COMP 540 Homework 3

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1: MAP and MLE parameter estimation

1. Estimate for θ using MLE

Solution. The maximum likelihood estimation of D given θ is that

$$MLE = l(D|\theta) = p(x^{(i)}) = \prod_{i=1}^{m} \theta^{x^{(i)}} (1 - \theta)^{1 - x^{(i)}}$$

take the NLL of MLE

$$NLL = -\sum_{i}^{m} [x^{(i)}log\theta + (1 - x^{(i)})log(1 - \theta)]$$

take the derivative of NLL and make it equal to zero

$$\frac{\partial NLL}{\partial \theta} = -\sum_{i=1}^{m} [x^{(i)} \frac{1}{\theta} - \frac{1}{(1-\theta)} (1-x^{(i)})] = 0$$

by computing this equation, we can get θ_{MLE}

$$\theta_{MLE} = \frac{1}{m} \sum_{i}^{m} x^{(i)}$$

2. Compare the MAP and MLE estimates of θ

Solution. If we add a conjugate prior and use both the D and this prior to make a estimation of θ , we can have this

$$MAP = l(D|\theta)Beta(D|a,b)$$

$$= [\prod_{i}^{m} \theta^{x^{(i)}} (1-\theta)^{1-x^{(i)}}] \theta^{a-1} (1-\theta)^{b-1}$$

take the derivative of θ and make it equal to zero, we can get θ_{MAP}

$$\theta_{MAP} = \frac{\sum_{i=1}^{m} x^{(i)} + a + 1}{m + a + b - 2}$$

if a = b = 1 then

$$\theta_{MAP} = \theta_{MLE} = \frac{1}{m} \sum_{i}^{m} x^{(i)}$$

2: Logistic regression and Gaussian Naive Bayes

1. For logistic regression, what is the posterior probability for each class, i.e., P(y=1|x) and P(y=0|x)? Write the expression in terms of the parameter θ and the sigmoid function.

Solution.

$$P(y = 1|x) = h_{\theta}(X) = \frac{1}{1 + e^{-\theta^T X}}$$

$$P(y = 0|x) = 1 - h_{\theta}(X) = \frac{e^{-\theta^T X}}{1 + e^{-\theta^T X}}$$

2. Derive the posterior probabilities for each class

Solution. The Gaussian distribution and Bernoulli distribution that we assume

$$P(y=1) = \gamma$$

$$p(x_j|y=1) = N(\mu_j^1, \sigma_j^2)$$

$$P(x_j|; y=0) = N(\mu_j^0, \sigma_j^2)$$

Naive Bayes model

$$p(x|y) = \prod_{j=1}^{d} P(x_j|y)$$

Bayes rule

$$P(y = 1|x) = \frac{P(y = 1)P(x|y = 1)}{\sum P(y = i)P(x|y = i)}$$

so these are what we can use now, then we will derive the posterior probabilities

$$\begin{split} p(y=1|x) &= \frac{P(y=1)P(x|y=1)}{P(y=0)P(x|y=0) + P(y=1)P(x|y=1)} \\ &= \frac{\gamma \prod_{j=1}^d N(\mu_j^1, \sigma_j^2)}{\gamma \prod_{j=1}^d N(\mu_j^1, \sigma_j^2) + (1-\gamma) \prod_{j=1}^d N(\mu_j^0, \sigma_j^2)} \\ &= \frac{1}{1 + \frac{1-\gamma}{\gamma} \prod_{j=1}^d exp(\frac{(x_j - \mu_j^1)^2 - (x_j - \mu_j^0)^2}{2\sigma_j^2})} \end{split}$$

the probability P(y=0|x) can be derived using the same method or just subtract it from 1.

3. part 3

Solution. Class 1 and class 0 are equally likely, that means $\gamma = \frac{1}{2}$ the probability equation can be written as

$$P(y=1|x) = \frac{1}{1 + \frac{1-\gamma}{\gamma} \prod_{j=1}^{d} exp(\frac{(x_j - \mu_j^1)^2 - (x_j - \mu_j^0)^2}{2\sigma_j^2})}$$

if we set $\mu_j^0 = -\mu_j^1$ then we have

$$P(y=1|x) = \frac{1}{1 + \prod_{j=1}^{d} e^{(\frac{2x_j \mu_j^0}{\sigma_j^2})}}$$

obviously it has the same form as logistic regression, if we see in this way

$$\theta = [\frac{2\mu_1^0}{\sigma_1^2}, \frac{2\mu_2^0}{\sigma_2^2}, ..., \frac{2\mu_d^0}{\sigma_d^2}]^T$$

$$X = [x_1, x_2, ..., x_d]^T$$

the equation can be rewritten using θ and X

$$P(y=1|x) = \frac{1}{1 + e^{-\theta^T X}}$$

3: Reject option in classifiers

1. part 1

Solution. The loss of choosing a class j is

$$loss = \lambda_s(1 - P(y = j|x))$$

then if we want to decide y = j, we need to make $loss \leq \lambda_r$, which means that

$$\lambda_s(1 - P(y = j|x)) \le \lambda_r$$

so

$$P(y=j|x) \ge 1 - \frac{\lambda_r}{\lambda_s}$$

2. When $\frac{\lambda_r}{\lambda_s}$ approches to 0, the cost of falsely classifying is extremely large, which means that the cost of rejects is small, so we tend to choose rejection. As $\frac{\lambda_r}{\lambda_s}$ increases, the cost of rejection increases. And when $\frac{\lambda_r}{\lambda_s} = 1$ we would never reject.

4: Kernelizing k-nearest neighbors

Solution. In knn classifier, we just need to compute the L2 distance of a test vector to all the training points. It can be kernelized in this form:

$$||x_i - x_i|| = \langle x_i, x_i \rangle + \langle x_i, x_i \rangle - 2 \langle x_i, x_i \rangle$$

where <> denote dot product. So computing L2 distance can be turned into a 3 kenerlized computations.

5: Constructing kernels

1. $k(x, x') = Ck_1(x, x')$

Solution.

$$Ck_1(x, x') = C\Phi_1(x)^T\Phi_1(x') = (\sqrt{C}\Phi_1(x)^T)(\sqrt{C}\Phi_1(x')^T)$$

if C > 0 then k(x, x') is a valid kernel.

2. $k(x, x') = f(x)k_1(x, x')f(x')$

Solution.

$$f(x)k_{1}(x, x^{'})f(x^{'}) = \langle f(x)\Phi_{1}(x), f(x^{'})\Phi_{1}(x^{'}) \rangle$$

since it satisfy the Mercer's theorem, $k(x, x^{'})$ is valid

3. $(x, x') = k_1(x, x') + k_1(x, x')$

Solution. Because $k_1(x, x^{'})$ and $k_2(x, x^{'})$ both are valid kernels, so by Mercer's theorem they both satisfy

$$\int_{d} k(x, x^{'}) f(x) f(x^{'}) \ge 0$$

then $k(x, x') = \Phi(x)^T \Phi(x')$ exists. So

$$k_{1}(x, x^{'} + k_{2}(x, x^{'})) = \int_{d} k_{1}(x, x^{'}f(x)f(x^{'})) + \int_{d} k_{2}(x, x^{'}f(x)f(x^{'}))$$

Because $\int_{d} k_{1}(x, x'f(x)f(x')) > 0$ and $\int_{d} k_{2}(x, x'f(x)f(x')) > 0$, so

$$k_1(x, x' + k_2(x, x') \ge 0$$

6: One vs all logistic regression

1. Implementing and predicting a one-vs-all classifier for the CIFAR-10 dataset.

```
one_vs_all on raw pixels final test set accuracy: 0.361700 [[463 58 22 24 19 35 26 60 203 90]
  69 464 17 34 23 31 44
                                 95 173]
                             50
 [123
     64 193 77
                  96 89 151
                             88
                                  72
      86 78 161 48 192 171
                             51
   65
      38 103 64 234 90 194 129
                                  35
                                     481
  50 64 81 128 81 274 113 87
  31 52 67 102 86 78 457
                             51 30 46]
  53
      62 51 46 69
                     84 66 407 48 114]
 [144 79
          8 25 10 34
                         22 19 544 115]
 59 208 14 22 23 29
                             56 110 420]]
                          59
```

Figure 1: confusion matrix of ova

2. Visualizing the learned parameter matrix. The visualized coefficients are shown in 2.

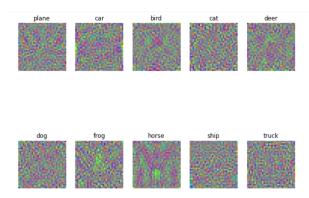


Figure 2: learned one-vs-all classifier visualization

The visualized coefficients are shown in 3.

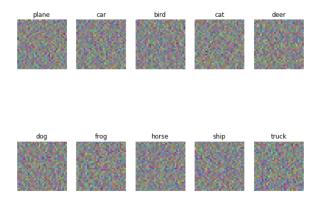


Figure 3: sklearn OVA classifier visualization

7: Softmax Regression

1. loss function for softmax regression (naive version)

Solution. When implemented in naive version, the softmax regression results are shown in below. naive loss: 2.352202e+00 computed in 14.744000s

2. loss function for softmax regression (vectorized version)

Solution. When implemented in naive version, the softmax regression results are shown in below.

Note it took 0.483000 seconds to compute the loss. vectorized loss: 2.352202e+00 computed in 0.483000s

Loss difference: 0.000000 Gradient difference: 0.000000

3. Compare naive version and vectorized version

Solution. It took 14.744000s for naive version to compute the result; and 0.483000 for vectorized version. The vectorized version is much faster than the naive version.

4. Using a validation set to select regularization lambda and learning rate for gradient descent.

Solution. We choose mini patch size 400, and set the number of iteration to 4000. The program result is: Best validation accuracy achieved during cross-validation: 0.418000

5. Evaluating the best softmax classifier on the test set and visualizing the coefficients

Solution. Using the selected best softmax model on testing dataset. The program result is: softmax on raw pixels final test set accuracy: 0.405100 Visualize the coefficients in Figure 4.

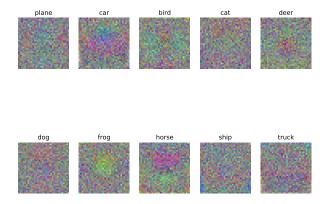


Figure 4: Visualization of the cofficients