HW2 Report

Part 1

Q1:

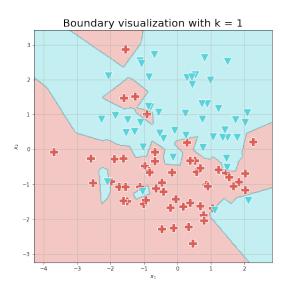


Figure 1: Boundary Visualization for KNN with $\mathbf{K}=1$

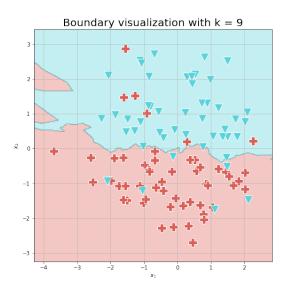


Figure 2: Boundary Visualization for KNN with K=9

As can be seen in figure 1, the model with k equal to 1 caused overfitting to occur, as for every new sample point, its prediction took into account only the closest point to it, thereby increasing the complexity of the model since every new point is predicted locally. On the other hand, as can be seen in figure 2, the model with k equal to nine takes into account the nine closest points to the sample being predicted, and therefore takes a broader outlook on the training data in-order to tdo the prediction, thereby lowering the complexity. But at the same time, we can see that a number of points were not classified correctly, and therefore this could indicate that this model is performing underfitting.

Q2:

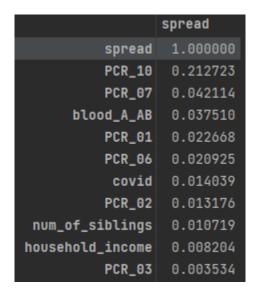


Figure 3: The 10 most correlated features to spread

Q3:

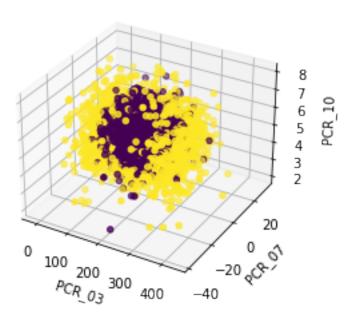


Figure 4: 3D Scatter plot of PCR_03, PCR_07, and PCR_10 according to spread

Q4:

The points with label -1 for spread are contained within an ellipsoid described by the following equation:

$$\frac{PCR_03}{a^2} + \frac{PCR_07}{b^2} + \frac{PCR_10}{c^2} = 1$$

where a, b, c are positive real numbers.

Q5:

The training accuracy for k = 11 is 81.333%.

Q6:

Z-score scaling scales the data of a feature by ensuring that they have zero mean and unit standard deviation, thereby causing the data to adhere to a normal distribution. Features scaled according to this technique have their outliers handled correctly, but no guarantee on the resulting range of the data is made, and the ranges of different features scaled according to Z-score may differ from each other. This technique is preferable in cases that have outliers and when the learning model assumes that the data adheres to a normal distribution. On-the-other-hand, the min-max technique involves scaling the data of a feature to a specific range (generally between 0 and 1). Contrary to the Z-score method, this technique guarantees a uniform range across features and maintains the original distribution of the data, but does not handle outliers well. Therefore, it would be preferable to use this technique only when the feature in question has no significant outliers and or the learning model to be used requires the feature data to fall within a certain range.

Q7:

The updated accuracy after normalizing the data is 88.5416%. This is an increase of approximately 7% over the accuracy from the non-normalized data. This can be explained by the fact that KNN is an algorithm that uses euclidean distances between data points for its predictions. Therefore, it is very sensitive to differences in scales between different features, and thus features that have undergone normalization, which equalizes or nearly equalizes the scales and ranges of different features, will do better in KNN because no single feature will dominate and have more influence over other features by way of having a larger scale.

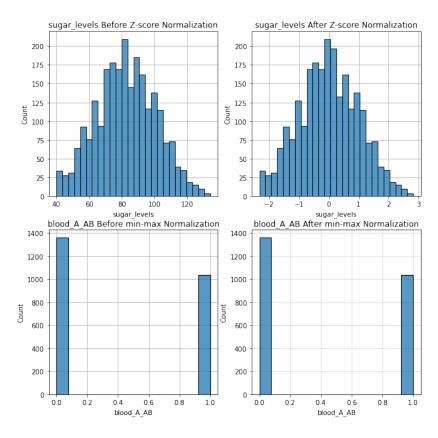


Figure 5: Histogram Plots of sugar_levels and blood_A_AB Before and After Normalization

sugar_levels We chose to do Z-score Normalization on sugar_levels since it already obeys an approximate normal distribution, as can be seen in figure 5, and therefore we would not be forcibly changing its distribution through the normalization process. Furthermore, we will not be using it in the KNN model, which works better with min-max normalized features, but rather with SGD SVM, which works better with Z-score normalized features, and a decision tree, which is mostly agnostic to Normalization.

blood_A_AB In contrast, we chose to do min-max normalization on blood_A_AB, since it does not obey a normal distribution, but rather already obeyed a binary distribution of 0 and 1, which we did not want to modify. Furthermore, it does not have outliers which could negatively affect the min-max normalization.

Q9:

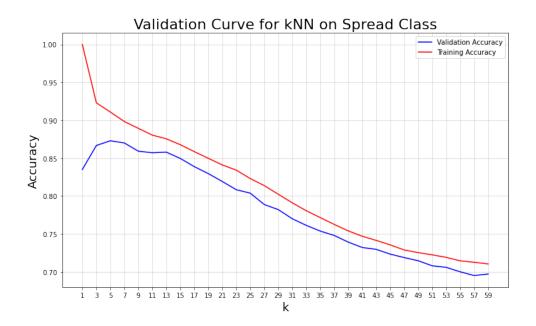


Figure 6: Validation Curve for kNN on spread Class

The optimal k value is 5, as can be seen by virtue of the fact that in figure 6, the highest point (peak) of the validation curve is at 5. The mean validation and training accuracies for k = 5 are 87.2916% and 91.083%, respectively.

Part 2

Q11:

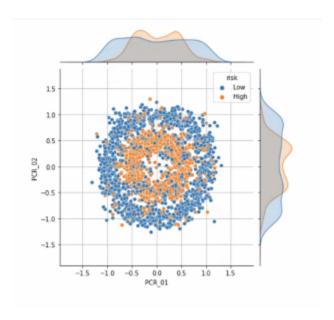


Figure 7: Joint plot of PCR_01 and PCR_02 according to risk

As can be seen from the scatter portion of the jointplot in figure 7, the plot is mostly separable into radiuses, and therefore it seems likely that PCR_01 and PCR_02 will be important in predicting the risk class.

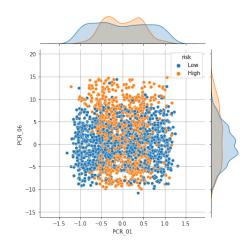


Figure 8: Joint plot of PCR_06 versus PCR_01 with respect to risk

Furthermore, from the scatter portion of the jointplot in figure 8 we noticed a mostly separable form similar to the letter "H" where the "H" itself is made up of a high proportion of points with low risk surrounded by clusters of points of high risk. Therefore, we can conclude that in addition, PCR_06 will be important in predicting the risk class.

Q12:



Figure 9: The 10 most correlated features to risk

Part 3

Q20:

In Q19, we used a 2 dimensional polynomial transform on the features within the primal objective of the SVM model. This approach is slower than the two dimensional kernel objective because it involves computing inner products between data points and the w vector in high dimensions, whereas in the kernel case, the dual objective is used, which only involves computing the kernel function for 2 dimensional polynomials which is of much lower dimension (and finding and storing the α s, which is not too computationally burdensome on the assumption that the support vectors are sparse).