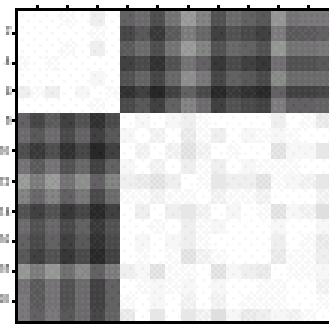
$$W = \begin{pmatrix} A & C \\ C^T & B \end{pmatrix}$$

# First Eigenvector of $W$

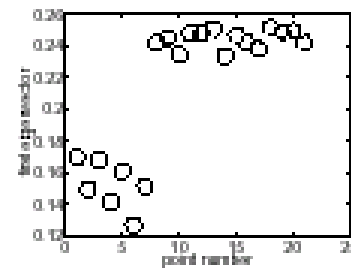
- Recall, vectors  $x_i$  satisfying  $Wx_i = \lambda_i x_i$
- Consider ordered by eigenvalues  $\lambda_i$ 
  - First eigenvector  $x_1$  has largest eigenvalue  $\lambda_1$
- Elements of first eigenvector serve as “index vector”
  - Selecting elements of highest affinity cluster



Points in plane



$W$

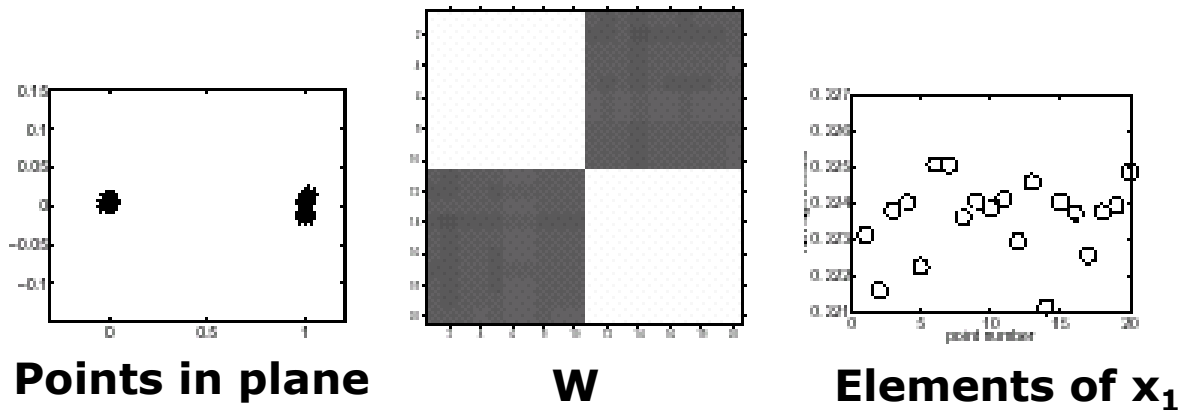


Elements of  $x_1$

Magnitude  
of elements

# Clustering

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



# Understanding Normalized Cuts

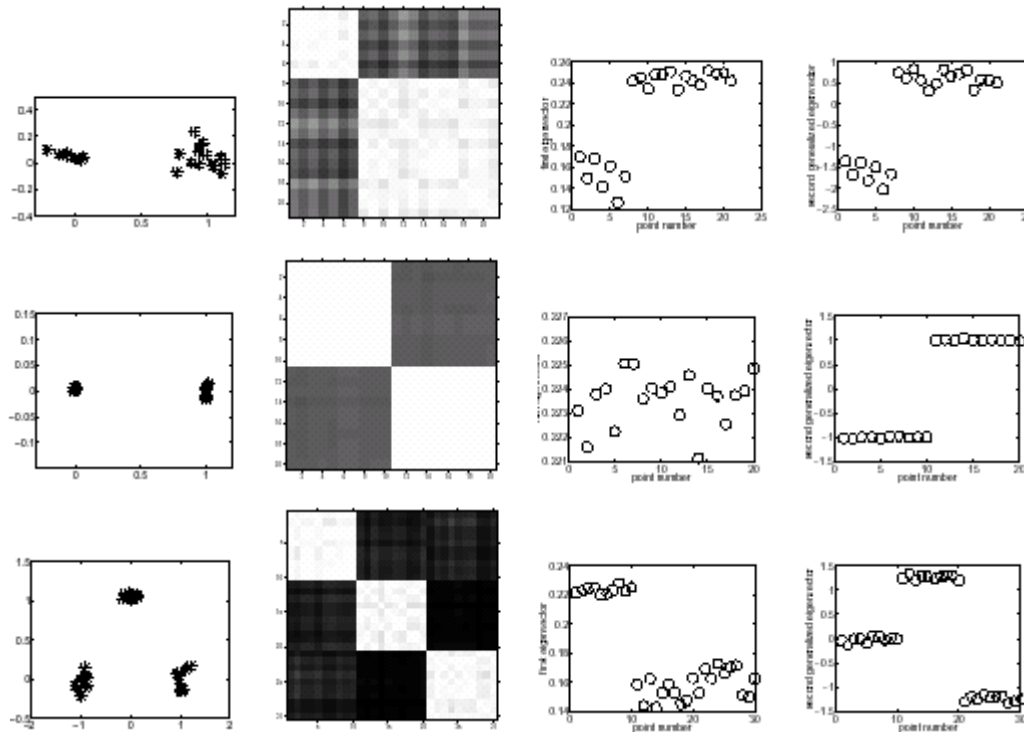
- Intractable discrete graph problem used to motivate continuous (real valued) problem
  - Find second *smallest* “generalized eigenvector”
$$(D-W)x_i = \lambda_i D x_i$$
  - Where D is (diagonal) degree matrix  $d_{ii} = \sum_j w_{ij}$
- Can be viewed in terms of first two eigenvectors of normalized affinity matrix
  - Let  $N = D^{-1/2} W D^{-1/2}$
  - 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  $D^{-1/2}x$  is a generalized eigenvector of  $W$  with eigenvalue  $1-\lambda$
  - The vector  $D^{-1/2}\mathbf{1}$  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

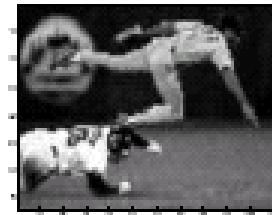
# Contrasting W and N

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

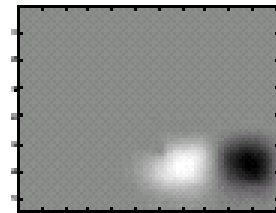
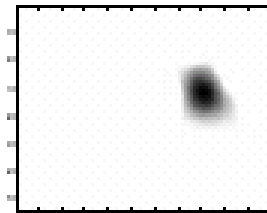
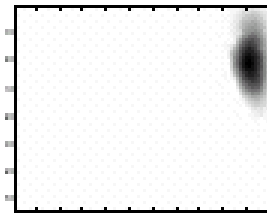
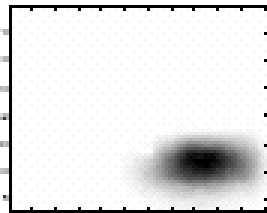


# Image Segmentation

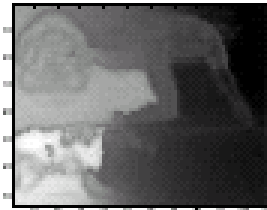
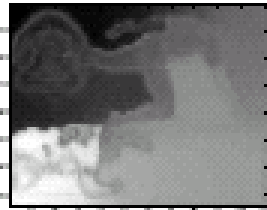
- Considering  $W$  and  $N$  for segmentation
  - Affinity a negative exponential based on distance in  $x, y, b$  space
- Eigenvectors of  $N$  more correlated with regions



**First 4  
eigenvectors  
of  $W$**



**First 4  
eigenvectors  
of  $N$**



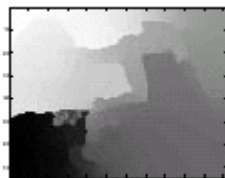
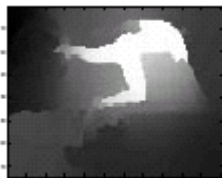


# Using More Eigenvectors

- 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  $Q=VV^T$ 
  - Each entry product of two unit vectors

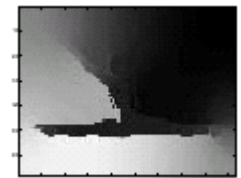
# Normalization and k Eigenvectors

- 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  $N$



Q

N



Q

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

# Mean Shift

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

# Mean Shift Procedure

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

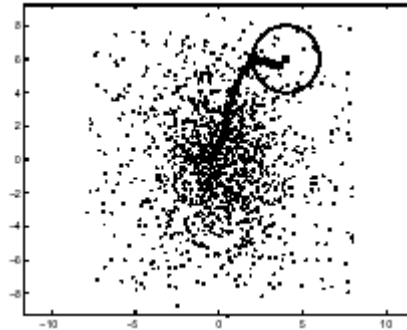
$$y_1 = x$$

$$y_{k+1} = \frac{1}{|S(y_k)|} \sum_{x \in S(y_k)} x$$

- 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

# Illustration of Mean Shift

- Path of successive values of  $y_k$  for given starting point  $x$



- Can be shown that converges to local density maximum

# 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  $\sigma$  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  $z_i = (u_i, v_i, B_i)$ 
  - Original spatial coords, mean shifted intensity

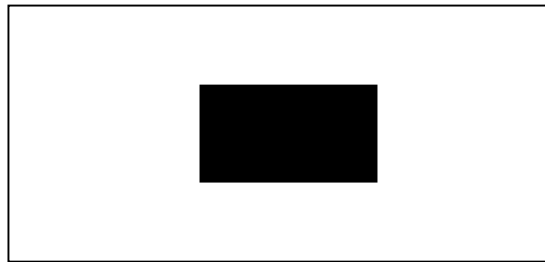
# Mean Shift Example





# 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 spatial-intensity space
- Mean shift finds local density maxima

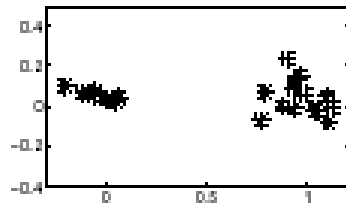


# Mean Shift Clustering

- 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