Kyle Benson CS 273A - Machine Learning: Fall 2013 Homework 2

Problem 1: Bayes Classifiers

(a) Probabilities for Naive Bayes Classifier:

p(y=1) = 0.4

p(Variable = 1 y = 1)	x_1	x_2	x_3	x_4	x_5	П
Probability	0.75	0.0	0.75	0.5	0.25	0
p(Variable = 1 y = -1)	x_1	x_2	x_3	x_4	x_5	\prod
Probability	0.5	0.833	0.667	0.833	0.33	.0756

(b) For $x = (00000), p(y = 1) \prod_{i} p(x_i = 0) = 0.4 * (0.25 * 1 * 0.25 * 0.5 * 0.75) = 0.009375$ and $p(y = -1) \prod_{i} p(x_i = 0) = 0.6 * (0.5 * 0.167 * 0.33 * 0.167 * 0.67) = 0.00185$. So the predicted class would be y = 1

For $x = (11010), p(y = 1) \prod_i p(X_i = x_i) = 0.4 * (0.75 * 1 * 0.25 * 0.5 * 0.75) = 0.028$ and $p(y = -1) \prod_i p(x_i = 0) = 0.6 * (0.5 * 0.833 * 0.333 * 0.833 * 0.667) = 0.046236114$. So the predicted class would be y = -1

(c)
$$p(y = 1|x = (11010)) = \frac{p(y=1)p(x=(11010)|y=1)}{p(x=(11010))} = \frac{0.4*(0.75*1*0.25*0.5*0.75)}{(0.75*1*0.25*0.5*0.75) + (0.5*0.833*0.333*0.833*0.667)} = 0.0087$$

- (d) Because then our probability table will have $O(F^2)$, rather than O(F), entries, where F is the number of features we are training on, in order to account for dependence among the feature variables.
- (e) We do not need to re-train the model. Because x_1 is independent of all other x_i , we can safely ignore x_1 entirely and make our predictions based on the classifier: $p(y) \prod_{i \neq 1} x_i$

Problem 2: Decision Trees

(a)
$$H(y) = -0.4 * log(0.4) - 0.6 * log(0.6) = 0.971$$

(b) Using the formula $H(y|x_i = 0) = p(y = 1|x_i = 0) \log p(y = 1|x_i = 0) + p(y = 0|x_i = 0) \log p(y = 0)$

	Variable	x_1	x_2	x_3	x_4	x_5
$0 x_i=0)$	Entropy $H(y x_i=0)$	1	0.4312	1.03	0.931	0.887
	Entropy $H(y x_i=1)$	0.811	0.22	0.701	0.72	1.028
	Info Gain	0.0465	0.61	0.006	0.0914	0.006

Therefore, we should split on feature x_2 first.

(c) After splitting on x_2 , we recompute the following info gains for each of the remaining features: $H(y|x_2=0)=0.722$

Variable	x_1	x_2	x_3	x_4	x_5
Info Gain	0.322	-	0.073	0.171	0.073

 $H(y|x_2=1)=0$, so no need to explore this branch

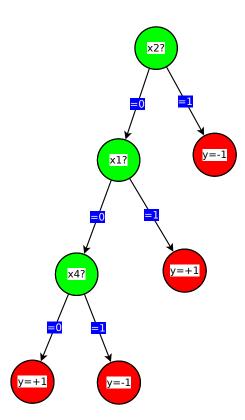
We therefore pick x_1 as the next feature to split. Again, we recompute the following info gains for each of the remaining features:

 $H(y|x_2=0, x_1=0)=1$

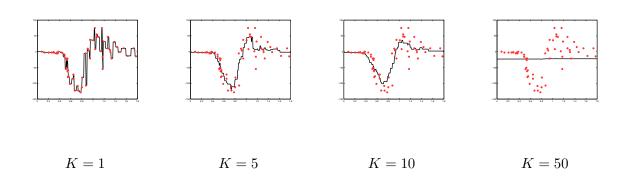
Variable	x_1	x_2	x_3	x_4	x_5
Info Gain	-	-	0	1	0

 $H(y|x_2=0,x_1=1)=0$, so no need to explore this branch

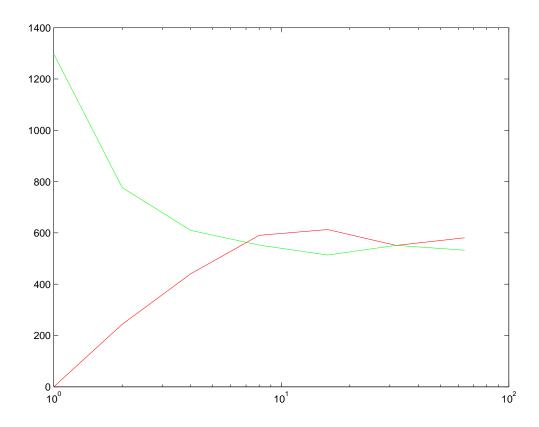
We therefore pick x_4 as the next feature to split and are done because the information gain is 1, meaning that feature x_4 perfectly explains this branch of the data $(x_1 = 0, x_2 = 0)$. This gives us the following decision tree:



Problem 2: K-Nearest Neighbors and Validation



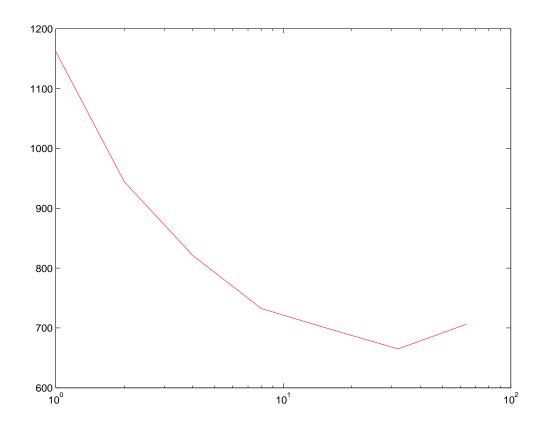
(a)



I would recommend choosing K=16 since the MSE appears to increase for the test set after this value.

(b)

Given the following MSEs:

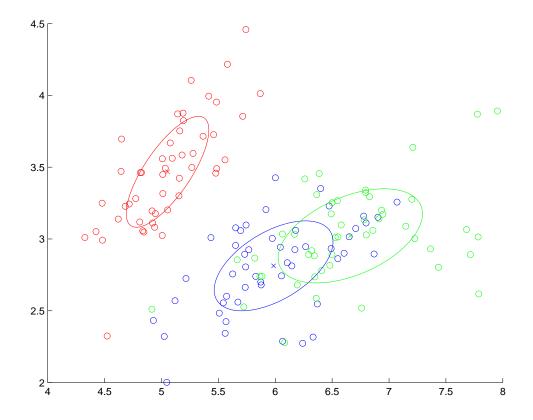


I would recommend choosing K = 32 since the MSE appears to increase for the test set after this value.

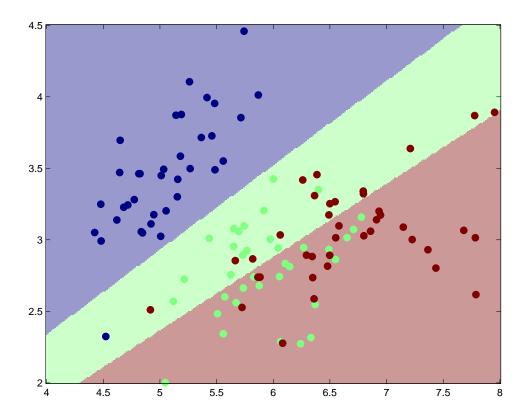
(c)

Problem 3: Bayes Classifiers

- (a) See plot in part c
- (b) See plot in part c



(c)



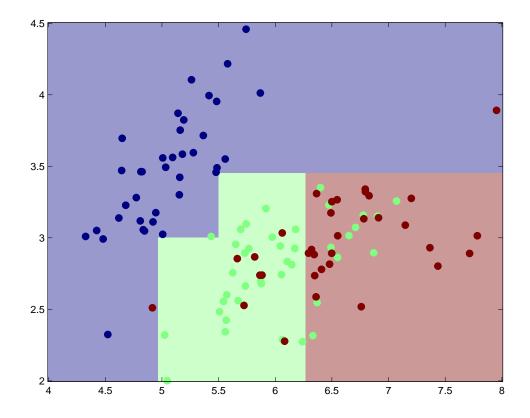
(d)

(e) training error rate = 0.2252test error rate = 0.5676

(f) training error rate = 0.0360test error rate = 0.0541

Problem 4: Decision Trees

(a) Done



(b)

- (c) For 2 features, I would choose to limit the depth to 3. After several runs, very few seem to improve the misclassification rate beyond a depth of 3 and many runs seem to actually worsen the misclassification rate beyond a depth of 3. Here, the test error comes to around 0.27.
- (d) For 4 features, I would choose to limit the depth to 3. Similarly to the 2 feature case, few runs seem to improve beyond this depth. Although very few runs actually get worse beyond this depth, many of them show an improvement between depth 2 and 3, so the extra complexity seems to usually be worth it. Here, the test error comes to between 0 and 0.1.