

### Machine Learning



### Activity recognition using performance metrics

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### Abstract

- Establish and evaluate a model, predicting performed activity type, based on a comprehensive set of performance metrics (features)
- Activities classes-labels (categorical type):
  - Running
  - Biking
  - Swimming
  - Hiking
  - Cardio workout
  - Strength training
- Classification algorithms will be used



# Training Data

 Garmin Connect application will be used, exporting training data in json file

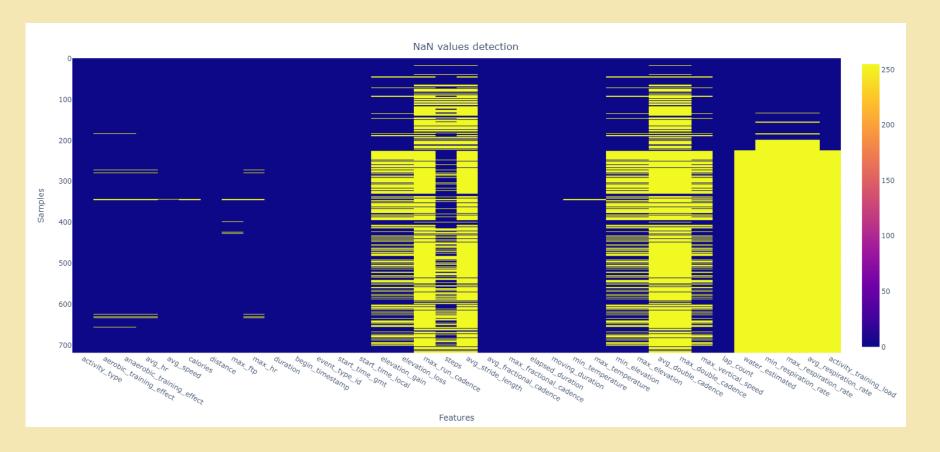
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"anaerobicTrainingEffect": 0.0,
"anaerobicTrainingEffectMessage": "NO_ANAEROBIC_BENEFIT_0",
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"avgHr": 68.0,
"avgSpeed": 0.0,
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        "volume": 0
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"uuidLsb": -7410265071826486760,
"uuidMsb": 1621105547867541023
```

### Data Set

- Data set consisted by about 750 activities will be used as training data
- This data set collected by personal use of Garmin Connect application for over a year
- 10 k-fold stratified cross-validation was used.
- Final train-test scores calculated by the average of each fold score.
- sklearn.model\_selection.GridSearchCV library was used

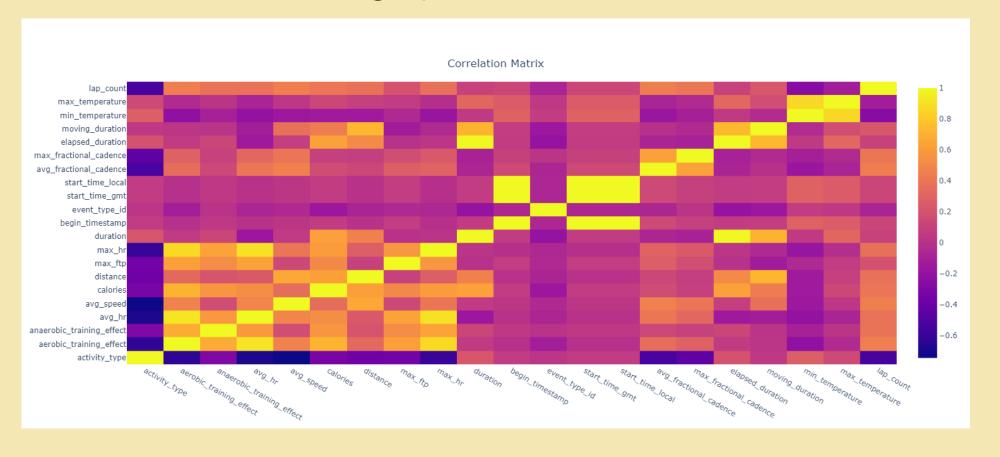
- 34 features were extracted for each example
- NaN value heapmap, correlation matrix and feature boxplots were implemented
- plotly library used

Removed features with multiple NaN values

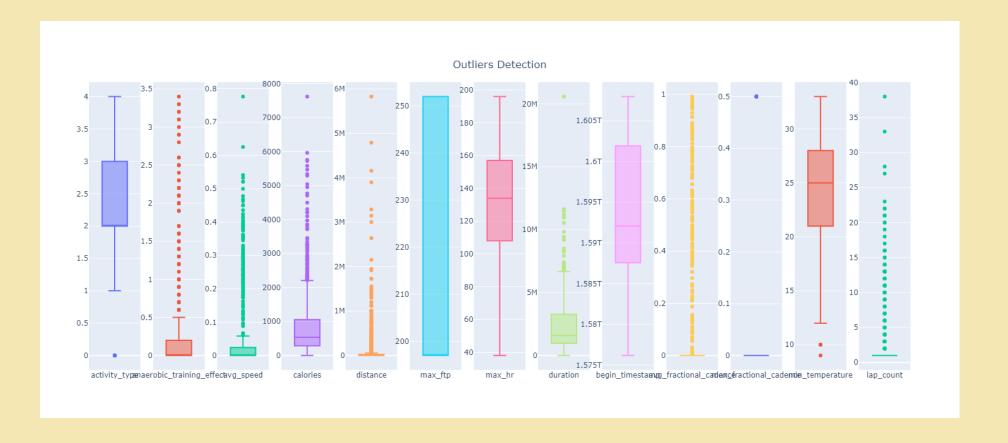




- Removed features with low correlation with activity type
- Removed features highly correlated with others



 Boxplots provide useful information about feature values distribution and outliers detection



# Data Preprocessing

- Removed samples containing NaN values
  - 1. changes range
- 2. preserves the shape of the original distribution
- 3. doesn't reduce the importance of outliers

- ssing library used to normalize the extracted
- preprocessing argorithms:

reduces the effects of outliers

MinMaxScaler: su divides by the ran

- . applied in rows and not in columns
- 2. each row values have a unit norm (if squared and summed, the total would equal 1)
- 3. transform all the features to values between -1 and 1

1. makes the mean of the distribution 0

- 2. about 68% of the values will lie be between -1 and 1
- 3. distorts the relative distances between the feature valuesZ3% VAIUE)

n from each value and then

then dividing by the interquartile

- ween-raindurascaler. Subtracts the mean and then scaling to unit variance
- Normalizer: L2 normalization
- Hybrid: Combination of previous algorithms

# Data Preprocessing

- After data preprocessing the following features were selected:
  - anaerobic\_training\_effect
  - avg\_speed
  - calories
  - distance
  - max\_ftp
  - max hr
  - Duration
  - begin\_timestamp
  - avg\_fractional\_cadence
  - max\_fractional\_cadence
  - min\_temperature
  - lap\_count
- Dataset size: 705 samples

# Classification Algorithms

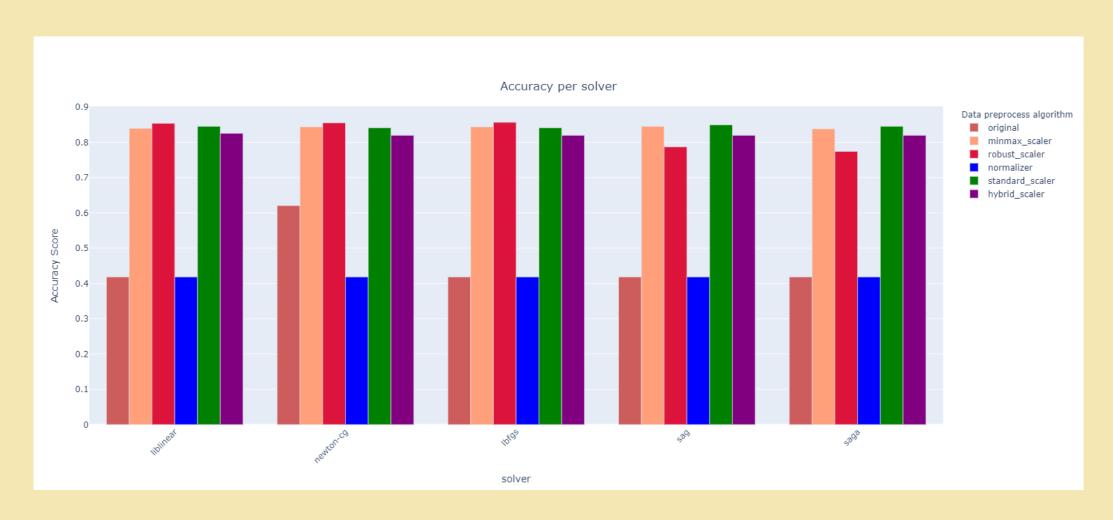
- Logistic Regression
- K-Nearest Neighbors
- Naive Bayes
- Decision Trees
- Support Vector Machine
- (Single-layer) Perceptron

# Logistic Regression

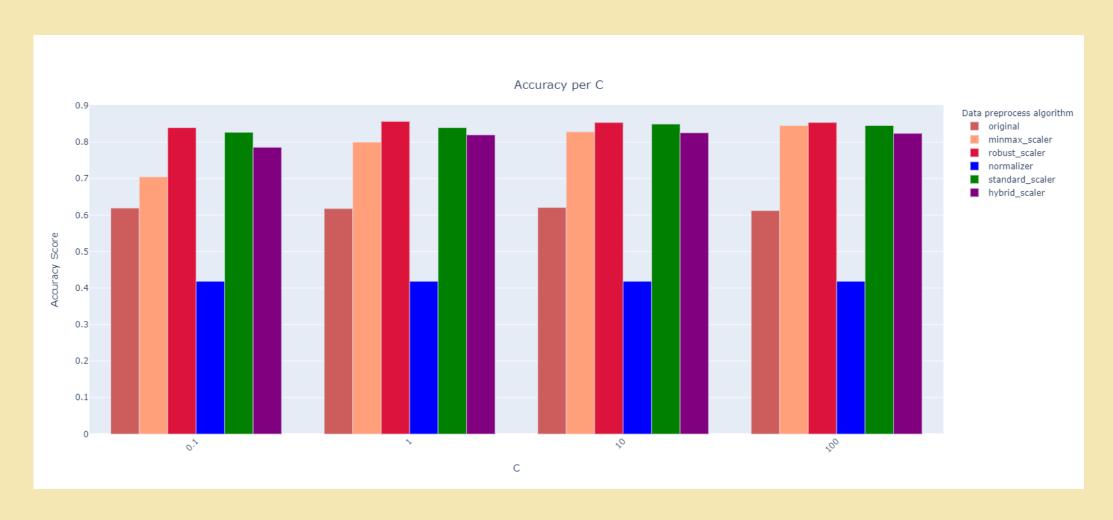
- the probabilities describing the possible outcomes of a single trial are modelled using a logistic function
- several classes involved multiclass logistic regression was preformed
- sklearn.linear\_model library was used
- Algorithm used to find the global likelihood function maximum at Gradient Ascent tric was used for trained model performance evaluation
  - eters:
  - solver: all available solvers compared
  - C: the penalty parameter, which represents misclassification or error term –
     ed C = 0.1, 1, 10, 100

low C -> Higher C ->

# Logistic Regression – Accuracy per solver



# Logistic Regression – Accuracy per C value





# Logistic Regression – Optimal Parameters

• solver: lbfgs

• C: 1.0

data preprocessing policy: robust\_scaler

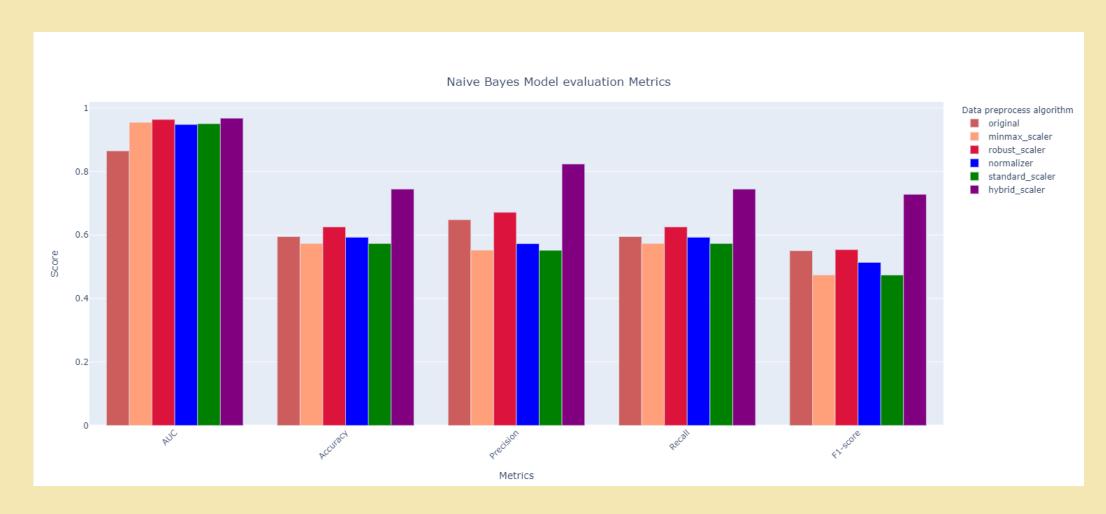
• Accuracy score: 0.8566



# Gaussian Naive Bayes

- based on Bayes' theorem
- assumption of independence between every pair of features -> highly correlated features should be removed
- assumption of Gaussian distribution of joint probability p(x | y)
- sklearn.naive\_bayes library was used
- Gaussian Naive Bayes classifier used
- stratified 10 fold cross-validation was performed
- accuracy, precision, recall, F1-measure metrics were used for trained model performance evaluation

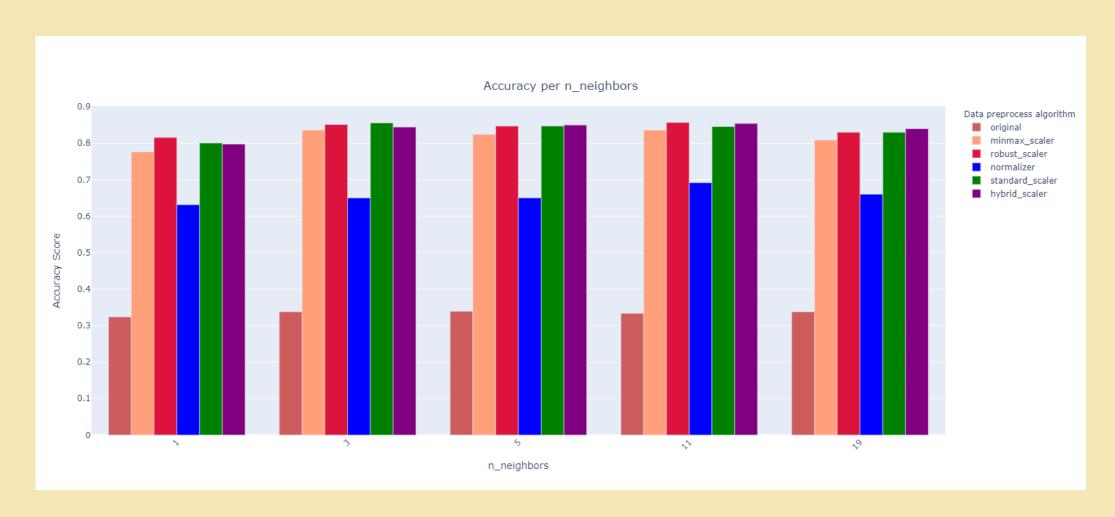
# Gaussian Naive Bayes



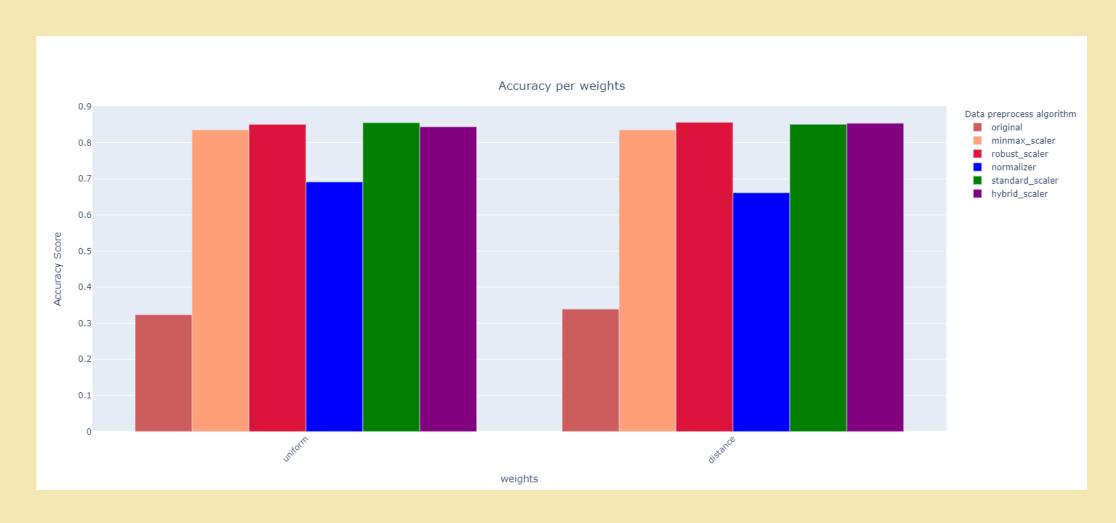
# k-Nearest Neighbors

- classification over k-nearest neighbors majority vote
- sklearn.neighbors library was used
- stratified 10 fold cross-validation was performed
- accuracy metric was used for trained model performance evaluation
- Hyper-parameters:
  - k\_neighbors: defines the number of nearest neighbors, used to classify the unknown point - used: k=1,3,5,11,19
  - weights:
    - uniform -> each neighbor 'vote' within the boundary carries the same weight
    - distance -> closer points will be more heavily weighted toward the decision
  - **metric**: the algorithm used to calculate the distance of neighboring points is chosen from the unknown point - used Euclidean and Manhattan distances

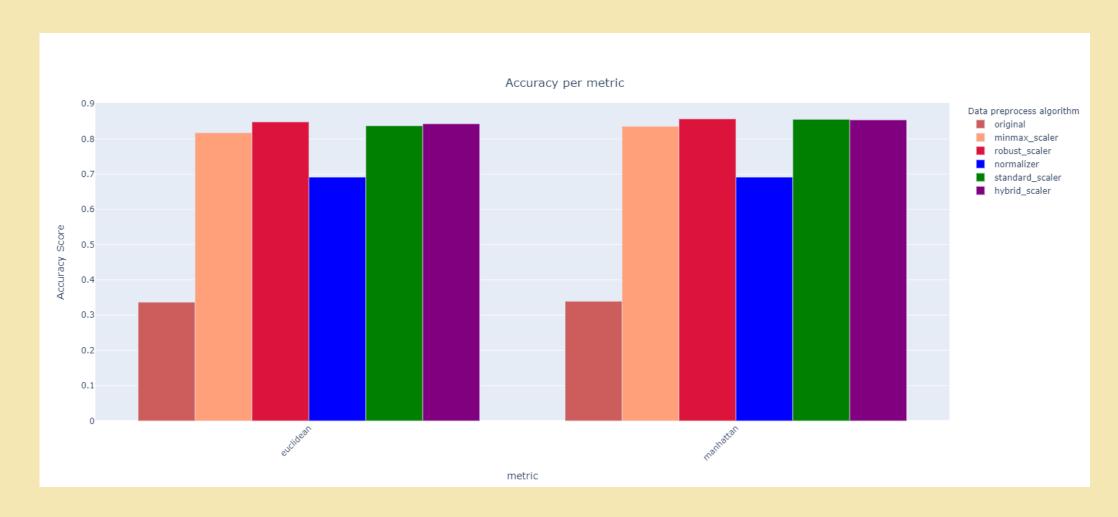
### k-Nearest Neighbors – Accuracy per n\_neighbors



### k-Nearest Neighbors – Accuracy per weights



### k-Nearest Neighbors – Accuracy per metric





# k-Nearest Neighbors – Optimal Parameters

• n\_neighbors: 11

weights: distance

metric: manhattan

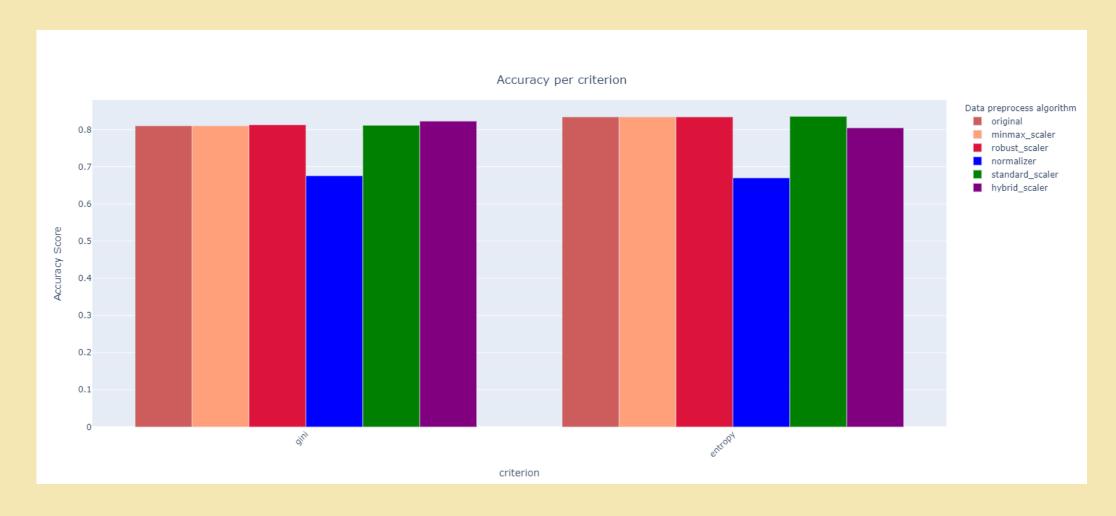
data preprocessing policy: robust\_scaler

• Accuracy score: 0.8568

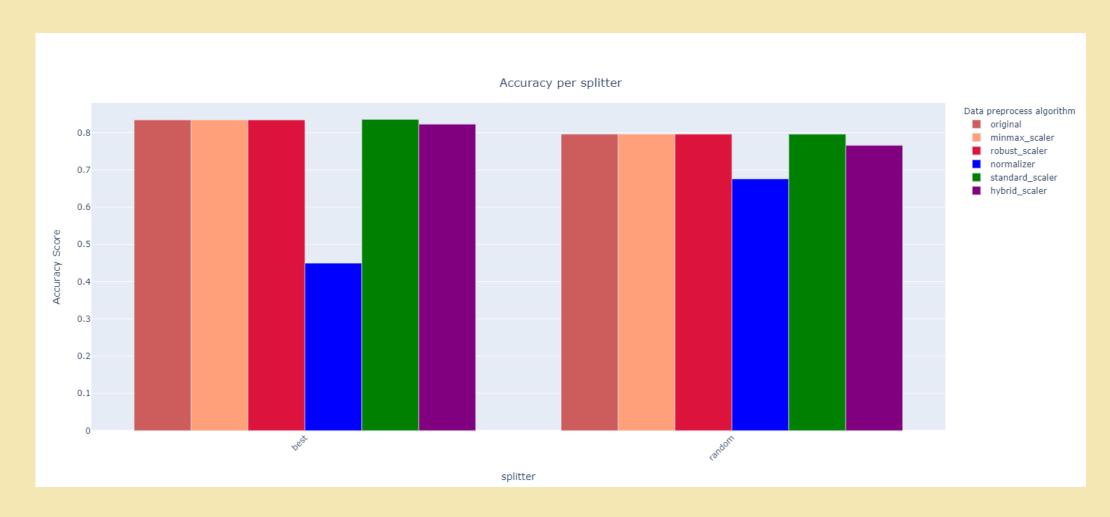
# **Decision Trees**

- Classification model built in the form of a tree structure
- **sklearn.tree** library was used
- stratified 10 fold cross-validation was performed
- accuracy metric was used for trained model performance evaluation
- Hyper-parameters:
  - criterion: defines the function to measure the information gain
     used: gini, entropy
  - splitter: the strategy used to choose the split at each node
    - best -> best split among all is chosen
    - random -> best split among a random subset is chosen
  - max\_depth: the maximum depth of the tree, if None -> nodes are expanded until all leaves are pure - used None, 1, 5, 10, 15, 20, 25

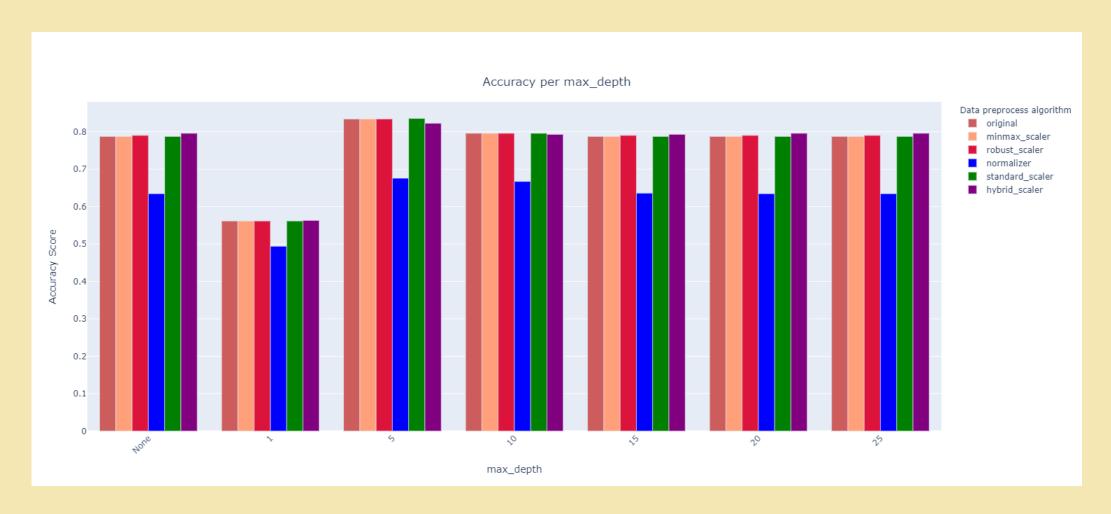
### Decision Trees – Accuracy per criterion



### Decision Trees – Accuracy per splitter



### Decision Trees – Accuracy per max\_depth





# Decision Trees – Optimal Parameters

criterion: entropy

• max\_depth: 5

• splitter: best

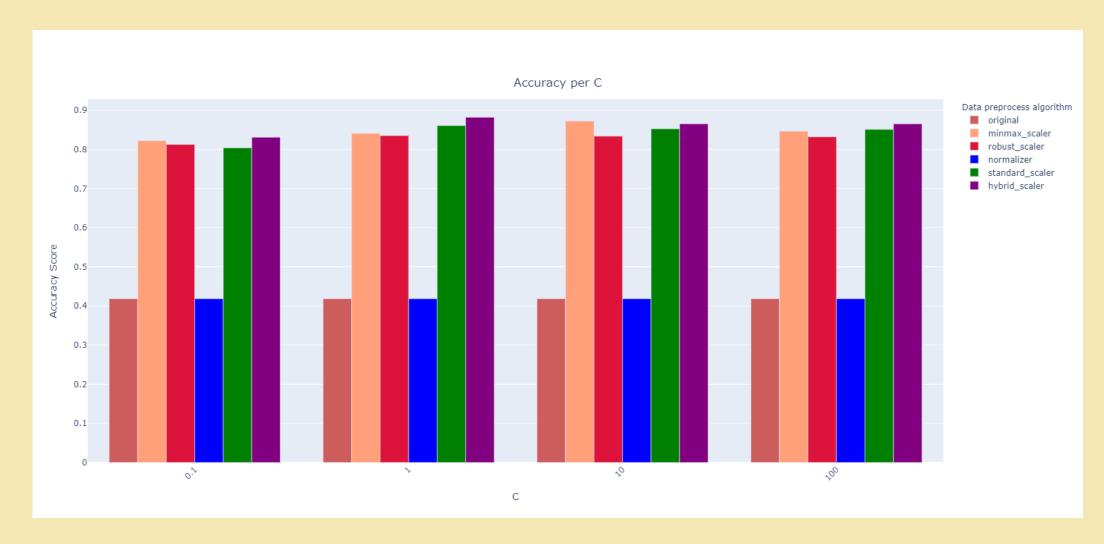
data preprocessing policy: standard\_scaler

• Accuracy score: 0.8358

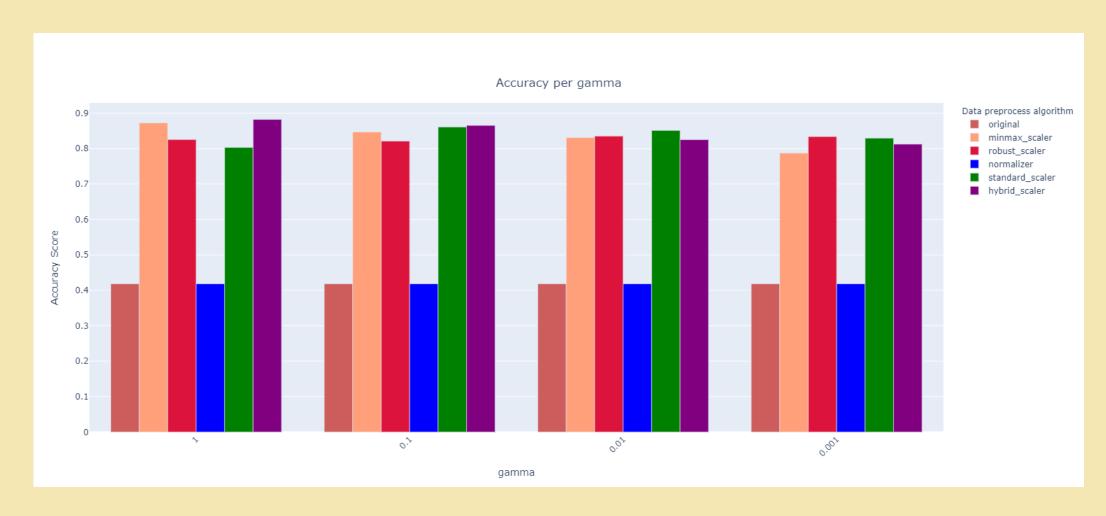
# Support Vector Machine (SVM)

- Training data represented as points in space
- Classification achieved by defining a decision surface that separates the classes
- Goal: maximize the margin between the data points of the two classes
- sklearn.svm.SVC library was used
- stratified 10 fold cross-validation was performed
- accuracy metric was used for trained model performance evaluation
- Hyper-parameters:
  - kernels: the function used to transform low dimensional input space into a higher-dimensional space
    - poly -> applying the polynomial combination of all the existing features (e.g.  $x^2$ )
    - rbf -> generates new features by measuring the distance between all other dots to a specific dot/dots centers
    - sigmoid -> sigmoid function is used to transform features
  - **C**: the penalty parameter, which represents misclassification or error term used C = 0.1, 1, 10, 100
  - **gamma**: defines how far influences the calculation of plausible line of separation used g = 1, 0.1, 0.01, 0.001

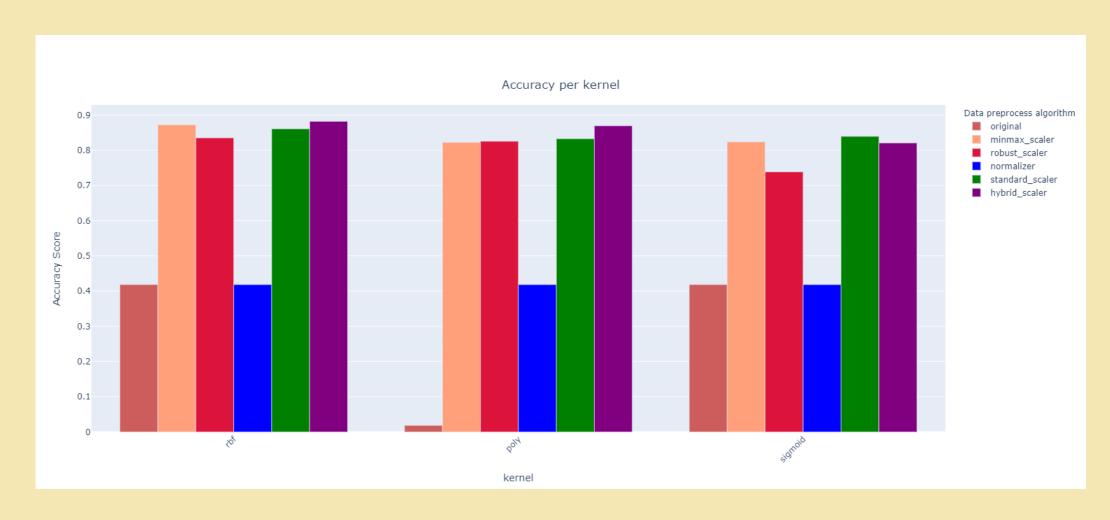
# SVM – Accuracy per C



# SVM – Accuracy per gamma



# SVM – Accuracy per kernel





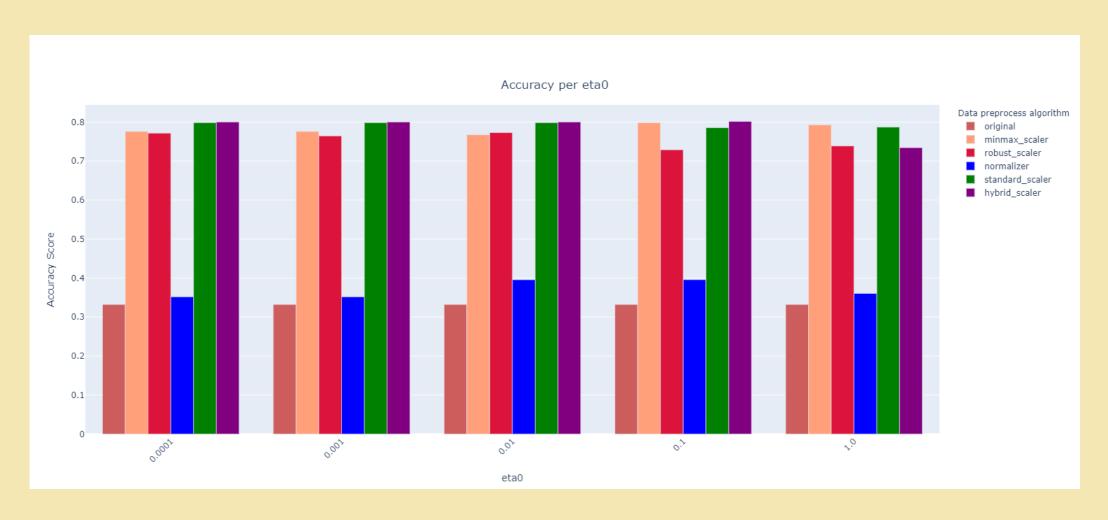
# SVM – Optimal Parameters

- C: 1
- gamma: 1
- kernel: rbf
- data preprocessing policy: hybrid\_scaler
- Accuracy score: 0.8822

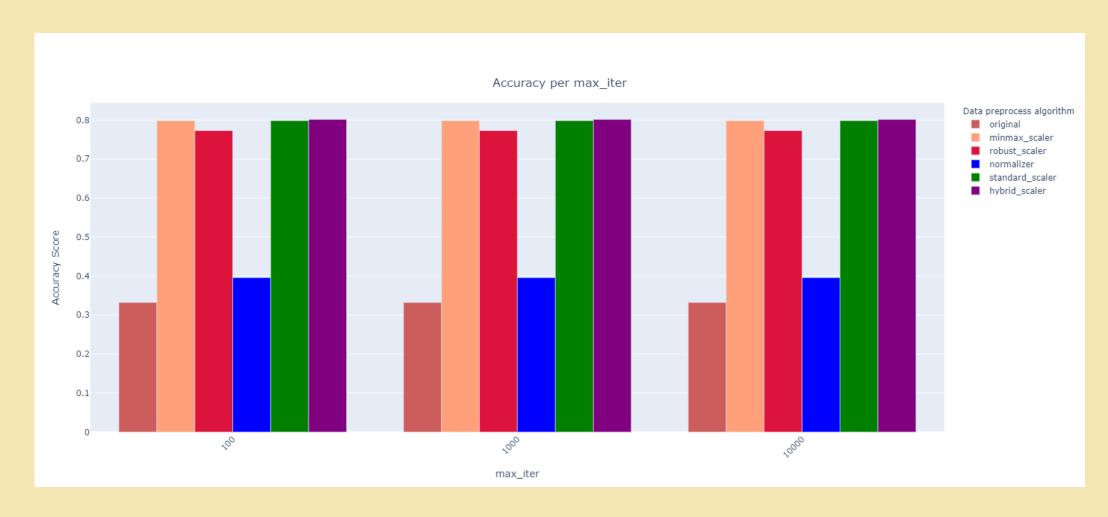
# (Single-layer) Perceptron

- single layer Perceptron is the basic unit of a neural network
- binary classification machine learning algorithm
- multi-class classification can be implemented by one-vs-all policy
- transforms the original feature space by multiplying feature vectors with the respective weights
- aims to define a decision hyperplane where the classes are fully separated (using an activation function -> sigmoid)
- is actually an SVM classifier without the soft-margin concept
- sklearn.linear\_model library was used
- stratified 10 fold cross-validation was performed
- accuracy metric was used for trained model performance evaluation
- Hyper-parameters:
  - learning rate (eta0):
    - high learning rate -> fast training ~ lower performance,
    - low learning rate -> slow training ~ better performance,
    - used 0.0001, 0.001, 0.01, 0.1, 1
  - max iterations: the number of epochs used to train the model used 100, 1000, 10000

### Perceptron – Accuracy per learning rate



### Perceptron – Accuracy per max iteration





# Perceptron – Optimal Parameters

• eta0: 0.1

• max\_iter: 100

data preprocessing policy: hybrid\_scaler

• Accuracy score: 0.8015

# Classifiers Comparison

 Optimal parameters and best data preprocess policy chosen per classifier

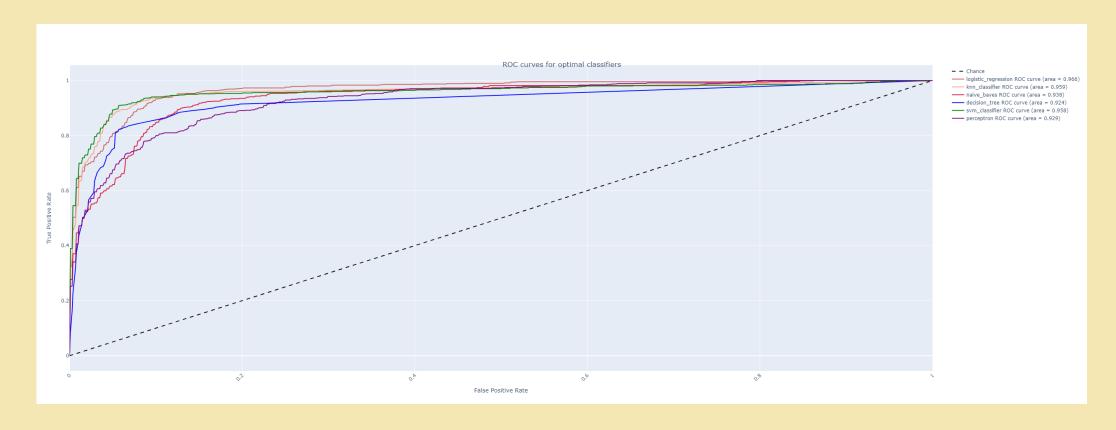
Implemented ROC curves per classifier

 Accuracy, Precision, Recall, F1-score and AUC scores compared

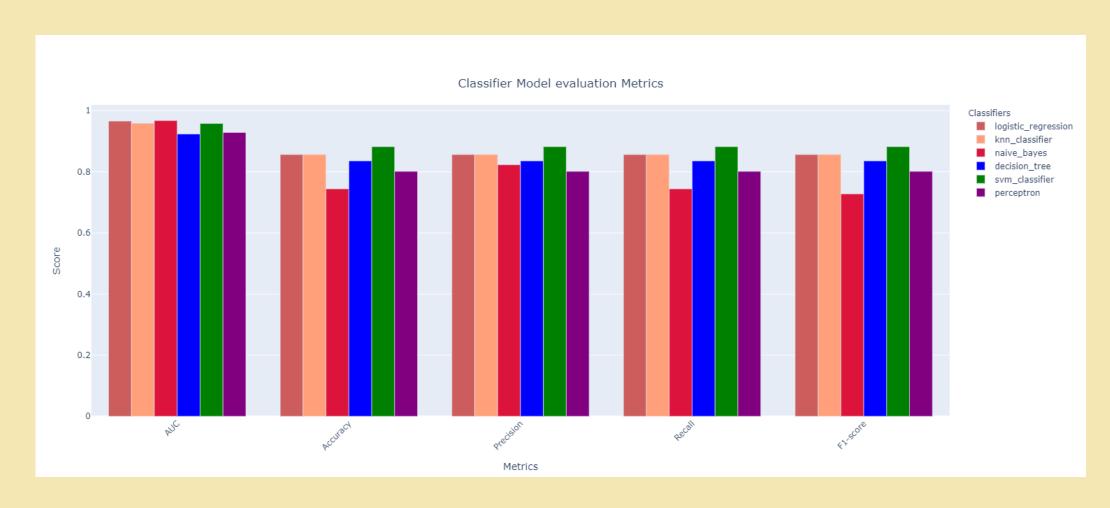
Model Fit time also compared



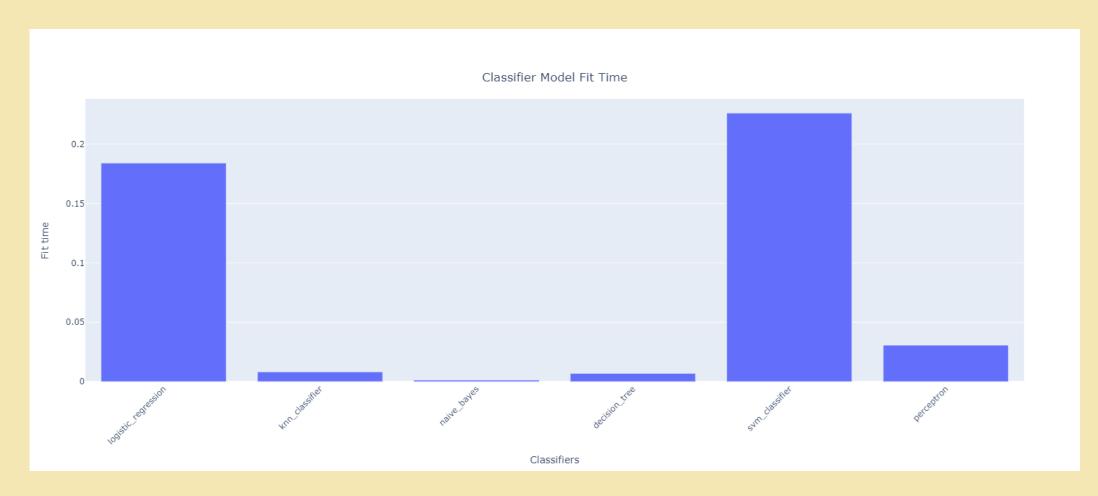
# Optimal Classifiers ROC curves



### Optimal Classifiers Evaluation Metrics



# Optimal Classifiers Model fit time



# Conclusion

- Overall best classifier (based on accuracy metrics):
   SVM
- Best hyper-parameters:
  - C: 1
  - gamma: 1
  - kernel: rbf
- Fastest model training: Gaussian Naïve Bayes
- Best tradeoff (accuracy~training time): kNN classifier