Lecture: k-means & mean-shift clustering

Juan Carlos Niebles and Ranjay Krishna Stanford Vision and Learning Lab

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Recap: Image Segmentation

• Goal: identify groups of pixels that go together

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Recap: Gestalt Theory

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

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

> **Max Wertheimer (1880-1943)**

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

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Recap: Gestalt Factors

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

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What will we learn today?

- K-means clustering
- Mean-shift clustering

Reading: [FP] Chapters: 14.2, 14.4

D. Comaniciu and P. Meer, Mean Shift: A Robust Approach toward Feature Space Analysis, PAMI 2002.

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Image Segmentation: Toy Example

- These intensities define the three groups.
- We could label every pixel in the image according to which of these primary intensities it is.
	- $-$ i.e., segment the image based on the intensity feature.
- What if the image isn't quite so simple?

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- Now how to determine the three main intensities that define our groups?
- We need to cluster.

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- Goal: choose three "centers" as the representative intensities, and label every pixel according to which of these centers it is nearest to.
- Best cluster centers are those that minimize Sum of Square Distance (SSD) between all points and their nearest cluster center c_i :

$$
SSD = \sum_{clusteri} \sum_{x \in clusteri} (x - c_i)^2
$$

Slide credit: Kristen Grauman Slide credit: Kristen Grauman

Clustering for Summarization

Goal: cluster to minimize variance in data given clusters

 $-$ Preserve information

Cluster center Data
\n
$$
c^*
$$
, $\delta^* = \arg \min_{c, \delta} \frac{1}{N} \sum_{j}^{N} \sum_{i}^{K} \delta_{ij} (c_i - x_j)^2$
\nWhether x_j is assigned to c_i

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Clustering

- With this objective, it is a "chicken and egg" problem:
	- $-$ If we knew the *cluster centers*, we could allocate points to groups by assigning each to its closest center.

 $-$ If we knew the *group memberships*, we could get the centers by computing the mean per group.

- 1. Initialize $(t = 0)$: cluster centers $c_1, ..., c_K$
- 2. Compute δ^t : assign each point to the closest center

$$
\delta^t
$$
 denotes the set of assignment for each x_j to cluster c_i at iteration t

$$
\delta^t = \underset{\delta}{\text{argmin}} \frac{1}{N} \sum_{j}^{N} \sum_{i}^{K} \delta_{ij}^{t-1} \Big(c_i^{t-1} \cdot x_j\Big)^2
$$

1. Computer c^t : update cluster centers as the mean of the points

$$
c^{t} = \operatorname*{argmin}_{c} \frac{1}{N} \sum_{j}^{N} \sum_{i}^{K} \delta_{ij}^{t} \left(c_{i}^{t-1} x_{j} \right)^{2}
$$

1. Update $t = t + 1$, Repeat Step 2-3 till stopped

- 1. Initialize $(t = 0)$: cluster centers $c_1, ..., c_K$
	- Commonly used: random initialization
	- Or greedily choose K to minimize residual

2. Compute δ^t : assign each point to the closest center

- Typical distance measure:
	- Euclidean $sim(x, x') = x^T x'$
	- Cosine $sim(x, x') = x^T x' / (\Vert x \Vert \cdot \Vert x' \Vert)$
	- Others
- 1. Computer c^t : update cluster centers as the mean of the points

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$$
c^{t} = \operatorname*{argmin}_{c} \frac{1}{N} \sum_{j}^{N} \sum_{i}^{K} \delta_{ij}^{t} \left(c_{i}^{t-1} x_{j} \right)^{2}
$$

- 2. Update $t = t + 1$, Repeat Step 2-3 till stopped
	- c^t doesn't change anymore.

• Java demo:

http://home.dei.polimi.it/matteucc/Clustering/tutorial_html/AppletKM.html

- Converges to a *local minimum* solution
	- $-$ Initialize multiple runs

Better fit for spherical data

Need to pick K (# of clusters)

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Segmentation as Clustering

Original image

2 clusters

3 clusters

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K-Means++

- Can we prevent arbitrarily bad local minima?
- 1. Randomly choose first center.
- 2. Pick new center with prob. proportional to $(x c_i)^2$
	- (Contribution of *x* to total error)
- 3. Repeat until *K* centers.
- Expected error = $O(\log K)^*$ optimal

Feature Space

- Depending on what we choose as the *feature space*, we can group pixels in different ways.
- Grouping pixels based on intensity similarity

Feature space: intensity value (1D)

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Feature Space

- Depending on what we choose as the *feature space*, we can group pixels in different ways. R=255
- Grouping pixels based on color similarity

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Feature Space

- Depending on what we choose as the *feature space*, we can group pixels in different ways.
- Grouping pixels based on texture similarity

Filter bank of 24 filters

Feature space: filter bank responses (e.g., 24D)

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Smoothing Out Cluster Assignments

Assigning a cluster label per pixel may yield outliers:

• How can we ensure they are spatially smooth?

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Segmentation as Clustering

- Depending on what we choose as the *feature space*, we can group pixels in different ways.
- Grouping pixels based on *intensity+position* similarity

 \Rightarrow Way to encode both *similarity* and *proximity.*

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K-Means Clustering Results

- K-means clustering based on intensity or color is essentially vector quantization of the image attributes
	- $-$ Clusters don't have to be spatially coherent

K-Means Clustering Results

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	- Clusters don't have to be spatially coherent

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• Clustering based on (r,g,b,x,y) values enforces more spatial coherence

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How to evaluate clusters?

• Generative

 $-$ How well are points reconstructed from the clusters?

- Discriminative
	- $-$ How well do the clusters correspond to labels?
		- Can we correctly classify which pixels belong to the panda?

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 $-$ Note: unsupervised clustering does not aim to be discriminative as we don't have the labels.

How to choose the number of clusters?

Try different numbers of clusters in a validation set and look at performance.

We can plot the objective function values for k equals 1 to $6...$

The abrupt change at $k = 2$, is highly suggestive of two clusters in the data. This technique for determining the number of clusters is known as "knee finding" or "elbow finding".

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K-Means pros and cons

- Pros
	- Finds cluster centers that minimize conditional variance (good representation of data)
	- Simple and fast, Easy to implement
- Cons
	- Need to choose K
	- Sensitive to outliers
	- **Prone to local minima**
	- All clusters have the same parameters (e.g., distance measure is nonadaptive)
	- $*$ Can be slow: each iteration is $O(KNd)$ for N d-dimensional points
- Usage
	- Unsupervised clustering
	- Rarely used for pixel segmentation

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What will we learn today?

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- Mean-shift clustering

Reading: [FP] Chapters: 14.2, 14.4

D. Comaniciu and P. Meer, Mean Shift: A Robust Approach toward Feature Space Analysis, PAMI 2002.

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Mean-Shift Segmentation

• An advanced and versatile technique for clusteringbased segmentation

http://www.caip.rutgers.edu/~comanici/MSPAMI/msPamiResults.html

D. Comaniciu and P. Meer, Mean Shift: A Robust Approach toward Feature Space Analysis, PAMI 2002.

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Mean-Shift Algorithm

• Iterative Mode Search

- 1. Initialize random seed, and window W
- 2. Calculate center of gravity (the "mean") of W :
- 3. Shift the search window to the mean
- 4. Repeat Step 2 until convergence

$$
\sum_{x \in W} xH(x)
$$

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Real Modality Analysis

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Mean-Shift Clustering

- Cluster: all data points in the attraction basin of a mode
- Attraction basin: the region for which all trajectories lead to the same mode

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Mean-Shift Clustering/Segmentation

- 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

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Mean-Shift Segmentation Results

http://www.caip.rutgers.edu/~comanici/MSPAMI/msPamiResults.html

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More Results

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More Results

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Problem: Computational Complexity

Many computations will be redundant.

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Speedups: Basin of Attraction

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Speedups

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Technical Details

Given *n* data points $x_i \in \mathbb{R}^d$, the multivariate kernel density estimate using a radially symmetric kernel¹ (e.g., Epanechnikov and Gaussian kernels), $K(\mathbf{x})$, is given by,

$$
\hat{f}_K = \frac{1}{nh^d} \sum_{i=1}^n K\left(\frac{\mathbf{x} - \mathbf{x}_i}{h}\right),\tag{1}
$$

where h (termed the *bandwidth* parameter) defines the radius of kernel. The radially symmetric kernel is defined as,

$$
K(\mathbf{x}) = c_k k(||\mathbf{x}||^2),\tag{2}
$$

where c_k represents a normalization constant.

Comaniciu & Meer, 2002

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Technical Details

where $g(x) = -k'(x)$ denotes the derivative of the selected kernel profile.

- Term1: this is proportional to the density estimate at x (similar to equation 1 from the previous slide).
- Term2: this is the mean-shift vector that points towards the direction of maximum density.

Comaniciu & Meer, 2002

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Technical Details

Finally, the mean shift procedure from a given point x_t is:

1. Computer the mean shirt vector m:

$$
\left[\frac{\sum\limits_{i=1}^{n} \mathbf{x}_i g\left(\left\|\frac{\mathbf{x}-\mathbf{x}_i}{h}\right\|^2\right)}{\sum\limits_{i=1}^{n} g\left(\left\|\frac{\mathbf{x}-\mathbf{x}_i}{h}\right\|^2\right)} - \mathbf{x}\right]
$$

2. Translate the density window:

$$
\mathbf{x}_i^{t+1} = \mathbf{x}_i^t + \mathbf{m}(\mathbf{x}_i^t).
$$

3. Iterate steps 1 and 2 until convergence.

$$
\nabla f(\mathbf{x}_i)=0.
$$

Comaniciu & Meer, 2002

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Summary Mean-Shift

- Pros
	- $-$ General, application-independent tool
	- $-$ Model-free, does not assume any prior shape (spherical, elliptical, etc.) on data clusters
	- $-$ Just a single parameter (window size h)
		- h has a physical meaning (unlike k-means)
	- $-$ Finds variable number of modes
	- $-$ Robust to outliers
- Cons
	- $-$ Output depends on window size
	- $-$ Window size (bandwidth) selection is not trivial
	- Computationally (relatively) expensive (~2s/image)
	- $-$ Does not scale well with dimension of feature space

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