Projectreport Machine Learning 2

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1. Introduction

Our general idea was to work with LGBT-related data. This was not as easy as expected, since it seems there are not a lot of datasets openly available that have that kind of information. Finally, we found a US survey by the CDC, that regularly monitors the country's youth in a lot of dimensions, but among other questions also asks for sexual experiences and identification.

2. The Dataset

"The *Youth Risk Behavior Survey (YRBS)* measures health-related behaviors and experiences that can lead to death and disability among youth and adults.[...] Some of the health-related behaviors and experiences monitored are: - Student demographics: sex, sexual identity, race and ethnicity, and grade - Youth health behaviors and conditions: sexual, injury and violence, bullying, diet and physical activity, obesity, and mental health, including suicide - Substance use behaviors: electronic vapor product and tobacco product use, alcohol use, and other drug use - Student experiences: parental monitoring, school connectedness, unstable housing, and exposure to community violence [1]. It is a national survey conducted by the CDC (Center for Disease Control and Prevention) and includes high school students from both private and public schools within the U.S. Data is collected from 1991 through 2021, we are only using the most recent data from 2021. If you want to learn more about the data there is an accompanying Data User Guide.[2].

2.1 Preprocessing of the dataset

To preprocess the dataset, we first ran a summary of our dataset. The number of NAs seems to depend very much on the question. The variable "orig_rec" only contained NAs and has therefore been removed, as well as the variable "site" which only contained "XX" entries. Variables q4 and q5 are already aggregated in "raceeth" and have also been deleted. The variable "record" seems to be an ID for the observations. This has to be considered later.

2.2 Missing Data

We will first exclude all the observations with NAs in all the target-related variables q25 to q29. Since we want to build our target variable on these questions, the target variable cannot all be empty. The amount of data available should be enough to just exclude these observations. After removing the observations that have NAs in all the variables, that are used to create our target variable, we still have around 13.7% NAs in the dataset.
What if we had just excluded every NA in the dataset? We will try and see if this is a viable option, since this woul not just be quick and easy, but we would also just have "real" answers. The exclusion of NAs leads to a severe reduction in the number of observations. The original data consisted of 17232 observations, after reducing the target-related NAs only, we have 11753 observations left. If we omit all NAs, the reduced dataframe still has 4334 observations.
In this case need to assess the loss of information foremost about our target variable. The important question is if there is a pattern to the missingness in our data, not just, but especially about our target variable.

2.2.2 Omitting NAs vs Data Imputation

If we can omit the NAs or if it may be necessary to impute the missing data points, depends on the type of missingness. If data is missing completely at random (MCAR), we can omit the NAs, if it is just missing at random (MAR) we would rather impute the data. If the data is missing not at random (MNAR), it would be a quite difficult problem because we cannot easily impute the missing data then. To find out if we can just omit the data, an MCAR test was applied.

We test the target-related variables q25 to q29 for potential pattern(s) in the missing data. This results in a p-value of 0, which means we can say for sure, that the data is not missing completely at random. Just omitting all NAs could be problematic and lead to bias.

```
## # A tibble: 1 x 4
##   statistic    df p.value missing.patterns
##       <dbl> <dbl>   <dbl>            <int>
## 1      452.    68       0               29
```

Because of this, we will use a rule base approach to create the target variable and impute the predictive variables afterwards. To ensure a good imputation, we need to impute our NAs before reducing the dataset to 2000 observations. To run the imputation properly we need to factorize our nominal and ordinal variables first.

2.3 Target Variable

As a target variable, we decided to calculate a score from 5 questions that reflect the suicide risk of the person (observation) in question. This score is aggregated with a rule-based approach.

After creating the target variable we need to exclude the variables q25 to q29, which were used for creating it, from our dataset.

2.4 Imputation

2.5 Reducing and balancing the dataset to 2000 observations

We need to reduce our data to the maximum allowed size of 2000 observations. To ensure the best possible data quality, we want to ensure that our dataset is balanced. Intuitively, we are considering if it is best to still use as much of the non-imputed data for our smaller dataset as possible, before filling it up with imputed data, since non-imputed data is usually of better quality. On the other hand the data seemingly shows patterns in the missingness so there are reasons to just do a stratified sampling over the imputed data as well. To do a proper statified sampling we need to identify the stratification variables. Therefore we will calculate the correlations with the target variable and see which variables are highly correlated to our target variable. These will then as well as the target variable be used as stratification variables.

2.6 Simple Synopsis of the Dataset

number of observations: 2000 number of variables: 100

datatypes: factor: 96 - nominal variables: 29 - ordinal variables: 67 numeric variables: 4 - discrete variables: 96 (here all factor variables) - continuous variables: 4 (here all numeric variables)

3. Additional Data Preparation

question: would it introduce information leakage to reduce the features before splitting the data?

3.1 Feature Reduction

Since our dataset has lots of variables, we decided to start by excluding some variables depending on the estimated feature importance.

3.1.1 Correlations

Unfortunately at this point we have to many variables to do a pairs plot or correlation plot with a visually usable outcome. We will therefore perform a correlation analysis only with respect to the target variable and in numeric format instead of any visual plot.

The 18 variables with high correlations (>0.25) with our target variable are:
q93
q85
q45
q35
q36
q39
q64
q46
q20
q30
q24
q19
q98
q34 q41 q43 q44 q47

3.1.2 Feature Selection Algorithm

Since the data still has a lot of variables, we need to use a feature selection technique to reduce the features before using a machine learning method. We chose to use model agnostic methods, because the feature selection should be valid for all methods that are later compared. In an earlier step the variables most correlated with the target variable were already identified. Unfortunately this captures only linear monotonous relationships in the data and does not work well for our nominal categorical features.

[maybe delete chi^2 test + text] We will also use a Chi^2 test between our variables and our target variable to assess their relationship with regards to independence. The variables that are found to have a significant relationship (p >= 0.05 %) with the target are kept.

To capture non-linear relationships as well, information gain between the target and the predictor variables is measured as well. The variables with high information gain with respect to the target variable are kept, because they can contribute more in predicting the target variable.

3.1.3 Feature Selection with Chi^2

3.1.4 Information Gain for Feature Selection

3.1.5 Domain Knowledge for Final Feature Selection

```
## [1] "q7orig" "psu"    "q40"    "q6orig" "q22"
```

Let's see what those variables actually stand for. "q7orig" and "q6orig" cannot be found in the data manual and will therefore be discarded. According to the data manual, "PSUs consist of counties, groups of smaller adjacent counties, or sub-areas of very large counties. "PSU" indicates the PSU the school the student attends was assigned to." (p.14). It is possible, that the district/locality of a school plays a role in the risk of

suicide. For example for queer students in a very religious place. Q22 is the variable that describes physical dating violence. Therefore q22 is also a valid choice as a predictor variable for our suicidal score target variable. Q40 encodes the range of age when a student first got into contact with drinking alcohol. This might be an indicator for a negligent social surrounding if someone is exposed to an alcoholic drink in an early age and therefore also could be a valid predictor variable in our case.

3.2 Naming the Variables

For an easy understanding of their values, the variables and levels are named according to their content. This is an optional step. It can lead to a better readability of our data frame though.

3.3 Splitting the Data

According to the project requirements we split our data in 60% Training, 20% Validation and 20% Testing Data. Since we want to do a cross validation we will split into 80% Training and 20% Testing Data.

4. Machine Learning Models

In the following, we compare two Supervised Learning methods on our reduced YOUTH AT RISK dataset to predict our suicidal_class target variable. Because we want to be able to also capture non-linear relationships in our data, we chose Support Vector Machines and Naive Bayes.

4.1 Short Mathematical Overview on the used Methods

4.1.1 Naive Bayes Classification

Naive Bayes Classifiers differs from a theoretically ideal Bayes Classification by assuming the independence of predictor variables.

4.1.2 Support Vector Machines (SVM)

The name Support Vector Machines already describes some elements of this method. A certain number of data points will define the (linear) boundary between two classification regions, these are called the support vectors. The support vectors are the datapoints (observations) that lie closest to our decision boundary. The boundary in two dimensional space is a line, in three dimensions a plane and in more than three dimensional space a hyperplane. For our dataset, we need a multidimensional hyperplane. The number of dimensions depend on the number of our predictor variables. We need to find the hyperplane, that separate our data into the classes of our target variable best. The best hyperplane is the one that maximizes the margin between the support vectors of the different classes. The margin is a strip on each of the boundaries sides. In the case of a hard classifier this strip does not contain any points. But we will have a soft classifier with the cost C as a hyperparameter. This cost C describes a budget that we allow for points within the margin or on the other side of the boundary. Depending on the position of the point, the amount it attributes to the total cost changes. A point on the correct side of the margin will not attribute to the total cost at all. If it is in the margin, but on the correct side of the boundary it will attribute between zero and one, if on the wrong side of the boundary but within the margin it will attribute with one to two and if on the wrong side of the margin it will cost more than two. Mathematically we need to maximize the Margin M with respect to α_0, α_i and ϵ_i in the following objective function:

$$y_i \left(\alpha_0 + \sum_{j=1}^{n} \alpha_i K(\mathbf{x}_i, \mathbf{x}_j) \right) \geq M(1 - \epsilon_i)$$

4

The following constraints are given: $\sum_{i=1}^{n} \alpha_i^2 = 1$, $\epsilon_i \geq 0$ and $\sum_{i=1}^{n} \epsilon_i < C$ with $C \geq 0$ and $i = 1, ..., n$
For our SVM approach, we will try different kernels K as hyperparameters:

A linear Kernel

$$K(u,v) = \langle \mathbf{u}, \mathbf{v} \rangle = \sum_{j=1}^{p} u_j v_j$$

a polynomial Kernel

$$K(u,v) = (c + \langle \mathbf{u}, \mathbf{v} \rangle)^d, \ with \ d > 1$$

and a radial Kernel

$$K(u,v) = \exp\left(-\gamma \sum_{i=1}^{p} (u_i - v_i)\right)$$

4.2 Preprocessing for SVM

svm in R tutorial + nice visualization: https://www.datacamp.com/tutorial/support-vector-machines-r

4.3 SVM Fitting and Hyperparameter Optimization

for SVM: try different available kernels and other hyperparameters and cross validate

Given the list of all resulting models, coming from the HPO of the SVM, we extract the best model for further inspection and results.

Mithilfe eines Likelihood-Ratio Tests wird überprüft welches Modell das bessere ist.

Das optimierte Modell ist signifikant besser als das erarbeitete zum festgelegten Signifikanzniveau, die Nullhypothese des LR-Tests, dass das optimierte Modell nicht besser ist als das komplexere, kann verworfen werden. Wir stellen außerdem fest: Beide Modelle weisen einen signifikanten Erklärungsgehalt auf, sie sind besser als das Nullmodell.

5. Comparison of the Models / Model's Performance on Test Data

if one model performs better, is this improvement significant to a usual significance level?

5.1 Quantitative

5.1.1 Confusion Matrix

5.1.2 Accuracy

5.1.2 Precision, Recall and F1-Score

5.2 Qualitative

5.3 Overfitting Check
compare training/validation and testing curves. . .

6. Visual Representation

Dimensionality Reduction Techniques: Sometimes, it's helpful to use dimensionality reduction techniques (like PCA) to identify the most significant variables or components and then focus the GAM analysis and visualization on these.

7. Final Discussion

8. References

[1] https://www.cdc.gov/healthyyouth/data/yrbs/overview.htm

[2] https: //www.cdc.gov/healthyyouth/data/yrbs/pdf/2021/2021_YRBS_Data_Users_Guide_508.pdf