Udacity Machine Learning Engineer Nanodegree program Final Project – Predicting heart disease

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I Definition

Project overview

Heart diseases are a huge problem in our modern society. It accounts for one third of the total of death cause in the U.S., yet Google and media coverage is about only 2-3%.

In order to make society more careful about this matter, we could use machine learning to predict whether people are prone to develop heart disease in the future and alert them in order to prevent deaths. This Dataset was used into the article "Heart Disease Prediction Using Machine learning and Data Mining Technique".

Problem statement

The goal is to create a prediction model that predicts with a minimum of 70% of accuracy whether a given person with determined attributes will develop a heart disease or not. The 14 attributes contained in the dataset are well defined and all measureable with a good degree of precision. Since the target variable is known, we will use supervised machine learning techniques to solve the problem.

Evaluation metrics

The target feature - 0 = no presence of heart disease; 1 = presence of heart disease is very well balanced (138 for "no presence of heart disease" and 165 for "presence of heart disease"), hence the evaluation metric is the accuracy score. As discussed in the section *II Analysis* – data exploration, the other features are not well balanced, so just in case, we will also explore the confusion matrix results.

II Analysis

Data and visualization exploration

In 1988 four institutes collected data from 303 people and made a dataset of 76 attributes about heart disease. The institutes are listed below

- 1. Hungarian Institute of Cardiology. Budapest: Andras Janosi, M.D.
- 2. University Hospital, Zurich, Switzerland: William Steinbrunn, M.D.
- 3. University Hospital, Basel, Switzerland: Matthias Pfisterer, M.D.
- 4. V.A. Medical Center, Long Beach and Cleveland Clinic Foundation: Robert Detrano, M.D., Ph.D.

Historically, from the 76 attributes, only a subset of 14 are been used and considered relevant for hear disease prediction studies, thus this capstone project will use the 14 features subset as the dataset. The target attribute refers to either 0 (no heart disease presence) or 1 (heart disease presence). This dataset was cited in this article "Heart Disease Prediction Using Machine learning and Data Mining Technique" but was collected from Kaggle.

All the features are shown below

- 1. Age (age): numerical
- 2. **Sex (sex):** categorical -0 = male and 1 = female
- 3. Chest pain type (cp): categorical from 1 = low pain to 4 = very painful
- 4. Resting blood pressure in mmHg (trestbps): numerical
- 5. **Serum cholestoral in mg/dl (chol):** numerical
- 6. Fasting blood sugar > 120 mg/dl (fbs): categorical 0 = False and 1 = True
- 7. Resting electrocardiographic results (restecg): categorical values 0, 1 and 2
- 8. Maximum heart rate achieved (thalach): numerical
- 9. Exercise induced angina (exang): categorical 0 = no and 1 = yes
- 10. ST depression induced by exercise relative to rest (Oldpeak): numerical
- 11. The slope of the peak exercise ST segment (slope): categorical values 0, 1 and 2
- 12. Number of major vessels colored by flourosopy (ca): categorical values 0, 1, 2, 3 and 4
- 13. **Thal:** categorical 1 = normal, 2 = fixed defect and 3 = reversible defect
- 14. **Target:** categorical 0 = no presence of heart disease and 1 = presence of heart disease

The table 1 shown below is just a glimpse on the first 10 entries of the dataset.

	age	sex	ср	trestbps	chol	fbs	restecg	thalach	exang	oldpeak	slope	ca	thal	target
0	63	1	3	145	233	1	0	150	0	2.3	0	0	1	1
1	37	1	2	130	250	0	1	187	0	3.5	0	0	2	1
2	41	0	1	130	204	0	0	172	0	1.4	2	0	2	1
3	56	1	1	120	236	0	1	178	0	0.8	2	0	2	1
4	57	0	0	120	354	0	1	163	1	0.6	2	0	2	1
5	57	1	0	140	192	0	1	148	0	0.4	1	0	1	1
6	56	0	1	140	294	0	0	153	0	1.3	1	0	2	1
7	44	1	1	120	263	0	1	173	0	0.0	2	0	3	1
8	52	1	2	172	199	1	1	162	0	0.5	2	0	3	1
9	57	1	2	150	168	0	1	174	0	1.6	2	0	2	1

Table 1 – A glimpse in the dataset

There are no missing values in the entire dataset. There are some outliers according to the 1.5*IQR rule and this will be addressed in the methodology section.

The image below shows the distribution of the numerical features and the scatter plot.

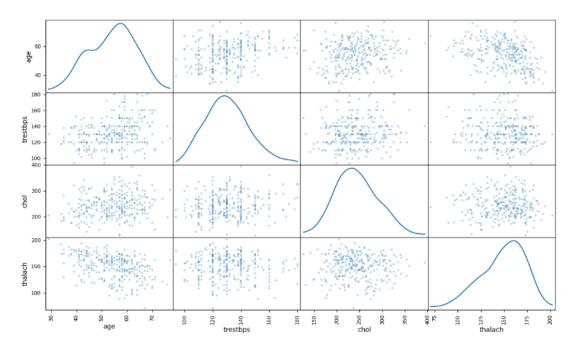


Figure 1 – Distribution and scatter plot of numerical features

It is possible to see that the distributions are skewed either to the left or to the right. Also, although not very clear, it is possible to see a correlation – or at least a trend - between "age" and "chol" and between "age" and "thalach" looking at the scatter plot. For the nine others categorical data, the distribution is also provided.

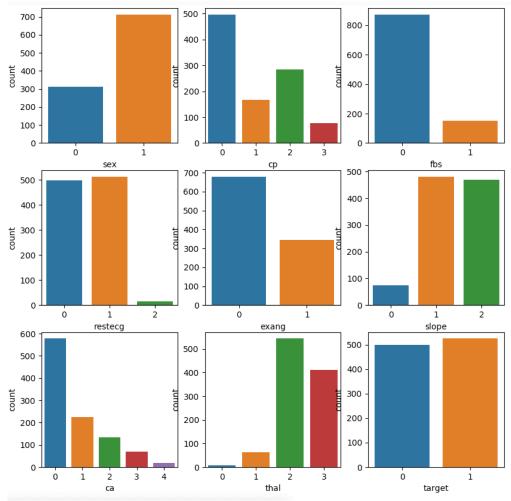


Figure 2 – Distribution of categorical features

As stated before, the target variable is well balanced. Even though the target variable will be not a problem, it would be nice to switch from accuracy to confusion matrix to see the performance of the model in others aspects. Since the goal is to determine if someone is prone to have a heart disease, it is more important for our model to be more accurate about "having the disease" than to "be health".

This dataset is being used for the last 30 years in studies and although the features are imbalanced, it may represent well the population in a statistic point of view. That said, the model may not present problems like the Amazon's known case, where their model to chose resumes was flaw because the features about people was imbalanced – and not statistically representative, preventing the model to suggest a "black man" or a "white woman" as a good candidate and only suggesting "white man" as good candidate.

Algorithms and techniques

The dataset has a target variable, so the solution is to use different supervised machine learning techniques and compare them to see which one would be more

accurate to solve the problem. After determining the fine-tuned supervised method, the trained model will be run against the test dataset for validation.

Since it is a categorical prediction problem, the supervised algorithms chose are:

- 1. Gradient Boosting Classifier
- 2. Random Forest Classifier
- 3. Logistic Regression
- 4. SVC

All these algorithms are contained in the Sklearn Python library as well as the metrics validation as accuracy, f1 score and confusion matrix.

In order to fine-tune the algorithms parameters, we will use a technique called Grid search, where some parameter values are passed in and it iterates these values using the algorithm. In other words, it would be like running the same algorithm over and over again, but in each time changing just one parameter to see how it changes the performance. This is a good way to find the optimal parameters for an algorithm given a specific model.

Benchmark model

The article "Heart Disease Prediction Using Machine learning and Data Mining Technique" reach a best result of 56.76% accuracy using Decision Tree with algorithm J48 with Reduced Error Pruning. This model is going to be used as the Benchmark model, since it is a categorical prediction model using the same dataset.

III Methodology

Data preprocessing

Processing the data does not mean that the end result of the machine learning will be better. That said, although some processing techniques will be applied, the machine learning processes will also handle the original dataset for comparison purposes.

The first step is to look for missing values. Fortunately, there is none in this dataset nor there is any mismatching type or similar problem.

The second step is to look for outliers. Applying the standard 1.5*IQR rule, we find out that 4.95% of the value fit this rule. This number is not to ignore. Let us tighten this rule making the threshold 2.0*IQR. Now the real outliers represent only 1.98% of the whole dataset, reducing the total number from 303 to 297 samples.

In figure 1 of the *II Analysis* section we saw that the distributions were either right or left skewed. Applying the natural log transformation, we get a slightly better curve as shown in the figure 3 below.

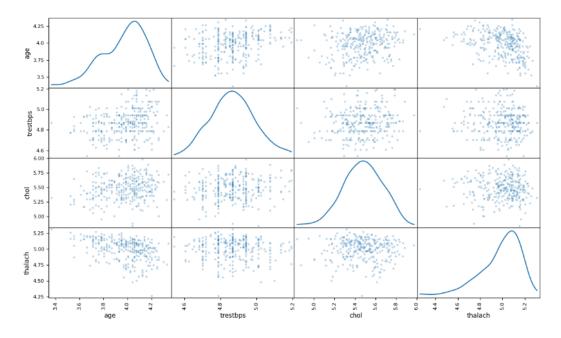


Figure 3 – Distribution and scatter plot of numerical features after log transformation

Implementation

The first step is to split the dataset in two parts, one for training the model and other to validate it. For that, the *train_test_split* function of Sklearn Python library will be used. In order to others be able to reproduce the results here presented, this *train_test_split* function will be used with a *random_state = 42* parameter, guaranteeing that everyone running the same algorithm with the same parameters will obtain the same results. The rate used here is 70% of the sample for model training purpose and 30% of the sample for testing purpose.

These training and test datasets will be running against 4 supervised machine learning algorithms using the Search Grid function. This methodology will be applied within 3 different datasets in that order:

- Original dataset
- Dateset with outliers removed
- Dataset with outliers removed and log transformation applied

The table below summarizes all the parameters and parameters values of grid search used in this work. The same algorithms, the same parameters and same parameters values are going to be used against all the 3 datasets listed above.

Algorithm	Parameter	Value		
		1		
	min complex loof	5		
Cundinut	min_samples_leaf	10		
Gradient Boosting		20		
Classifier		20		
Classifier	n_estimators	50		
	ii_estiiiatois	100		
		150		
		1		
	min_samples_leaf	5		
	iiiii_saiiipies_ieai	10		
Random Forest		20		
Classifier		20		
	n_estimators	50		
	ii_estiiiatois	100		
		150		
	penalty	L1		
	penalty	L2		
Logistic		0.1		
Regression	С	0.5		
	C	1.0		
		5.0		
		Linear		
	kernel	Rbf		
SVC		Sigmoid		
	С	0.5		
	C	1.0		

Table 2 – Algorithms, parameters and parameters values to be used in the analysis

IV Results

Model evaluation e validation

The table 3, 4 and 5 below summarizes the best result of each algorithm, which parameters led to that result and which dataset it came from.

Dataset: Original									
		Final result							
Algorithm	Parameters	Target Value	Precision	Recall	F1 score	Accuracy			
Gradient Boosting	min_samples_leaf = 1	0	0.84	0.78	0.81	0.84			
Classifier	n_estimators = 100	1	0.83	0.88	0.85				
Random Forest	min_samples_leaf = 1	0	0.84	0.78	0.81	0.84			
Classifier	n_estimators = 10	1	0.83	0.88	0.85				
Logistic Pagrossian	Penalty = L2	0	0.80	0.78	0.79	0.81			
Logistic Regression	C = 1.0	1	0.82	0.84	0.83				
SVC	Kernel = Rbf	0	0.80	0.80	0.80	0.82			
370	C = 1.0	1	0.84	0.84	0.84				

Table 3 – Best results per algorithm for Original dataset

Dataset: Outliers removed									
		Final result							
Algorithm	Parameters	Target Value	Precision	Recall	F1 score	Accuracy			
Gradient Boosting	min_samples_leaf = 1	0	0.76	0.74	0.75	0.79			
Classifier	n_estimators = 100	1	0.81	0.83	0.82				
Random Forest	min_samples_leaf = 1	0	0.79	0.71	0.75	0.80			
Classifier	n_estimators = 10	1	0.80	0.87	0.83				
Logistic Rogression	Penalty = L2	0	0.87	0.71	0.78	0.83			
Logistic Regression	C = 1.0	1	0.81	0.92	0.86				
SVC	Kernel = Rbf	0	0.84	0.71	0.77	0.82			
370	C = 1.0	1	0.81	0.90	0.85				

Table 4 – Best results per algorithm for dataset with outliers removed

Dataset: Outliers removed and log transformation applied									
		Final result							
Algorithm	Parameters	Target Value	Precision	Recall	F1 score	Accuracy			
Gradient Boosting	min_samples_leaf = 1	0	0.78	0.74	0.76	0.80			
Classifier	n_estimators = 100	1	0.81	0.85	0.83	0.80			
Random Forest	min_samples_leaf = 1	0	0.73	0.84	0.78	0.80			
Classifier	n_estimators = 10	1	0.87	0.77	0.82	0.80			
Logistic Bograssian	Penalty = L2	0	0.81	0.68	0.74	0.80			
Logistic Regression	C = 1.0	1	0.79	0.88	0.84				
SVC	Kernel = Rbf	0	0.88	0.74	0.80	0.04			
370	C = 1.0	1	0.83	0.92	0.87	0.84			

Table 5 – Best results per algorithm for dataset with outliers removed and log transformation applied

For the original dataset, the best result was obtained using the either the Gradient Boosting Classifier or the Random Forest Classifier, both resulting in a F1 score of 0.81 for predicting health people and 0.85 for predicting people with heart disease and an overall accuracy of 0.84.

Considering the dataset with the outliers removed, the best algorithm is the Logistic Regression resulting in a F1 score of 0.78 for predicting health people and 0.86 for predicting people with heart disease and an overall accuracy of 0.83.

Considering the dataset with the outliers removed and the log transformation applied, the best algorithm is the SVC resulting in a F1 score of 0.80 for predicting health people and 0.87 for predicting people with heart disease and an overall accuracy of 0.84.

That said, the best alternative is considering the outlier removal, log transformation and applying the SVC algorithm which give us the 87% accuracy in predicting that someone has heart disease.

It is also possible to see that all the results in all three conditions lad to good results. This shows that changes in the dataset does not seem to greatly change the result.

Justification

Comparing to the benchmarking model, our model outstands. Even our Random Forest algorithm did way better than the benchmarking model, which is a Decision Tree with algorithm J48 with Reduced Error Pruning resulting in 56.76% accuracy.

The authors of the benchmarking paper didn't preprocessing the data, but the interesting point is that even our untreated data, i.e. the original dataset, performed way better with a accuracy of 81% to 84%.

V Conclusion

Although the complete dataset has 76 features, in the last 30 years only 14 features, including the target variable has proven enough for the academic studies. The benchmark paper got an accuracy that is not good enough to be used as a tool in the real world. With the result achieved in this work of 87% accuracy to predict a true positive, one can say that this model can be used in the real world as a predictor tool or, at least, be a trial guide for further investigation if one really got a heart disease.

Reflection

One interesting point in this work is that although Neural Networks and deep learning are the top-notch trends nowadays, simple supervised machine learning models are still very powerful to solve a wide range of important problems as this one.

The second point to be highlighted here is the fact that the features are very imbalanced and still it was not a bit deal to our model and results. That is because as stated before, the imbalanced features really represents imbalanced characteristics of the population. That said, our model was trained considering just some people with severe chest pain and in the real life just a few people really have a severe chest pain. Again, that is a different scenario from the Amazon case previously mentioned in this work, where "white man" was highly present in the training dataset as a good candidate, while in the real population, half of the people are woman and a considerably amount of this same population is black.

Improvement

Further studies can be done with unsupervised learning clusterization algorithms to try to determine groups with similar characteristics and split the original dataset in smaller datasets and then training supervised models against them. This can be good to improve the accuracy of the F1 score and can define a real life tool where the subject is first filtered accordingly to the cluster group and then analyzed by the supervised model to get a more trustworthy result.

It would be also interesting to use this dataset in neural networks algorithms to compare performance.