Lab course Image Analysis I ST 2023

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Lab 1: Bayesian Classifiers





Content of the 1st Lab

Goal:

- Automatic extraction of features from images for classification task
 - By probabilistic image interpretation → requires models/classifiers

Tasks:

- 1. Implement your methods to extract features from images
- 2. Implement and train Bayesian classifiers to
 - classify real images using extracted features
- 3. Generate synthetic datasets drawn from Gaussian distribution and
 - apply the implemented classifiers on that toy dataset
- 4. Discussion and evaluation (Visually and Quantitively)





- Hand-crafted Features
- Bayesian Classification
- Generative Probabilistic Classifiers:
 - Normal Distribution (Gaussian) Classifier
 - Gaussian Mixture Model Classifier
- Quality Metrics



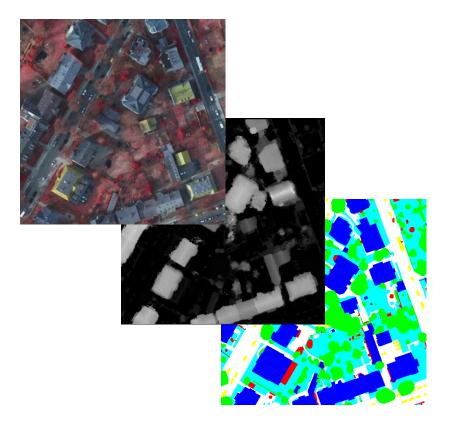
Feature Extraction

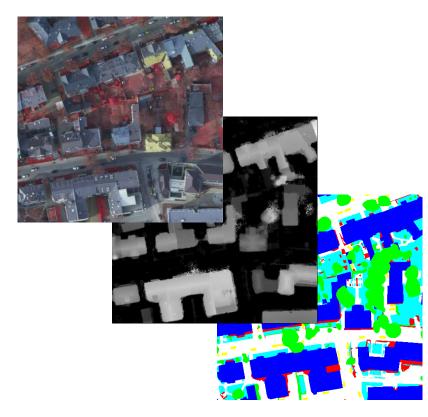
- Two patches (training / testing) with:
 - Channels: Near infrared, red and green (NIR,R,G)
 - Normalized Digital Surface Model (NDSM)
 - Reference labels (per pixel, 5 classes)
 - Classes are: 'STREET', 'HOUSE', 'LOW VEG.', 'HIGH VEG.', 'CAR'
- Additional feature: NDVI = $\frac{NIR-R}{NIR+R}$ (if NIR = R = 0 \rightarrow NDVI = 0) Note that -1 < NDVI > +1
- We will not use class, clutter' → merge with class, low vegetation'
- Normalization of channels required!
 - Shift and scale IR, R and G so that 0 will be mapped to -1.0 and 255 will be mapped to 1.0



Dataset

Training Testing





Objective: Classification Task: USE extracted features, fit a probabilistic classifier

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Classification task: What is the class label *L* for a feature **x**?

MAP-Criterion: Classification based on maximum of the **posterior probability** $p(C \mid \mathbf{x})$

Generative approach:

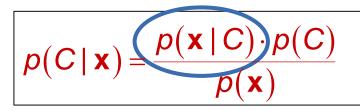
- Often easier to model the causal relation between object type and observed features:
 the observed features are a function of the object type
- $p(C \mid \mathbf{x})$ is modelled indirectly according to the Theorem of Bayes:

$$p(C \mid \mathbf{x}) = \frac{p(\mathbf{x} \mid C) \cdot p(C)}{p(\mathbf{x})}$$

 The theorem of Bayes allows inverse reasoning: derive information about the cause (the class) from the effect (the observed features).



• $p(\mathbf{x}|C)$: Likelihood



- Probability to observe x if it is known to belong to class C
- For each class C^k there is a model for $p(\mathbf{x}|C=C^k)$ that describes the distribution of the features for this class
- Determined from training data, but ?????? HOW ??????
- Non-parametric Models: direct determination of p(x | C) from the training data
- Parametric Models: Based on the assumption of an analytical model for $p(\mathbf{x} \mid C)$ whose parameters are estimated from the training data \rightarrow Machine Learning





• p(C): Prior probability

 $p(C \mid \mathbf{x}) = \frac{p(\mathbf{x} \mid C) \cdot p(C)}{p(\mathbf{x})}$

- Corresponds to bias for the occurrence of C
- If no information is available: Uniform Distribution (e.g for 5 classes: p(C1)=1/5, p(C2)=1/5, p(C3)=1/5, p(C4)= /5, p(C5)=1/5)
 → MAP becomes Maximum Likelihood (ML) criterion
- -p(C) can be determined iteratively:
 - 1) Classification under the assumption of a uniform distribution of the occurrence of the individual classes
 - 2) Determination of p(C) from the relative frequencies of occurrence of the individual classes C^k
 - 3) Classification according to the theorem of Bayes





p(x): Probability of the data

- $p(C \mid \mathbf{x}) = \frac{p(\mathbf{x} \mid C) \cdot p(C)}{p(\mathbf{x})}$
- Equal for all values of C because it does not depend on C
 - \Rightarrow MAP can also be applied without knowing $p(\mathbf{x})$:

$$p(C|\mathbf{x}) \propto p(\mathbf{x}|C) \cdot p(C)$$
 implies that $\max(p(C|\mathbf{x})) = \max(p(\mathbf{x}|C) \cdot p(C))$

- $-p(\mathbf{x})$ ensures that $p(C|\mathbf{x})$ can be interpreted as a probability and can be used as such in further probabilistic processes
- $-p(\mathbf{x})$ can be determined as the marginal distribution of $p(\mathbf{x},C)$: $p(\mathbf{x})$ independent of C

$$\rho(\mathbf{x}) = \sum_{k} \rho(\mathbf{x} \mid C = L^{k}) \cdot \rho(C = L^{k})$$





Summary: Bayesian Classification

• Given:

$$p(C \mid \mathbf{x}) = \frac{p(\mathbf{x} \mid C) \cdot p(C)}{p(\mathbf{x})}$$

- Models for the likelihoods $p(\mathbf{x}|C=L^k)$ of all classes L^k
- Priori probabilities $p(C=L^k)$ of all classes L^k
- ??Question??: A feature vector **x** to be classified
- Wanted: class C_{map} of **x** according to the MAP criterion: Maximize posterior probability $p(C \mid x)$
- Procedure: Compute the posterior and take the class label to be the index of the maximum (ref. slide 9)
 - 1) For all L^k : calculate $p(\mathbf{x}, C=L^k) = p(\mathbf{x}|C=L^k) \cdot p(C=L^k)$
 - 2) Calculate $p(\mathbf{x}) = \sum_{k} p(\mathbf{x} \mid C = L^{k}) \cdot p(C = L^{k})$
 - 3) For all L^k : calculate $p(C=L^k|\mathbf{x}) = p(\mathbf{x}, C=L^k) / p(\mathbf{x})$
 - 4) C_{map} is the label L^k for which $p(C=L^k|x)$ is a maximum



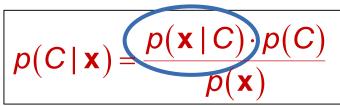


- Hand-crafted Features
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- Generative probabilistic classifiers:
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Generative Probabilistic Classifiers

How to model the likelihood? → Learning from data!



Parametric Methods:

- Analytical model for the probability density $p(\mathbf{x}|C)$ is assumed
- The probability density function $p(\mathbf{x}|C)$ also depends on parameters θ , i.e. $p(\mathbf{x}|C) = p(\mathbf{x}|C,\theta)$
- Parameters are learned from training data
 → Training samples are required for each class C=L^k to determine the parameters θ_k of p(x|C=L^k,θ_k)
- BUT which parametric models do we USE?

(Non-parametric methods are another option, but not covered in this lab)





1. Single Gaussian Model

• Frequent assumption: Multivariate normal distribution

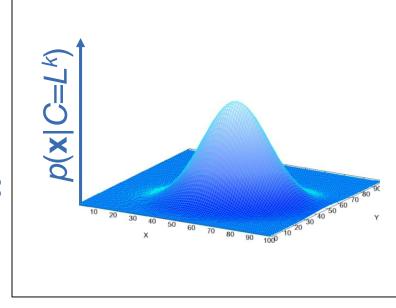
$$\rho(\mathbf{x} \mid C = L^{k}) = \frac{1}{(2\pi)^{\frac{D}{2}} \cdot \|\mathbf{\Sigma}_{k}\|^{\frac{1}{2}}} \cdot e^{-\frac{1}{2}(\mathbf{x} - \mathbf{\mu}_{k})^{T} \cdot \mathbf{\Sigma}_{k}^{-1} \cdot (\mathbf{x} - \mathbf{\mu}_{k})}$$

- Extension of Univariate case, x here is a feature Vector
- Maximum (Log-) Likelihood estimation of the parameters:

$$\Sigma_i \ln p(\mathbf{x}_{ik}|\mathbf{\theta}_k) \rightarrow \max$$



$$\Rightarrow \text{Result for } \Sigma_k : \quad \Sigma_k = \frac{1}{N_k - 1} \cdot \sum_i (\mathbf{x}_{ik} - \mathbf{\mu}_k) \cdot (\mathbf{x}_{ik} - \mathbf{\mu}_k)^T$$



$$\theta_k == (\mu_k \text{ and } \Sigma_k)$$
:

Learned parameters for a Single Gaussian Model

Prerequisite: L^k must only correspond to one cluster in feature space (c./e.g. Slide 48/47)

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2. Gaussian Mixture Model

• In the case of N_j clusters for the class $C = L_k$, every cluster is described by a normal distribution

$$\theta_k == (\pi_{j}, \mu_{kj} \text{ and } \Sigma_{kj})$$
:

The total probability density is obtained from the weighted sum

Learned parameters for a Mixture of Gaussian Model

of the components:

$$p(\mathbf{x} \mid C = L^k) = \sum_{i=1}^{N_j} \pi_j \cdot N(\mathbf{x} \mid \mathbf{\mu}_{kj}, \mathbf{\Sigma}_{kj})$$

with π_i

Mixture coefficient for cluster j, corresponding to the prior probability for j

 μ_{kj} ...

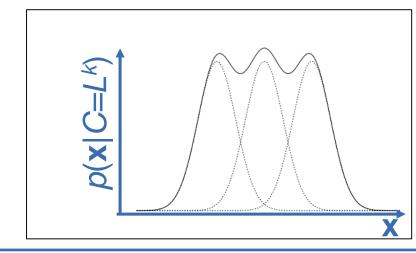
Mean value for cluster *j*

 Σ_{kj} ...

Covariance matrix for cluster *j*

 $N(\mathbf{x}/\mu_{ki}, \Sigma_{ki})$...

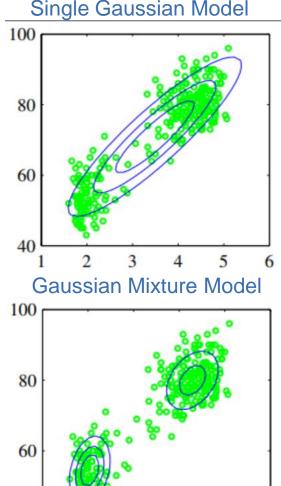
Probability density of the normal distribution for cluster *j*





Gaussian Mixture Model: Example

Single Gaussian Model



- One feature x, two Gaussian components (class index k is omitted):
 - Component 1 (j = 1): $\mu_1 = 3.5$, $\sigma_1 = 1.0$, $\pi_1 = 0.3$

$$\mu_1 = 3.5$$
,

$$\sigma_1 = 1.0,$$

$$\pi_1 = 0.3$$

- Component 2 (j = 2): $\mu_2 = 7.0$, $\sigma_2 = 0.8$, $\sigma_2 = 0.7$

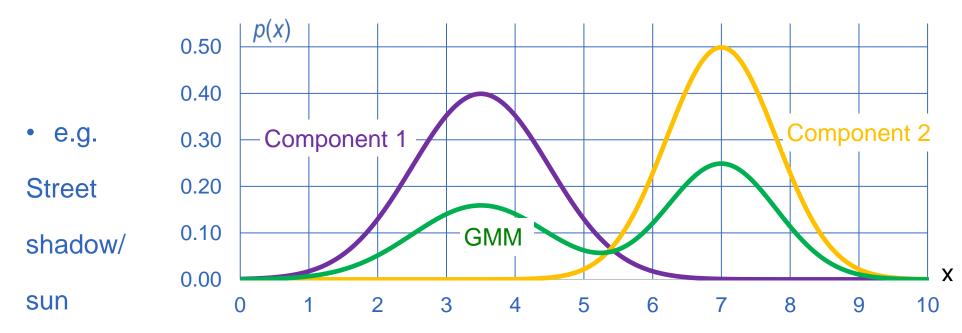
$$\mu_2 = 7.0$$

$$\sigma_2 = 0.8$$

$$\pi_2 = 0.7$$

- GMM: $p(x) = \pi_1 \cdot N(x|\mu_1, \sigma_1) + \pi_2 \cdot N(x|\mu_2, \sigma_2)$

Linear combination of Gaussian distributions



Summary: Gaussian Mixture Model

- Issue: We doesn't know which data points came from which latent component/cluster
- Parameters to be estimated: π_i , μ_{kj} , Σ_{kj} (on set of these params. per cluster j and class k)
- Training of the mixture model requires cluster analysis of the feature space → unsupervised classification
- Method: "Expectation Maximisation" (EM)
 - → see lecture "Unsupervised Classification"
 - E-step: For each data point, compute the prob. of being generated by each component
 - M-step: Adjust the parameters to maximize the likelihood of the data given those params.
- EM requires the number of clusters N_i to be known in advance





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- Quality metrics are used to evaluate the performance of a classifier
- Evaluation requires a reference against which to compare the results of a classifier
- Class-wise metrics:
 - Precision (aka User's Accuracy, Correctness)
 - Recall (aka Producer's Accuracy, Completeness)
 - Intersection over Union (aka Quality)

[python script]

- F1 score
- ...
- Global metrics:
 - Overall Accuracy
 - Mean F1 score
 - ...

```
def calculate_quality_metrics(predicted, true_labels, classes)
   TP = # True Positives
   FP = # False Positives
   FN = # False Negatives

   p = TP/(TP+FP)
   r = TP/(TP+FN)
   f1 = 2*p*r/ (p + r)

   return p, r, f1
```





Terminology:

- True Positives (TP_k):
 - Number of features correctly classified as class k
- False Positives (FP_k):
 - Number of features wrongly classified as class k
- False Negatives (FN_k):
 - Number of features wrongly classified as not class k



Precision:

$$P_k = \frac{TP_k}{TP_k + FP_k}$$

All samples classified as k

Recall:

$$R_k = \frac{TP_k}{TP_k + FN_k}$$

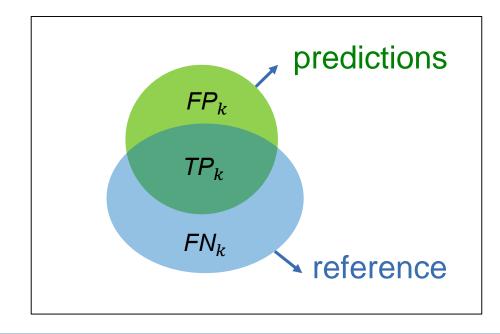
All samples that belong to class *k*

• I.o.U.:

$$IoU_k = \frac{TP_k}{TP_k + FP_k + FN_k}$$

• F1 score:

$$F_k = \frac{2 \cdot P_k \cdot R_k}{P_k + R_k}$$





Overall accuracy:

$$OA = \frac{1}{N} \cdot \sum_{k=1}^{M} TP_k$$

Mean F1 score:

$$mF1 = \frac{1}{M} \cdot \sum_{k=1}^{M} F_k$$

With:

• N: Number of samples

• M: Number of classes

Submission of Results

- Assignment
- → <u>Jupyter Notebook</u> (only digital)
 - Run every cell and save the notebook
 - Use meaningful variable names (refer to IAI_23_Lab_Python_Basics.ipynb)
 - Write comments if required
 - Answer concisely but completely
 - Consider acceptance rules
 - Consider IAI_23_Lab_Technical_Details.pdf





Submission of Results

- Submission deadline: 12. June 2023 before 11:00 am
- There are no resubmissions!
- If you failed one or more labs in the last semester, you have to hand in ALL labs again!

Introduction to the 2nd Lab: 12. June. 2023 at 1:00pm

