

## TSBB15 Computer Vision Lecture 6 Clustering and Learning



# Why learning?

- Learning is a very important part of Computer Vision:
  - Parameter tuning
     Adaptation to changing conditions
     Finding patterns in data



- Most Computer Vision systems are complex pieces of software.
- The more complex a system is, the more parameters it has.



- Most Computer Vision systems are complex pieces of software.
- The more complex a system is, the more parameters it has. E.g. filter sizes, thresholds for detection etc. These need to be tuned!





• Tuning in brief:

 Give examples of the desired behaviour of an algorithm.
 Look for the parameters that produce the desired behaviour.

If you let the computer look for the parameters, tuning becomes learning.



• Example:

Automatically decide which motion vectors are  $good(v \in G)$  and which are  $bad(v \in B)$ .



• Look for tracker parameters that maximise:  $J(p_1,...,p_N) = |G|/(|G|+|B|)$ 



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- Computer Vision systems that are deployed in live situations face changing conditions. E.g. different illumination at night and during the day.
- In order to cope with changes, a vision system needs to be adaptive.
- Example: Background models introduced later in this lecture.



### Finding patterns in data

- Recognition and matching (LE 8) uses learned features (or tuned).

 Applications such as: object recognition, object tracking, image captioning etc.





### Learning systems

• Batch learning: learn once, use forever

Online learning: learn continuously



### Learning systems

- Batch learning: *learn once, use forever* Is used to automatically tune parameters, features, classifiers etc.
- Online learning: *learn continuously* Is used to automatically adapt e.g. classifiers and trackers to changing conditions.



## Today's topics

- Learning paradigms
- K-means clustering
- Mixture models and EM
- Background models
- Meanshift clustering
- Generalised Hough Transforms (GHT)
- Channel clustering



• Different learning situations/paradigms:

Supervised learning Reinforcement learning Unsupervised learning

 Covered in depth in: TBMI26 Neural Networks and Learning Systems



• Different learning situations/paradigms:

Supervised learning Reinforcement learning Unsupervised learning ←this lecture

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Supervised learning
 learn y=f(x) from examples {x<sub>n</sub>,y<sub>n</sub>}<sup>1</sup>

 function approximation





- Unsupervised learning learn y=f(x) from examples {x<sub>n</sub>}<sup>1</sup>N =manifold learning or clustering
  - Manifold learning finds low dimensional representations of high dimensional data.
     E.g. coordinates on a surface in nD.



- Unsupervised learning learn y=f(x) from examples {x<sub>n</sub>}<sup>1</sup>
   =manifold learning or clustering
  - Manifold learning finds low dimensional representations of high dimensional data.
     E.g. coordinates on a surface in nD.
- This lecture is mainly about clustering.
- $y \in \mathbb{N}$ , i.e. each sample  $\mathbf{x}_n$  is assigned a cluster *label*.





– Our input is a set of data points  $\{\mathbf{x}_n\}_1^N$ 





- Each data point  $\{\mathbf{x}_n\}_1^N$  is assigned a cluster label  $y \in [1 \dots K]$ , and a prototype  $\{\mathbf{p}_k\}_1^K$ 





 A good clustering has small distances between prototypes and samples within that cluster:

$$J(\mathbf{p}_1, \dots, \mathbf{p}_K) = \sum_{k=1}^K \sum_{n=1}^N \delta[y_n = k] ||\mathbf{x}_n - \mathbf{p}_k||^2$$



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- NP-complete problem.
- K-means clustering [MacQueen'67] is a useful heuristic.



#### K-means clustering

- 1. Pick random sample points as cluster prototypes.
- 2. Assign cluster labels  $\{y_n\}_1^N$  to samples  $\{\mathbf{x}_n\}_1^N$ according to prototype distances  $d_k^2 = ||\mathbf{x}_n - \mathbf{p}_k||^2$
- 3. Assign prototypes as averages of samples within cluster:  $\mathbf{p}_{k} = \frac{1}{|\{y_{n} = k\}|} \sum_{n=1}^{N} \delta[y_{n} = k] \mathbf{x}_{n}$
- 4. Repeat 2-3 until labels stop changing.



#### K-means clustering

- K-means finds a *local min* of the cost:  $J(\mathbf{p}_1, \dots, \mathbf{p}_K) = \sum_{k=1}^K \sum_{n=1}^N \delta[y_n = k] ||\mathbf{x}_n - \mathbf{p}_k||^2$
- Issue 1:Bad repeatability:



Issue 2:What is the value of K?



## Fuzzy K-means clustering

- Fix (partial) for repeatability:
- Replace binary indicator function  $\delta[y_n = k]$ with a continuous weight,  $w_{kn}$ , for each sample.

$$J(\mathbf{p}_{1}, \dots, \mathbf{p}_{K}) = \sum_{k=1}^{n} \sum_{n=1}^{n} w_{kn} ||\mathbf{x}_{n} - \mathbf{p}_{k}||^{2}$$

- Smoother cost fcn  $\Rightarrow$  fewer local min.
- Called *fuzzy k-means* or *fuzzy c-means*.



## Fuzzy K-means clustering

- 1. Pick random sample points as cluster prototypes.
- 2. Assign weights,  $w_{kn}$ , to samples  $\{\mathbf{x}_n\}_1^N$ according to  $w_{kn} = 1/(||\mathbf{x}_n - \mathbf{p}_k||^2 + \epsilon)$
- 3. Assign prototypes as weighted averages of samples:  $\mathbf{p}_{k} = \frac{1}{\sum_{n=1}^{N} w_{kn}} \sum_{n=1}^{N} w_{kn} \mathbf{x}_{n}$
- 4. Repeat 2-3 until labels stop changing.



#### K-means problems

- Fix for the local min problem:
  - Run the algorithm many times, and pick the solution with the lowest *J*.
- Steps 2,3 can be seen as special cases of the EM-algorithm [Dempster et al. 77]
- more on this soon.
- First we need to introduce *mixture models*.



- A *generative model* for data that may come from several distributions.
- E.g. pixel values at a step edge with uncertain location:



k=1



#### Mixture models

• We model the probability density of pixel intensity / as:  $p(I) = \sum_{k}^{K} p(I|\Gamma_{k})P(\Gamma_{k})$ 



 We model the probability density of pixel intensity I as:  $p(I) = \sum_{k=1}^{N} p(I|\Gamma_k) P(\Gamma_k)$ k=1• *Mixture probabilities*:  $(\Gamma_k) = 1$ k=1e.g.  $P(\Gamma_1) = P(\Gamma_2) = 0.5$ gives this p(I): 0.5



- We model the probability density of pixel intensity / as:  $p(I) = \sum_{k}^{K} p(I|\Gamma_k) P(\Gamma_k)$
- Mixture components:

e.g.  

$$p(I|\Gamma_k) = \frac{1}{\sqrt{2\pi}\sigma_k} e^{-0.5(I-\mu_k)^2/\sigma_k^2}$$

Gaussian mixture model



k=1

 $p(I|\Gamma_k)$ 



• Gaussian mixture components:

$$p(I|\Gamma_k) = \frac{1}{\sqrt{2\pi\sigma_k}} e^{-0.5(I-\mu_k)^2/\sigma_k^2}$$

• Notation conditioned on the parameters:

$$p(I|\mu_k, \sigma_k) = \frac{1}{\sqrt{2\pi\sigma_k}} e^{-0.5(I-\mu_k)^2/\sigma_k^2}$$

• Also the mixture probabilities are parameters:

$$P(\Gamma_k) = \pi_k$$
, where  $\sum_k \pi_k = 1$ 



• Given a set of measurements,  $\{I_n\}_1^N$ how do we estimate the parameters of the mixture distribution p(I)?

$$p(I) = \sum_{k=1}^{K} p(I|\Gamma_k) P(\Gamma_k)$$



• Given a set of measurements,  $\{I_n\}_1^N$ how do we estimate the parameters of the mixture distribution p(I)?

$$p(I|\{\pi_k, \mu_k, \sigma_k\}_1^K) = \sum_{k=1}^K \pi_k p(I|\mu_k, \sigma_k)$$

- This can be done with the EM algorithm.
- Note similarities with K-means below.



- 1. Postulate a mixture distribution.
- 2. E: Compute partial memberships,  $w_{kn}$ , with  $\sum_{k=1}^{K} w_{kn} = 1$  to samples  $\{I_n\}_1^N$ , using the mixture distribution.
- 3. M: Use partial memberships to estimate mixture distribution parameters.
- 4. Repeat 2-3 until convergence.



- For the mixture:  $p(I|\{\pi_k, \mu_k, \sigma_k\}_1^K) = \sum_{k=1}^K \pi_k p(I|\mu_k, \sigma_k)$
- The E-step becomes:

$$\tilde{w}_{kn} = \pi_k p(I_n | \mu_k, \sigma_k)$$
$$w_{kn} = \tilde{w}_{kn} / \sum_{l=1}^K \tilde{w}_{ln}$$


- For the mixture:  $p(I|\{\pi_k, \mu_k, \sigma_k\}_1^K) = \sum_{k=1}^K \pi_k p(I|\mu_k, \sigma_k)$
- The E-step becomes:

$$\begin{split} \tilde{w}_{kn} &= \pi_k p(I_n | \mu_k, \sigma_k) \\ w_{kn} &= \tilde{w}_{kn} / \sum_{l=1}^K \tilde{w}_{ln} \\ \end{split} \\ \end{split} \\ \end{split} \\ \end{split} \\ \end{split}$$



• The M-step becomes:

$$\pi_k = P(\Gamma_k) = \frac{1}{N} \sum_{n=1}^N w_{kn}$$

• and, assuming a Gaussian mixture:

$$\mu_{k} = \frac{1}{\sum_{n=1}^{N} w_{kn}} \sum_{n=1}^{N} w_{kn} I_{n}$$
$$\sigma_{k}^{2} = \frac{1}{\sum_{n=1}^{N} w_{kn}} \sum_{n=1}^{N} w_{kn} (I_{n} - \mu_{k})^{2}$$



- Generalizes to higher dimensions.
- e.g. in 2D we have 5 parameters in each mixture component:

$$\mu = \begin{pmatrix} \mu_1 \ \mu_2 \end{pmatrix} \quad \mathbf{\Sigma} = \begin{pmatrix} \sigma_{11} & \sigma_{12} \ \sigma_{12} & \sigma_{22} \end{pmatrix}$$

 Just like K-means, EM also finds a local min.



• Demo for 2D case:





- A popular application of mixture models is background modelling (SHB 16.5.1):
  - Estimate a mixture model for the image *in* each pixel.
  - Pixel values far from the mixture are seen as foreground pixels.
  - Popular way track e.g. people and cars in stationary surveillance cameras.
  - Fast compared to motion estimation.



Background modelling+shadow detection



CVL Master thesis of John Wood 2007



- Samples now arrive one at a time.
- EM uses a batch update:

$$\mu_{k} = \frac{1}{\sum_{n=1}^{N} w_{kn}} \sum_{\substack{n=1\\N}}^{N} w_{kn} I_{n}$$
$$\sigma_{k}^{2} = \frac{1}{\sum_{n=1}^{N} w_{kn}} \sum_{n=1}^{N} w_{kn} (I_{n} - \mu_{k})^{2}$$

• On-line update is needed



- Samples now arrive one at a time.
- On-line update:

$$\mu_k[n] = (1 - \alpha)\mu_k[n - 1] + \alpha I_n$$
  

$$\sigma_k^2[n] = (1 - \alpha)\sigma_k^2[n - 1] + \alpha (I_n - \mu_k[n - 1])^2$$
  

$$\pi_k[n] = (1 - \alpha)\pi_k[n - 1] + \alpha w_{kn}$$

• How to design  $\alpha(w_{kn}, \pi_k)$  can be investigated in project 1.



## Mean-shift Clustering

- A proper solution to the local min problem is to find *all* local minima.
- Two steps:
  - Mean-shift filter (mode seeking)
  - Clustering



## Kernel density estimate

• For a set of sample points  $\{\mathbf{x}_n\}_1^N$ we define a continuous PDF-estimate







# Kernel density estimate

• For a set of sample points  $\{\mathbf{x}_n\}_1^N$ we define a continuous PDF-estimate

as:  

$$p(\mathbf{x}) = \frac{1}{Nh^d} \sum_{n=1}^N K\left(\frac{\mathbf{x}_n - \mathbf{x}}{h}\right)$$

- K() is a kernel, e.g.  $K(\mathbf{x}) = c \exp\left(-\mathbf{x}^T \mathbf{x}/2\right)$
- h is the kernel scale.



# Mode seeking

- By *modes* of a PDF, we mean the local peaks of the kernel density estimate.
  - These can be found by gradient ascent, starting in each sample.
  - If we use the Epanechnikov kernel,

 $K_E(\mathbf{x}) = \begin{cases} c(1 - \mathbf{x}^T \mathbf{x}) & \text{if } \mathbf{x}^T \mathbf{x} \leq 1 \\ 0 & \text{otherwise.} \end{cases}$ a particularly simple gradient ascent is possible.



### Mean-shift filtering

- Start in each data point,  $\mathbf{m}_n = \mathbf{x}_n$
- Move to position of local average  $\mathbf{m}_n \leftarrow \operatorname{mean}_{\mathbf{x}_n \in S(\mathbf{m}_n)}(\mathbf{x}_n)$
- Repeat step 2 until convergence.





### Mean-shift clustering

 After convergence of the mean-shift filter, all points within a certain distance (e.g. h) are said to constitute one cluster.





### Pose estimation

- Mean-shift can be used for "continuous voting" in pose estimation.
- Each local invariant feature (e.g. SIFT or MSER) will cast a vote (sample point)

$$\mathbf{x} = egin{pmatrix} x_0 & y_0 & lpha & s & arphi & heta & extsf{type} \end{pmatrix}^T$$







#### Mean-shift

Choice of kernel scale affects results





#### Mean-shift

- For the Epanechnikov kernel, the algorithm is quite fast.
- The Gaussian kernel is another popular choice.
- There is also a scale adaptive version of meanshift, that works in a manner similar to EM in each iteration (slower).



# **Generalised Hough Transform**

- Another way to find modes of a PDF is to quantize the parameter space into accumulator cells.
- Each sample then casts a vote in one or several cells.
- This is called the *Generalised Hough Transform* (GHT).



# **Generalised Hough Transform**

Non-iterative ⇒ constant time complexity.





# **Generalised Hough Transform**

 Quantisation can be dealt with by increasing the number of cells, and blurring.





- A similar technique is to use averaging in channel representation.
  - By first quantizing, and then blurring, we are actually introducing aliasing of the PDF.
  - Better to directly sample the kernel density estimate at regularly sampled positions.
  - Density of samples is regulated by the kernel scale.



Channel encoding





Channel encoding
 Channel value



Channel decoding

$$\hat{x} = \operatorname{dec}(\mathbf{x})$$



 A local decoding is necessary in order to decode a multi-valued channel representation.

$$x_{1}=2.5 r_{1}=1 \Rightarrow 0 1 2 3 4 5 6 7 8 9$$

$$x_{2}=7 r_{2}=0.5$$

That is

 $\hat{x}_1 = \operatorname{dec}(x_1 \dots x_3)$   $\hat{x}_2 = \operatorname{dec}(x_6 \dots x_8)$ - Decoding formula depends on the kernel.



## **Channel Clustering**

- Channel encode data points,  $\mathbf{x}_n = \operatorname{enc}(x_n)$
- Average channel vectors  $\bar{\mathbf{x}} = \frac{1}{N}$
- Compute all decodings  $(\hat{x}, \hat{r})$





# **Channel Clustering**

- The decoding step computes *location*, *density*, and *standard deviation* at mode.
- Optimal decoding is expensive, but fast heuristic decodings exist.
- It can be shown [Forssén 04] that averaging in channel representation is equivalent to a regular sampling of a kernel density estimator.



# Summary

- This was a quick overview of clustering, and related techniques.
- The main purpose with learning is to make Computer Vision systems adapt to data.
- The alternative, to manually tune parameters, works for small static problems, but does not scale and cannot adapt to changes.



### Course events this week

- Thursday: Lab1
   Material on the course web page.
   Preparation is necessary to finish on time.
- Friday: Projects start Introductory lecture Assignments into groups (4/5 per group)