HW 03: Classification

due Thursday, Oct. 6 at 11:59pm

Exercise 1

Suppose I have data about patients who have Melanoma disease. My response Y is binary status, where Y = 0 implies a patient died, and Y = 1 means the patient is alive. I have the following predictors: a binary predictor ulcer (1 = has an ulcer, 0 = no ulcer), and continuous predictor thickness of the tumour (mm). I fit the model in R and obtain the following results.

```
##
## Call:
  glm(formula = status ~ ulcer + thickness, family = "binomial",
##
       data = melanoma)
##
##
  Deviance Residuals:
##
       Min
                 10
                      Median
                                    30
                                            Max
  -2.0725
                                         1.7332
##
           -1.0787
                      0.5229
                               0.6453
##
## Coefficients:
##
               Estimate Std. Error z value Pr(>|z|)
                2.07388
                           0.31013
                                      6.687 2.28e-11 ***
##
  (Intercept)
## ulcer1
               -1.39220
                           0.37229
                                    -3.740 0.000184 ***
## thickness
                                    -2.500 0.012425 *
               -0.15782
                           0.06313
##
## Signif. codes: 0 '***' 0.001 '**' 0.05 '.' 0.1 ' ' 1
##
##
   (Dispersion parameter for binomial family taken to be 1)
##
##
       Null deviance: 232.84
                              on 190
                                      degrees of freedom
## Residual deviance: 198.98
                              on 188
                                      degrees of freedom
## AIC: 204.98
##
## Number of Fisher Scoring iterations: 4
```

- a) Write down the fitted logistic regression model based on this summary output from R.
- b) Interpret the coefficients $\hat{\beta}_0$, $\hat{\beta}_1$, and $\hat{\beta}_2$.
- c) Given that someone has an ulcer and the melanoma tumour has a thickness of 10mm, what is their probability of dying from melanoma?
- d) The following displays the contingency table for predictions on the training data using the model. Obtain the following three rates: overall error rate, false positive rate, and false negative rate.

```
## truth
## pred 0 1
## 0 19 12
## 1 38 122
```

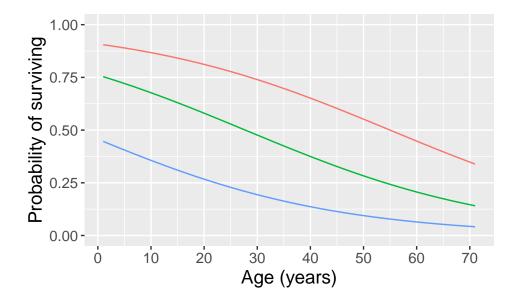
Exercise 2

The sinking of the Titanic is one of the most infamous shipwrecks in history. We have data about the survival status of the passengers (0 = died, 1 = survived), as well as some other demographic and socioeconomic information. I fit a logistic regression model to the survival status of the passengers, regressed on age (years) and ticket class (categorical with three levels: 1st, 2nd, 3rd).

After fitting the model, I have the following estimates coefficients:

```
## (Intercept) Age class1 class3
## 1.1584789 -0.0417545 1.1375334 -1.3320273
```

The following plot displays the estimated probability of survival across age, at the three different ticket classes.



- a) Based on this information, match each class to one of the lines in the plot. Briefly explain your reasoning.
- b) Based on your decision in (a), approximately how old would someone in third class have to be in order to have the same probability of survival as a 70-year old person in first class?

Exercise 3

Recall that the Naive Bayes classifier assumes that the p features are independent, i.e. $f_k(x) = f_{k1}(x_1) \times f_{k2}(x_2) \times \cdots \times f_{kp}(x_p)$. For categorical predictors, an estimate for each of the f_{kj} is simply the proportion of training observations for the j-th predictor corresponding to each class k.

Assume in this problem that we have n=200 training observations with p=2 features $(X_1$ and $X_2)$, and the response has K=3 possible classes. Further, assume that X_1 and X_2 are both qualitative, where X_1 has three levels and X_2 has two levels.

The following displays the "aggregated" or summary data. That is, you are provided with the total counts within each class and predictor combination, as opposed to the individual, raw observations. For example, 22 observations have $X_1 = \text{brown}$ and Y = class 1. Similarly, 8 observation have $X_2 = \text{tree}$ and Y = class 2.

```
##
## X1
             1
                 2
                    .3
##
     brown 22 16 33
##
            11
                6 23
     red
     white 29 23 37
##
                  2
## X2
               1
     ground 48 37 71
##
##
     tree
             14
                  8 22
```

- a) What are the number of observations within each class of Y? Call them n_1, n_2, n_3 .
- b) Using your answer from (a), what are $f_{kj}(x_i)$ for k = 1, 2, 3 and j = 1, 2?
- c) Assume that $\hat{\pi}_1 = 0.3$, $\hat{\pi}_2 = 0.3$, and $\hat{\pi}_3 = 0.4$. Suppose that we wish to classify a new observation $x^* = (\text{red, ground})'$. Under the Naive Bayes classifier, calculate $\Pr(Y = k | X = x^*)$ for each k. Based on your answer, what would you classify this new observation as?
- d) Now let's say that our new observation is $x^* = (\text{green}, \text{ground})'$. For this new observation, what would $\Pr(Y = k | X = x^*)$ be for each k = 1, 2, 3, and why? This is known as the *zero-frequency* problem.

Exercise 4

When the number of features p is large, there tends to be a deterioration in the performance of KNN and other local approaches that perform prediction using only observations that are near the test observation for which a prediction must be made. Organizing/clustering/searching data often relies on detecting spaces where objects form groups with similar properties. When p is large, we will see that the objects (features) will seem very dissimilar. This phenomenon is known as the curse of dimensionality, and it ties into the fact that non-parametric approaches often perform poorly when p is large. We will now investigate this curse (which is common in many disciplines, e.g. genetics).

- a) Suppose that we have a set of observations, each with measurements on p=1 feature, X. We assume that X is uniformly (evenly) distributed on [0,1]. Associated with each observation is a response value Y. Suppose that we wish to predict a test observation's response using only observations that are within 10% of the range of X closest to that test observation. For instance, in order to predict the response for a test observation with X=0.6, we will use observations with associated X in the range [0.55, 0.65]. On average, what fraction of the available observations will we use to make the prediction?
- b) Now suppose that we have a set of observations, each with measurements on p=2 features, X_1 and X_2 . We assume that (X_1,X_2) are uniformly distributed on $[0,1] \times [0,1]$. We wish to predict a test observation's response using only observations that are within 10% of the range of X_1 and within 10% of the range of X_2 closest to that test observation. For instance, in order to predict the response for a test observation with $X_1=0.6$ and $X_2=0.35$, we will use observations in the range [0.55,0.65] for X_1 and in the range [0.3,0.4] for X_2 . On average, what fraction of the available observations will we use to make the prediction?
- c) Now suppose that we have a set of observations on p=100 features. Again the observations are uniformly distributed on each feature, and again each feature ranges in value from 0 to 1. We wish to predict a test observation's response using observations within the 10% of each feature's range that is closest to that test observation. What fraction of the available observations will we use to make the prediction?
- d) Using your answers to parts (a)–(c), argue that a drawback of KNN when p is large is that there are very few training observations "near" any given test observation.

e) Now suppose that we wish to make a prediction for a test observation by creating a p-dimensional hypercube centered around the test observation that contains, on average, 10% of the training observations. For p = 1,2, and 100, what is the length of each side of the hypercube? Comment on your answer.

Note: A hypercube is a generalization of a cube to an arbitrary number of dimensions. When p = 1, a hypercube is simply a line segment, when p = 2 it is a square, and when p = 100 it is a 100-dimensional cube.

BONUS (each part is worth extra points)

Suppose that you wish to classify an observation into a toxic or nontoxic class, based on a predictor X. You fit a logistic regression model and find that

$$\widehat{\Pr}(Y = \text{toxic}|X = x) = \frac{e^{\widehat{\beta}_0 + \widehat{\beta}_1 x}}{1 + e^{\widehat{\beta}_0 + \widehat{\beta}_1 x}}$$

I also fit a logistic regression model to the same data, but I use the softmax formulation and find that

$$\widehat{\Pr}(Y = \text{toxic}|X = x) = \frac{e^{\hat{\alpha}_{0,\text{toxic}} + \hat{\alpha}_{1,\text{toxic}}x}}{e^{\hat{\alpha}_{0,\text{toxic}} + \hat{\alpha}_{1,\text{toxic}}x} + e^{\hat{\alpha}_{0,\text{nontoxic}} + \hat{\alpha}_{1,\text{nontoxic}}x}}$$

(I am writing my model with α 's to not confuse with your model's β 's.)

- a) In my model, what are the log-odds of toxic vs nontoxic?
- b) Suppose that in your model, $\hat{\beta}_0 = -2$ and $\hat{\beta}_1 = 3$. What are the coefficient estimates in my model? Be as specific as possible.
- c) Now suppose that we fit the same two models on a different data set. This time, I estimate the coefficients $\hat{\alpha}_{0,\text{toxic}} = -1.5$, $\hat{\alpha}_{1,\text{toxic}} = 2$, $\hat{\alpha}_{0,\text{nontoxic}} = 1$, and $\hat{\alpha}_{1,\text{nontoxic}} = -1.25$. What are the coefficient estimates in your model?
- d) Finally, suppose you apply both models from (c) to a data set with 2000 test observations. What fraction of the time do you expect the predicted class labels from your model to agree with those from my model? Explain your answer.
- e) In the binary response case, would you prefer the usual logistic regression with the logit coding, or the softmax coding? Why?

Submission

Upload your assignment as a PDF file to Canvas. Please show all work!