Project Assignment 1

TEK5020/9020 - Pattern Recognition

Autumn 2024

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1. Introduction

In this project it is implemented three different classifiing algorithms being

- · Nearest Neighbors (NN)
- Least Squares (LS)
- · Minimum error rate (MER)

A classifier is then trained for each three datasets being splitted in train and test sets as described in the project description.

For the first part each classifier is implemented and trained on the datasets 1 for using all features.

For the code implementaition of the classifiers, refer to the GitHub repository module.

For the second part, for each datasett nearest neighbors classifier is used to find the error rate for all combinations of features as described in section five of the project description.

Then for the combination of features with the lowest error rate, a ranking of the the classifiers is done based on the error rate estimate.

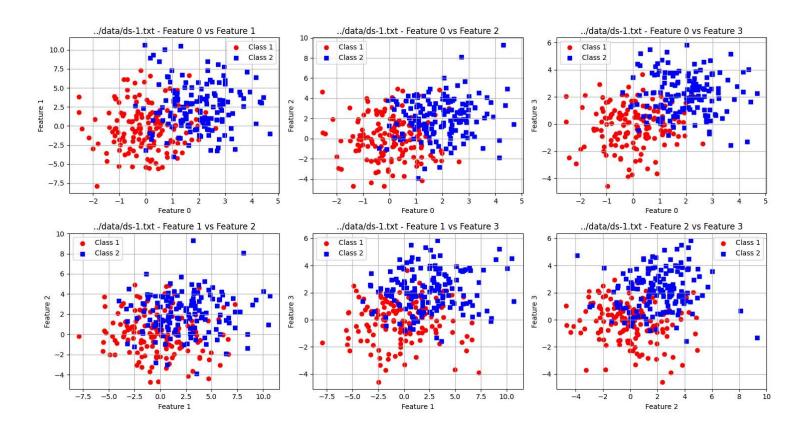
2. Datasets

The datasets used are:

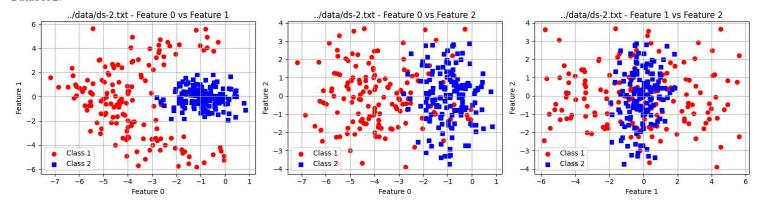
- Dataset 1: 300 objects with 4 features.
- Dataset 2: 300 objects with 3 features.
- Dataset 3: 400 objects with 4 features.

Each dataset is split into training and test sets, with odd-numbered objects in the training set and even-numbered objects in the test set.

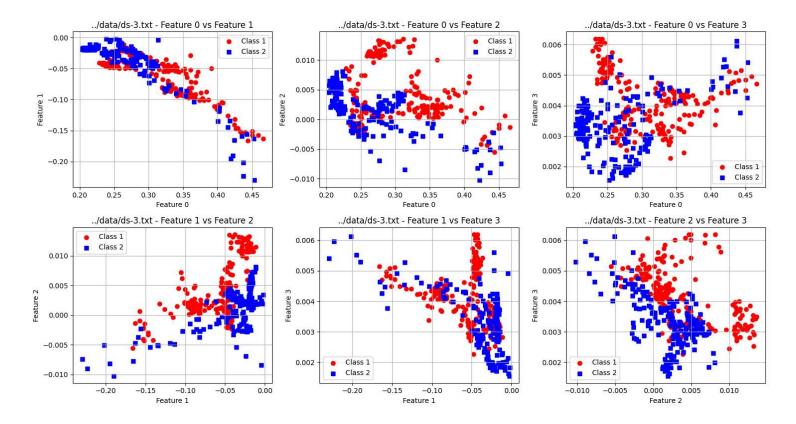
Dataset 1:



Dataset 2:



Dataset 3:



3. Theory

3.1 Minimum Error Rate Classifier

This classifier assigns an object with feature vector \vec{x} to class ω_k if

$$P(\omega_k|ec{x}) = \max_i P(\omega_i|ec{x}), \quad i=1,\ldots,c$$

where c is the number of classes. Assuming normally distributed class-conditional density functions, we obtain the discriminant functions

$$g_i(ec{x}) = ec{x}^T W_i ec{x} + ec{w}_i^T ec{x} + w_{i0}$$

where

$$W_i = -\frac{1}{2}\Sigma_i^{-1} \tag{1}$$

$$\vec{w}_i = \Sigma_i^{-1} \vec{\mu}_i \tag{2}$$

and

$$w_{i0} = -\frac{1}{2}\vec{\mu}_i^T \Sigma_i^{-1} \vec{\mu}_i - \frac{1}{2} \ln |\Sigma_i| + \ln P(\omega_i)$$
(3)

Here, Σ_i and $\vec{\mu}_i$ are the covariance matrix and the mean vector of class ω_i , respectively. These quantities are usually unknown and must be estimated from the training data. Here, the maximum likelihood method should be used to estimate these quantities. This means that the mean vector and covariance matrix of class ω_i are estimated as

$$\hat{ec{\mu}}_i = rac{1}{n_i} \sum_{i=1}^{n_i} ec{x}_{ij}$$

and

$$\hat{\Sigma}_i = rac{1}{n_i} \sum_{i=1}^{n_i} (ec{x}_j - \hat{ec{\mu}}_i) (ec{x}_j - \hat{ec{\mu}}_i)^T$$

where n_i is the number of objects in the training set for class ω_i and \vec{x}_i is the feature vector of object j in class ω_i .

Because we are only considering two-class problems here, we can use

$$g(\vec{x}) = g_1(\vec{x}) - g_2(\vec{x})$$

as the discriminant function. With this, \vec{x} is assigned to class ω_1 if $g(\vec{x}) > 0$ and to class ω_2 otherwise.

As part of the project assignment, you should create a function that trains a minimum error rate classifier as outlined above. You need to create a function that, based on the training objects, calculates the matrices W_i , vectors \vec{w}_i , and constants w_{i0} as given in equations (1), (2), and (3) for i=1,2. Since the prior probabilities $P(\omega_i)$ included in equation (3) are unknown here, they must be estimated based on the number of training objects from each class in the dataset.

The function should be used to construct classifiers for selected feature combinations from all three datasets.

3.2 Least Squares Method

This method yields a linear classifier with a discriminant function given by

$$g(\vec{x}) = \vec{a}^T \vec{y} \tag{4}$$

where $\vec{a} = [a_0, a_1, a_2, \dots, a_n]^T$ is the extended vector of weights and $\vec{y} = [1, x_1, x_2, \dots, x_n]^T$ is the extended vector of features. Let \vec{Y} be a matrix of dimension $n \times (d+1)$, which contains all the extended training vectors stored row-wise, i.e.,

$$ec{Y} = egin{bmatrix} 1 & x_{11} & x_{12} & \cdots & x_{1n} \ 1 & x_{21} & x_{22} & \cdots & x_{2n} \ dots & dots & dots & \ddots & dots \ 1 & x_{n1} & x_{n2} & \cdots & x_{nn} \end{bmatrix}$$

Further, we define an n-dimensional vector $\vec{b}=[b_1,b_2,\ldots,b_n]^T$ where each element b_i is given as $b_i=1$ for \vec{y}_i in class ω_1 and $b_i=-1$ otherwise.

The vector \vec{a} is now chosen such that the length of the error vector $\vec{e} = \vec{Y}\vec{a} - \vec{b}$ is minimized. That is, the cost function:

$$J(\vec{a}) = \vec{e}^T \vec{e} = (\vec{Y}\vec{a} - \vec{b})^T (\vec{Y}\vec{a} - \vec{b})$$

is minimized. This can be done by solving the normal equation

$$\vec{Y}^T \vec{Y} \vec{a} = \vec{Y}^T \vec{b}$$

In other words, you should write a function that, based on the **training objects**, calculates the vector \vec{a} for an arbitrary combination of features, so that the classifier is then given by the discriminant function $g(\vec{x}) = \vec{a}^T \vec{y}$ (equation 4).

3.3 Nearest Neighbor Classifier

This is a computationally simple classifier that simultaneously provides good and reliable classifications. For each object \vec{x} to be classified, the distance to all training objects in the training set is calculated, and \vec{x} is assigned the same class as the nearest training object \vec{x}_k , where the distance is given by

$$\|ec{x}-ec{x}_k\|=\min_i\|ec{x}-ec{x}_i\|$$

4. Training of Classifiers

The classifiers were trained on the training sets and evaluated on the test sets. The evaluation metric used is the error rate estimate.

5. Results

5.1 Feature Combination Rankings

For each dataset, the nearest neighbor classifier was used to estimate the error rate for all combinations of features. The best feature combinations for each dataset are as follows:

Dataset 1:

- Best combination (d=1): Feature 0 with NN error rate 0.24
- Best combination (d=2): Features 0, 3 with NN error rate 0.1667
- Best combination (d=3): Features 0, 1, 3 with NN error rate 0.1

• Best combination (d=4): Features 0, 1, 2, 3 with NN error rate 0,0933

• Dataset 2:

- Best combination (d=1): Feature 0 with NN error rate 0.18
- Best combination (d=2): Features 0, 1 with NN error rate 0.0133
- Best combination (d=3): Features 0, 1, 2 with NN error rate 0.02

• Dataset 3:

- Best combination (d=1): Feature 1 with NN error rate 0.31
- Best combination (d=2): Features 1, 2 with NN error rate 0.095
- Best combination (d=3): Features 1, 2, 3 with NN error rate 0.075
- Best combination (d=4): Features 0, 1, 2, 3 with NN error rate 0.095

5.2 Classifier Rankings

For the best feature combinations within each possible feature dimension, the classifiers were ranked based on their error rates.

Dataset 1:

- o Best classifier (d=1): MER with error rate 0.18
- Best classifier (d=2): MER with error rate 0.1133
- o Best classifier (d=3): LS with error rate 0.093
- o Best classifier (d=4): LS with error rate 0.0733

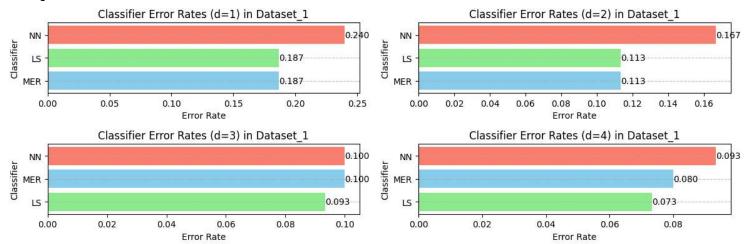
Dataset 2:

- o Best classifier (d=1): MER/LS with error rate 0.107
- o Best classifier (d=2): NN with error rate 0.0133
- Best classifier (d=3): NN with error rate 0.02

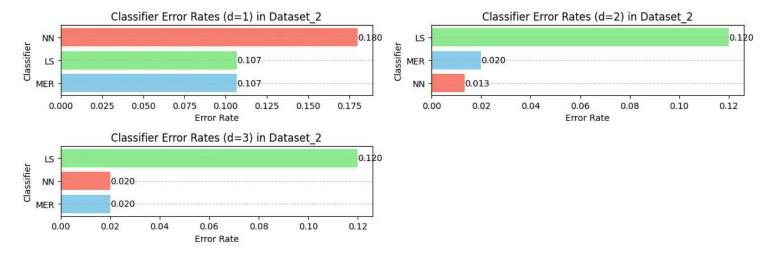
• Dataset 3:

- Best classifier (d=1): MER with error rate 0.225
- Best classifier (d=2): NN with error rate 0.095
- Best classifier (d=3): NN with error rate 0.075
- o Best classifier (d=4): MER with error rate 0.07

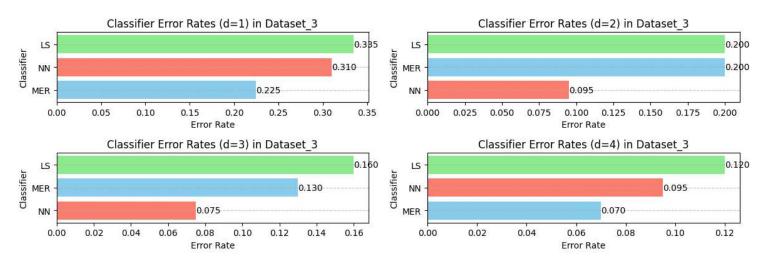
Ranking models for Dataset 1:



Ranking models for Dataset 2:



Ranking models for Dataset 3:



5.3 Discussion

The results demonstrate that the best-performing feature combinations and classifiers vary notably across datasets and feature dimensions. This variability highlights the importance of understanding dataset characteristics and choosing appropriate classifiers accordingly.

For dataset 2, we see that NN performs best for using two features instead of whole dataset that is for the feature combination (0, 1), This outcome can be attributed to the well-defined class boundarie observable in the feature plots of dataset 2. These plots show a distinct separation between Class 1 and Class 2, enabling the NN classifier to accurately categorize samples based on their proximity to training samples. This emphasizes the effectiveness of NN in scenarios where data points from different classes are spatially separated. NN doesn't require linear boundaries and can adapt to complex shapes, making it suitable for dataset 2's using features 0 and 1. adding full dataset to the feature combination, the error rate increases. This is becasue the added features intermingle the classes that the NN classifier struggles to separate. MER does very well for two and three features that is full dataset as it relies on the probabilistic approach to model the class distributions.

In contrast, dataset 1 demonstrates the limitations of the NN classifier. Here, NN consistently has the highest error rates across all feature combinations. The feature plots of dataset 1 reveal a lack of clear boundaries between the two classes, with significant overlap between the data points of Class 1 and Class 2. This overlap leads to higher misclassification rates for NN, as proximity-based classification struggles in cases where class separation is ambiguous. The results underscore that NN classifiers are less effective when classes are not distinctly partitioned in feature space, as seen in dataset 1.

For dataset 3, LS appear to perform worst for all feature combinations classifiers. This is due to the non-linear decision boundaries present in the dataset, which LS struggles to capture effectively. The feature plots of dataset 3 show complex class distributions with overlapping regions, making it challenging for LS to find a linear decision boundary that separates the classes accurately.

We can observe that the class distributions in dataset 3 are sufficiently distinct, allowing MER to estimate the class likelihoods more effectively. As the number of features increases, the error rate for MER decreases significantly.

Across all datasets, Least Square (LS) displays mixed results, performing moderately well in certain cases but failing to match the effectiveness of NN in dataset 2 or the robustness of MER in dataset 3 with higher dimensions. LS may struggle in cases where the underlying decision boundaries are non-linear or irregular, as seen in dataset 1 and dataset 3.

6. Conclusion

1. Why is it reasonable to use the nearest neighbor classifier to find favorable feature combinations?

The nearest neighbor classifier is non-parametric and makes minimal assumptions about the data distribution. This flexibility allows it to effectively evaluate the impact of different feature combinations on classification performance, identifying combinations that result in lower error rates without being constrained by predefined model structures.

2. Why might it be reasonable in a practical application to find a linear or quadratic classifier to replace the nearest neighbor classifier? Linear and quadratic classifiers are computationally more efficient, especially with large datasets, and often require less memory. Additionally, they provide explicit decision boundaries that are easier to interpret and implement in real-time applications. These classifiers can generalize better to unseen data and are less sensitive to noise compared to nearest neighbor classifiers.

3. Why is it unfavorable to use the same dataset both for training and evaluation of a classifier?

Using the same dataset for both training and evaluation can lead to overfitting, where the classifier learns the noise and specific patterns of the training data instead of the underlying generalizable trends. This results in artificially low error rates during evaluation, which do not reflect the classifier's performance on new, unseen data.

4. Why does a linear classifier give poor results for dataset 2?

A linear classifier performs poorly on dataset 2 because the data is poorly linearly separable, meaning that no straight line can effectively separate the classes. The classes overlap and intermingle, making it difficult for a linear classifier to accurately distinguish between them, leading to poor classification performance. On the other hand, the nearest neighbor classifier performed well on this dataset, having a very low error rate compared to the other classifiers. See Ranking models for Dataset 2.