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May 30, 2019

Predicting Presence of Heart Disease using Machine Learning

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Problem Description

According to the World Health organization, coronary heart disease is the undisputed global cause of death (2018). Based on the most recent global cause of death release from the World Health Organization; 10 million deaths were due to heart disease in 2016. Heart disease has remained the leading cause of death in the world for the past 18 years. The goal of this experiment is to use the Heart Disease dataset from the Machine Learning repository accessed through driven data to train models for heart disease prediction. We are trying to identify one or a combination of models that would be able to predict heart disease and help save lives.

Dataset

We will be building the heart disease classification models using the heart disease data from the Cleveland Clinic. The dataset contains data from 270 people with datapoints recorded for each of individual across 13 attributes or features pertaining to heart disease. Below are the features of the dataset and the parameters that define them.

1. age : age in years

2. sex : 1=male, 0= Female

3. chest pain type : chest pain type (4 types)

4. blood_pressure : resting blood pressure mm/hg5. cholesterol : serum cholesterol in mg/Dl

6. blood sugar : fasting blood sugar > 120 mg/Dl (1=true, 2=false)

7. rest_ecg : resting electrocardiographic results

8. max_heart_rate : maximum heart rate achieved

9. exercise induced angina : 1=yes, 0=no

10. st depression : ST depression induced by exercise relative to rest

11. st slope : Slope of Peak exercise ST segment

12. num major vessels : colored by fluoroscopy (0-3)

13. thalassemia : 3=normal, 6= fixed defect, 7=reversible defect

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May 30, 2019

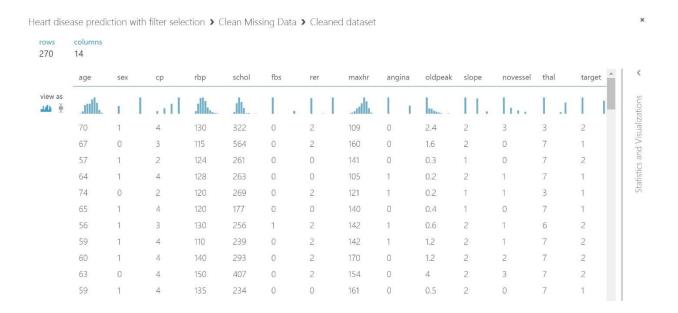
The Experiment – Background:

For this experiment, the dataset will be running through five different algorithms, all of which would be analyzed for accuracy, precision and recall; a recommendation of most predictively potent model(s) for this particular case would be provided based on ensuing results. The five algorithms we will using to predict heart disease are: Two Class Boosted Decision Tree, Two Class Decision Forest, Two Class Naïve Bayes Model, Support Vector Machine and Two Class Logistic Regression.

Cleaning the Data

Before running our algorithms, we needed to make sure the data was clean and free of missing values. After cleaning the data, a snapshot statistical summary of datapoints was obtained. The purpose of this exercise is to give us a standard range for applicable features. For example, with age, we wanted to look at the age distribution of this dataset as it correlates to the real world because heart disease is not just a problem for older adults. Based on our data summary we could tell that individuals in our data set had minimum age of 29 and maximum age of 77 with mean age of 54 and std deviation of 7.5 years. We found the age distribution as plausible and representative of the real world. We applied this same logic to the "sex" feature. With 0 being female and 1 being male, a mean of .67 tells us that the data is more skewed towards men than women. These considerations have been noted and will be made available to the beneficiaries of this experiment. Shown below are graphics of the cleaned data and its statistical summary.

Cleaned Data Set



Cloud Project - Azure ML Studio

May 30, 2019

Data Summary

Heart disease prediction with filter selection > Summarize Data > Results dataset rows columns 14 23 Unique Missing Mean Feature Count Value Value Min Max Mean Deviation Count Count view as 14b ± 270 0 29 77 54.433333 7.505185 age 41 0 0 1 270 0.677778 0.43679 270 4 0 1 4 3.174074 0.789218 ср 94 131.344444 rbp 270 47 0 200 13.829959 schol 270 144 0 126 564 249.659259 39.024088 fbs 2 0 0 1 0.148148 0.252401 270 rer 270 3 0 0 2 1.022222 0.992263 maxhr 90 0 71 202 149.677778 18.71572 0 0 1 0.32963 0.441948 angina 2

Feature Selection

oldpeak

270

39

For this experiment we chose to focus on the 10 most important features within our dataset. It is our goal to prioritize the features that are most linked to heart disease. We want to know which of the list of features tell us the most about the dependent variable; the "target"; identifying presence of heart disease or not. The importance of this feature filter for us is that it translates to statistical significance. We want to load our algorithms with significant data in order to make impactful deductions and increase the potency of our models.

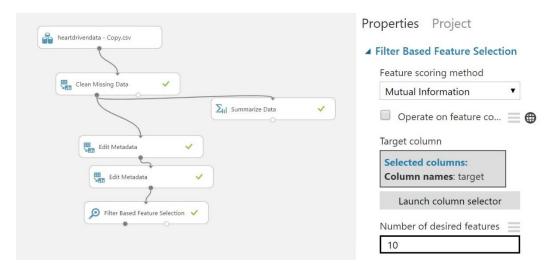
0

0

6.2

1.05

0.918889

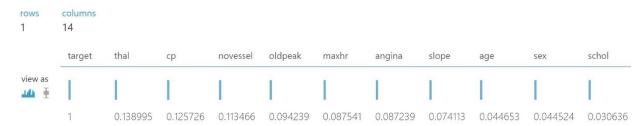


Cloud Project - Azure ML Studio

May 30, 2019

We selected mutual information as feature scoring method because we are dealing with a lack of numeric uniformity in how features are quantified. Mutual Information measures how much information the presence or absence of any feature contributes to making the correct classification decision, in this case; is heart disease present or not ("target" column). When we ran the filter feature selection the results were as follows:

Heart disease prediction with filter selection > Filter Based Feature Selection > Features



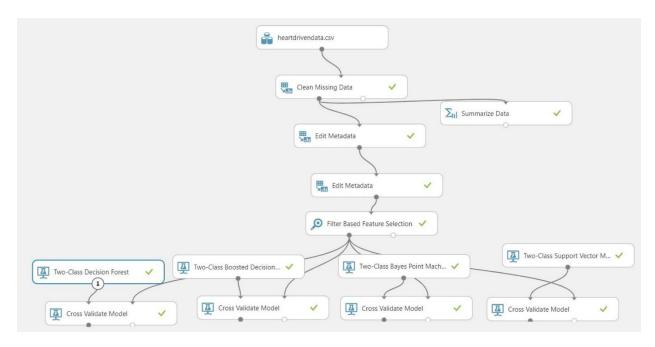
Thalassemia at about 14% was our most correlated feature to heart disease with Cholesterol levels being the least out of the ten features selected at about 3%. The rest of the features; which are fasting blood sugar, resting ecg and resting blood pressure all had a correlation factor below 2% when it comes to predicting the predicting the presence of heart disease.

Cross Validation Technique

Since our goal of this experiment is predictability of heart disease given a combination of features, we chose to not individually score and train the models but rather cross validate them. Cross Validation, we believe is ideal for this experiment because it helps mitigate the problem of overfitting (a situation whereby the models are trained to learn about the actual dataset so closely the models "memorize" and mimic the original dataset). Overfitting would be catastrophic to our efforts to solve a global problem that is heart disease. We cross-validated all five models to ensure that our models do not learn about our data to "memorize" it but rather to "understand" the data so they can predict results given new information. We believe that with cross validation technique, the predictive power of our models has been optimized, giving us more confidence about the resulting accuracy and precision that we obtain. Below is a graphic showing cross-validation of the models. Since we have a small dataset and we could end up getting sufficiently big differences in quality or different optimal parameters between folds. We choose k=10, as a general rule; empirically k=10 tends to yield test error estimates that suffer neither from excessively high bias nor high variance.

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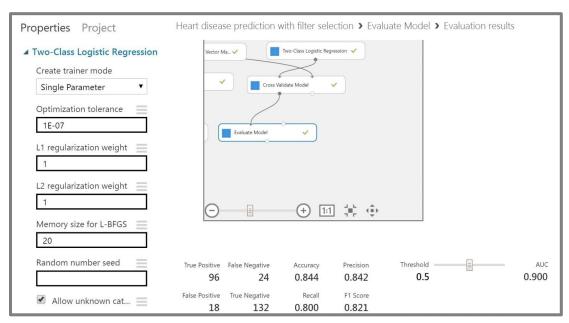
May 30, 2019



Models and Model Evaluations

Two Class Logistic Regression:

The first Machine Learning algorithm we ran was the two-class logistic regression because it is the go-to method for predicting binary classification outcomes. Our ultimate goal is to classify people under two categories of "heart disease not present" or "heart disease present" with both classes represented by a "1" or "2" respectively. Below are the properties that we ran the model with and the ensuing evaluation results.



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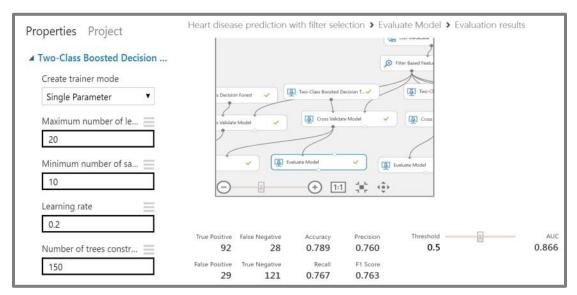
May 30, 2019

We knew the independent parameter that we are testing against (target column) so we chose single parameter logistic regression and left memory size at 20 because the data set is not large enough to require longer training for optimization.

Overall this model had an accuracy and precision level of about 84% with a recall level of 80%. Overall this model will be able to predict heart disease with an accuracy of 84%. This is considered significant for us and viable for future application.

Two Class Boosted Decision Tree:

The decision tree model predicts the value of a target variable based on several input variables. Since we have a target variable we are trying to predict, we chose this algorithm as our second stop. We ran it with 100 trees and then 150 trees at a learning rate of 0.2 with a single parameter as shown below. Note, 50 extra trees on the second iteration had negligible effect on the results. By increasing the number of trees, the algorithm took a few seconds longer to run. In principle this helps reduce variance and increase accuracy. Given the size of the data set the results were not as impressive as the logistic regression model as shown below:



This model had an accuracy score of 78.9% and Precision 76% with a recall of 76%. These numbers are satisfactory but not impressive by our standards. Whether the boosted decision is recommended for future use with this dataset will depend on the results of remaining three models.

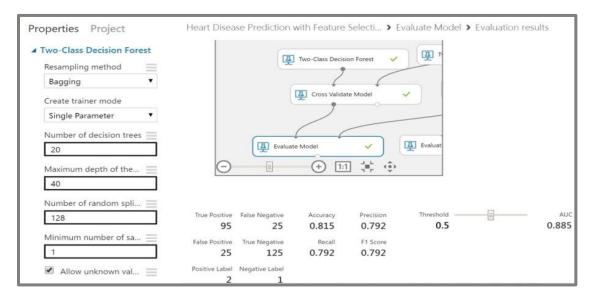
Two Class Decision Forest:

The third algorithm we ran was a two-class decision forest. By using a decision forest which is, in essence an aggregation of multiple decision trees. We hope to see a better accuracy compared to the boosted decision tree. To further solidify the model accuracy, we increased the number to trees to twenty (20) and chose to run the algorithm with the "Bagging" method.

Cloud Project - Azure ML Studio

May 30, 2019

Bagging allows for each tree to be grown from a new sample randomly generated from the original dataset. Bagging enables the model to accommodate variables with wide ranges and varied distributions. The properties of the algorithm and its performance were as follows:



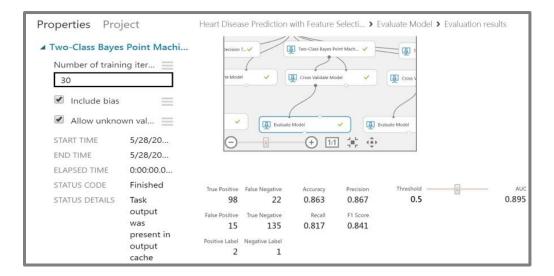
The decision forest had an accuracy score of 81.5% and Precision 79.2% with a recall of 79.2%. The numbers from the decision forest are higher than the decision tree as we envisioned, on average, accuracy, precision and recall all increased by 3%. Given the size of the heart dataset, we saw minimal increase in accuracy and a longer training time for the algorithm when we increased the number of trees.

Two Class Naïve Bayes Model:

We chose to run the two class Naïve Bayes model because we believe the independence assumption holds in terms of our features. While all the futures could contribute to heart disease they do not directly cause or depend on each other e.g age or sex doesn't lead to high cholesterol and vice-versa. This is just one example among 15 others that could be drawn from the dataset. We ran the Naïve Bayes algorithm with the default 30 iterations. We found 30 iterations to be enough because of the rather small dataset and efficient model training time. Through our feature selection we recorded that Thalassemia had the highest correlation factor to heart disease at about 14%. Since the correlation of the features was low with regards to the target we did not deem it necessary to change model properties. Shown below are the properties of the Naïve Bayes model and resulting performance.

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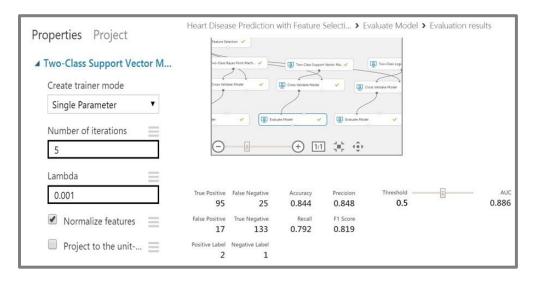
May 30, 2019



The two-class Naïve Bayes algorithm resulted in 86% accuracy and about 87% precision. Higher than any other model that we trained and evaluated. Along with the Logistic Regression model, both scored a recall of about 80%. The Naïve Bayes model was the best performing algorithm in the experiment.

Two-Class Locally Deep Support Vector Machine:

We chose to run the Support Vector Machine (SVM) model because we wanted to see how this model would perform compared to decision trees, decision forest and logical regression models. SVM is capable of doing both classification and regression but tends to take a longer time to run. SVMs excel at identifying complex boundaries, we hoped that this algorithm would set better boundaries within the parameters of the dataset and predict directly what factors lead to heart disease. To further optimize the boundary setting power of the SVM, we ran the algorithm with 5 iterations. The model took longer to run but resulted in better accuracy than all the decision tree and regression models. Shown below are the SVM algorithm properties and performance.



Cloud Project – Azure ML Studio

May 30, 2019

The Support Vector Machine was the second most accurate algorithm in the experiment behind Naïve Bayes model.

Model Analysis and Recommendations

The 'Add Rows' module was used to combine all the accuracy rates of the predictions of all the models we used in the project. It appends a set of rows from an input dataset to the end of another dataset. The results in the 'Evaluate Model' Module was connected to the 'Add Rows' module to create a column-row result with all the accuracy rates from the models

Below is table summarizing the performance of the 5 models in terms of accuracy, precision recall and f-1 score.

Model (Two-Class)	ACCURACY	PRECISON	RECALL	F1 SCORE
Logistic Regression	0.844	0.842	0.800	0.821
Decision Tree (Boosted)	0.789	0.760	0.767	0.763
Decision Forest	0.815	0.792	0.792	0.792
Naive Bayes	0.863	0.867	0.817	0.841
Support Vector Machine	0.844	0.848	0.792	0.819

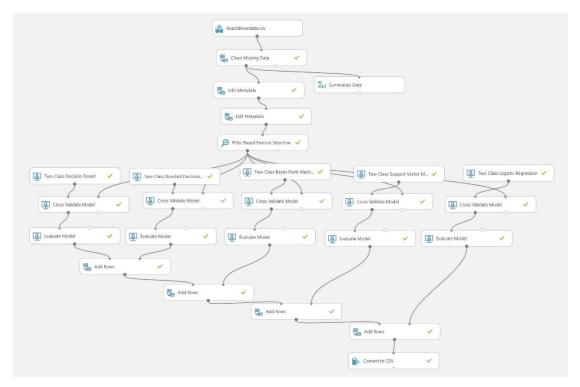
For this experiment we have chosen to evaluate the performance of our models based on their f-1 score. We believe that the F-1 score, which is a weighted average between precision and recall is more useful than accuracy, precision or recall if individually used when analyzing model performance.

Our dataset is unevenly distributed. We are dealing with identifying heart disease in people, the problem that we are trying to solve directly involves human life. The cost of having false positives and false negatives in this experiment needs to be mitigated as much as possible. The nature of this real world problem we are trying to solve makes it almost impossible to solely just look at model accuracy. For this reason, we believe the F1 score gives us a better judgement of model performance. Based on this premise, we believe that the Naïve Bayes model with an F1 score of 0.819 performed the best in our experiment, followed by Support Vector machine with a score of 0.891. We would like to recommend these two models for this data because the other remaining models all fall under the .800 mark or 80th percentile in terms of F1 Score. All models scored relatively high but given the size of the dataset, we are choosing to recommend the two most proficient models which are Naïve Bayes and Support Vector Machine.

Cloud Project - Azure ML Studio

May 30, 2019

Appendix I – Screenshot of Entire Model



Appendix II - Dashboard of Performances from our Azure ML Algorithms

