# CPSC 340 Assignment 5 (due Friday March 20 at 11:55pm)

## Instructions

Rubric: {mechanics:5}

IMPORTANT!!! Before proceeding, please carefully read the general homework instructions at https://www.cs.ubc.ca/~fwood/CS340/homework/. The above 5 points are for following the submission instructions. You can ignore the words "mechanics", "reasoning", etc.

We use blue to highlight the deliverables that you must answer/do/submit with the assignment.

## 1 Kernel Logistic Regresion

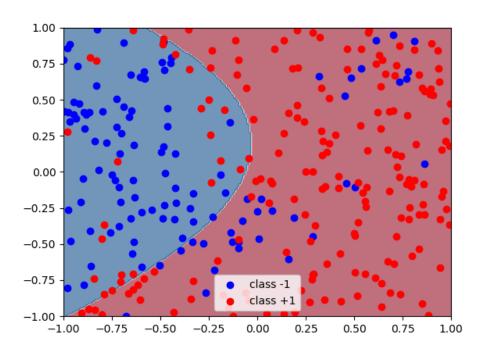
If you run python main.py -q 1 it will load a synthetic 2D data set, split it into train/validation sets, and then perform regular logistic regression and kernel logistic regression (both without an intercept term, for simplicity). You'll observe that the error values and plots generated look the same since the kernel being used is the linear kernel (i.e., the kernel corresponding to no change of basis).

## 1.1 Implementing kernels

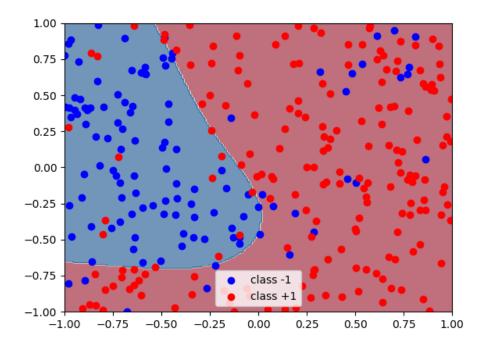
Rubric: {code:5}

Implement the polynomial kernel and the RBF kernel for logistic regression. Make sure to include the code you have written in your pdf GradeScope submission. Report your training/validation errors and submit the plots generated for each case. You should use the hyperparameters p=2 and  $\sigma=0.5$  respectively, and  $\lambda=0.01$  for the regularization strength.

Answer:
Polynomial:
Training error 0.193
Validation error 0.150



Answer: RBF: Training error 0.137 Validation error 0.100



```
elif question == "1.1":
    dataset = load_dataset('nonLinearData.pkl')
    X = dataset['X']
    y = dataset['y']

    Xtrain, Xtest, ytrain, ytest = train_test_split(X,y,random_state=0)

# kernel logistic regression with a poly kernel
    lr_kernel = kernelLogRegL2(kernel_fun=logReg.kernel_poly, lammy=1, p=2)
    lr_kernel.fit(Xtrain, ytrain)

print("Training error %.3f" % np.mean(lr_kernel.predict(Xtrain) != ytrain))
    print("Validation error %.3f" % np.mean(lr_kernel.predict(Xtest) != ytest))

utils.plotClassifier(lr_kernel, Xtrain, ytrain)

utils.savefig["logRegPolyKernel.png"]

# kernel logistic regression with a RBF kernel
    lr_kernel = kernelLogRegL2(kernel_fun=logReg.kernel_RBF, lammy=1, sigma=0.5)
    lr_kernel.fit(Xtrain, ytrain)

print("Training error %.3f" % np.mean(lr_kernel.predict(Xtrain) != ytrain))
    print("Training error %.3f" % np.mean(lr_kernel.predict(Xtrain) != ytrain))
    print("Validation error %.3f" % np.mean(lr_kernel.predict(Xtest) != ytest))

utils.plotClassifier(lr_kernel, Xtrain, ytrain)
    utils.savefig("logRegRBFKernel.png")
```

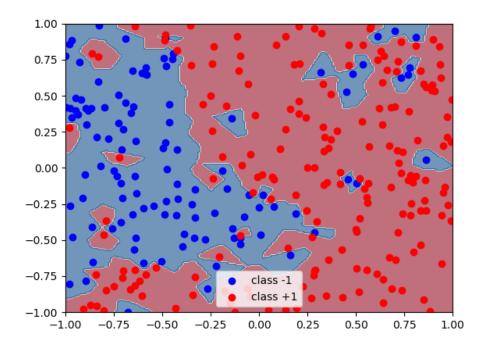
### 1.2 Hyperparameter search

Rubric: {code:3}

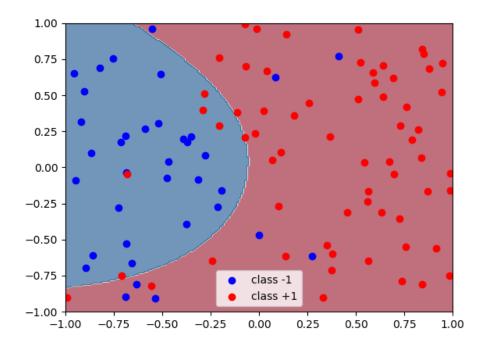
For the RBF kernel logistic regression, consider the hyperparameters values  $\sigma = 10^m$  for  $m = -2, -1, \dots, 2$  and  $\lambda = 10^m$  for  $m = -4, -3, \dots, 0$ . In main.py, sweep over the possible combinations of these hyperparameter values. Make sure to include the code you have written in your pdf GradeScope submission. Report the hyperparameter values that yield the best training error and the hyperparameter values that yield the best validation error. Include the plot for each.

Note: on the job you might choose to use a tool like scikit-learn's GridSearchCV to implement the grid search, but here we are asking you to implement it yourself by looping over the hyperparameter values.

Answer: Training error 0.000, lamda = 0.010, sigma = 0.000



Answer: Validation error 0.130, lamda = 1.000, sigma = 0.100



#### 1.3 Reflection

Rubric: {reasoning:1}

Briefly discuss the best hyperparameters you found in the previous part, and their associated plots. Was the training error minimized by the values you expected, given the ways that  $\sigma$  and  $\lambda$  affect the fundamental tradeoff?

Answer: Yes, when the value of hyperparameter increases, traing error increases and validation error decreases.

## 2 MAP Estimation

Rubric: {reasoning:8}

In class, we considered MAP estimation in a regression model where we assumed that:

- The likelihood  $p(y_i \mid x_i, w)$  is a normal distribution with a mean of  $w^T x_i$  and a variance of 1.
- The prior for each variable j,  $p(w_j)$ , is a normal distribution with a mean of zero and a variance of  $\lambda^{-1}$ .

Under these assumptions, we showed that this leads to the standard L2-regularized least squares objective function,

$$f(w) = \frac{1}{2} ||Xw - y||^2 + \frac{\lambda}{2} ||w||^2,$$

which is the negative log likelihood (NLL) under these assumptions (ignoring an irrelevant constant). For each of the alternate assumptions below, show how the loss function would change (simplifying as much as possible):

1. We use a Laplace likelihood with a mean of  $w^T x_i$  and a scale of 1, and we use a zero-mean Gaussian prior with a variance of  $\sigma^2$ .

$$p(y_i \mid x_i, w) = \frac{1}{2} \exp(-|w^T x_i - y_i|), \quad p(w_j) = \frac{1}{\sqrt{2\pi}\sigma} \exp\left(-\frac{w_j^2}{2\sigma^2}\right).$$

Answer:

$$-\sum \log(p(w_j)) = \frac{1}{2\sigma^2} \sum_{i=1}^n w_j^2 = \frac{1}{2\sigma^2} ||w||^2$$

$$f(w) = \frac{1}{2} ||Xw - y||_1 + \frac{1}{2\sigma^2} ||w||^2$$

2. We use a Gaussian likelihood where each datapoint has its own variance  $\sigma_i^2$ , and where we use a zero-mean Laplace prior with a vairance of  $\lambda^{-1}$ .

$$p(y_i \mid x_i, w) = \frac{1}{\sqrt{2\sigma_i^2 \pi}} \exp\left(-\frac{(w^T x_i - y_i)^2}{2\sigma_i^2}\right), \quad p(w_j) = \frac{\lambda}{2} \exp(-\lambda |w_j|).$$

You can use  $\Sigma$  as a diagonal matrix that has the values  $\sigma_i^2$  along the diagonal.

Answer:

$$-\sum_{i=1}^{n} \log(\exp(-\frac{(w^T x_i - y_i)^2}{2\sigma_i^2})) = \frac{1}{2\sigma^2} \sum_{i=1}^{n} (w^T x_i - y_i)^2 = \frac{1}{2\sigma^2} ||Xw - y||^2$$

$$f(w) = \frac{1}{2\sigma^2} ||Xw - y||^2 + \lambda ||w||_1$$

3. We use a (very robust) student t likelihood with a mean of  $w^T x_i$  and  $\nu$  degrees of freedom, and a zero-mean Gaussian prior with a variance of  $\lambda^{-1}$ ,

$$p(y_i|x_i, w) = \frac{\Gamma\left(\frac{\nu+1}{2}\right)}{\sqrt{\nu\pi}\Gamma\left(\frac{\nu}{2}\right)} \left(1 + \frac{(w^Tx_i - y_i)^2}{\nu}\right)^{-\frac{\nu+1}{2}}, \quad p(w_j) = \frac{\sqrt{\lambda}}{\sqrt{2\pi}} \exp\left(-\lambda \frac{w_j^2}{2}\right).$$

where  $\Gamma$  is the "gamma" function (which is always non-negative).

Answer:

$$-\sum_{i=1}^{n} \log p(y_i|x_i, w) = -\sum_{i=1}^{n} \log \left(\frac{\Gamma\left(\frac{\nu+1}{2}\right)}{\sqrt{\nu\pi}\Gamma\left(\frac{\nu}{2}\right)} \left(1 + \frac{(w^T x_i - y_i)^2}{\nu}\right)^{-\frac{\nu+1}{2}}\right)$$

$$= -\sum_{i=1}^n \log \frac{\Gamma\left(\frac{\nu+1}{2}\right)}{\sqrt{\nu\pi}\Gamma\left(\frac{\nu}{2}\right)} + \frac{\nu+1}{2}\sum_{i=1}^n \log \left(1 + \frac{(w^Tx_i - y_i)^2}{\nu}\right)$$

$$f(w) = \frac{v+1}{2} \sum_{i=1}^{n} \log \left( 1 + \frac{(w^{T} x_{i} - y_{i})^{2}}{\nu} \right) + \frac{\lambda}{2} ||w||^{2}$$

4. We use a Poisson-distributed likelihood (for the case where  $y_i$  represents counts), and we use a uniform prior for some constant  $\kappa$ ,

$$p(y_i|w^Tx_i) = \frac{\exp(y_iw^Tx_i)\exp(-\exp(w^Tx_i))}{y_i!}, \quad p(w_j) \propto \kappa.$$

(This prior is "improper" since  $w \in \mathbb{R}^d$  but it doesn't integrate to 1 over this domain, but nevertheless the posterior will be a proper distribution.)

Answer:

$$f(w) = -\sum_{i=1}^{n} \log p(y_i|x_i, w) = \sum_{i=1}^{n} [-y_i w^T x_i + \exp(w^T x_i)]$$

## 3 Principal Component Analysis

Rubric: {reasoning:3}

Consider the following dataset, containing 5 examples with 2 features each:

$$\begin{array}{c|cc} x_1 & x_2 \\ \hline -4 & 3 \\ 0 & 1 \\ -2 & 2 \\ 4 & -1 \\ 2 & 0 \\ \end{array}$$

Recall that with PCA we usually assume that the PCs are normalized (||w|| = 1), we need to center the data before we apply PCA, and that the direction of the first PC is the one that minimizes the orthogonal distance to all data points.

- 1. What is the first principal component?
- 2. What is the reconstruction loss (L2 norm squared) of the point (-3, 2.5)? (Show your work.)
- 3. What is the reconstruction loss (L2 norm squared) of the point (-3, 2)? (Show your work.)

Answer:

- 1. 1st PCA:  $w = [1 \ 0]$
- 2.  $z = w^T x = \begin{bmatrix} 1 \ 0 \end{bmatrix}^T \begin{bmatrix} -3 \ 2.5 \end{bmatrix} = \begin{bmatrix} -3 \ 0 \end{bmatrix} f(W, Z) = \|w^T z x\| = \|\begin{bmatrix} 1 \ 0 \end{bmatrix}^T \begin{bmatrix} -3 \ 0 \end{bmatrix} \begin{bmatrix} -3 \ 2.5 \end{bmatrix}\| = \|\begin{bmatrix} 0 \ -2.5 \end{bmatrix}\| = 6.25$

3. 
$$z = w^T x = \begin{bmatrix} 1 & 0 \end{bmatrix}^T \begin{bmatrix} -3 & 2 \end{bmatrix} = \begin{bmatrix} -3 & 0 \end{bmatrix} f(W, Z) = \|w^T z - x\| = \|\begin{bmatrix} 1 & 0 \end{bmatrix}^T \begin{bmatrix} -3 & 0 \end{bmatrix} - \begin{bmatrix} -3 & 2 \end{bmatrix}\| = \|\begin{bmatrix} 0 & -2 \end{bmatrix}\| = 4$$

Hint: it may help (a lot) to plot the data before you start this question.

## 4 PCA Generalizations

## 4.1 Robust PCA

Rubric: {code:10}

If you run python main -q 4.1 the code will load a dataset X where each row contains the pixels from a single frame of a video of a highway. The demo applies PCA to this dataset and then uses this to reconstruct the original image. It then shows the following 3 images for each frame:

- 1. The original frame.
- 2. The reconstruction based on PCA.
- 3. A binary image showing locations where the reconstruction error is non-trivial.

Recently, latent-factor models have been proposed as a strategy for "background subtraction": trying to separate objects from their background. In this case, the background is the highway and the objects are the cars on the highway. In this demo, we see that PCA does an OK job of identifying the cars on the highway in that it does tend to identify the locations of cars. However, the results aren't great as it identifies quite a few irrelevant parts of the image as objects.

Robust PCA is a variation on PCA where we replace the L2-norm with the L1-norm,

$$f(Z,W) = \sum_{i=1}^{n} \sum_{j=1}^{d} |\langle w^j, z_i \rangle - x_{ij}|,$$

and it has recently been proposed as a more effective model for background subtraction. Complete the class pca.RobustPCA, that uses a smooth approximation to the absolute value to implement robust PCA. Make sure to include the code you have written in your pdf GradeScope submission. Briefly comment on the results.

Note: in its current state, pca.RobustPCA is just a copy of pca.AlternativePCA, which is why the two rows of images are identical.

Hint: most of the work has been done for you in the class *pca.AlternativePCA*. This work implements an alternating minimization approach to minimizing the (L2) PCA objective (without enforcing orthogonality). This gradient-based approach to PCA can be modified to use a smooth approximation of the L1-norm. Note

that the log-sum-exp approximation to the absolute value may be hard to get working due to numerical issues, and a numerically-nicer approach is to use the "multi-quadric" approximation:

$$|\alpha| \approx \sqrt{\alpha^2 + \epsilon},$$

where  $\epsilon$  controls the accuracy of the approximation (a typical value of  $\epsilon$  is 0.0001).

Answer:

```
class RobustPCA:
       def __init__(self, k):
    self.k = k
      def fit(self, X):
    n,d = X.shape
    k = self.k
    self.mu = np.mean(X,0)
    X = X - self.mu
             # Randomly initial Z, W
z = np.random.randn(n*k)
             w = np.random.randn(k*d)
             checkRPCAGrad(self, z, w, X, k)
              f = np.sum(np.abs(np.dot(z.reshape(n, k), w.reshape(k, d))-X))
              for i in range(50):
                    f_old = f
                    z = find_min(self._fun_obj_z, z, 10, False, w, X, k)
w = find_min(self._fun_obj_w, w, 10, False, z, X, k)
f = np.sum(np.abs(np.dot(z.reshape(n, k), w.reshape(k, d))-X))
print('Iteration {:2d}, loss = {}'.format(i, f))
                     if f_old - f < 1e-4:
             self.W = w.reshape(k, d)
return self
       def compress(self, X):
             n,d = X.shape

k = self.k

X = X - self.mu

# We didn't enforce that W was orthogonal so we need to optimize to find Z
             z = np.zeros(n*k)
z = find_min(self._fun_obj_z, z, 100, False, self.W.flatten(), X, k)
Z = z.reshape(n,k)
      def expand(self, Z):
   X = np.dot(Z, self.W) + self.mu
   return X
```

```
def expand(self, Z):
    X = np.dot(Z, self.W) + self.mu
    return X
def _fun_obj_z(self, z, w, X, k, epsilon=0.0001):
    n,d = X.shape
    Z = z.reshape(n,k)
      = w.reshape(k,d)
    R = np.dot(Z, W) - X
    D = np.sqrt(R**2 + epsilon)
       np.sum(D)
    R[:] = R[:] / D[:]
    g = np.dot(R, W.T)
    return f, g.flatten()
def _fun_obj_w(self, w, z, X, k, epsilon=0.0001):
    n,d = X.shape
    Z = z.reshape(n, k)
    W = w.reshape(k, d)
    R = np.dot(Z, W) - X
      = np.sqrt(np.power(R, 2) + epsilon)
       np.sum(D)
    R[:] = R[:] / D[:]
    g = np.dot(Z.T, R)
      turn f, g.flatten()
```

#### 4.2 Reflection

Rubric: {reasoning:3}

1. Briefly explain why using the L1 loss might be more suitable for this task than L2.

Answer: L1 can pin some coefficient values to 0.

2. How does the number of video frames and the size of each frame relate to n, d, and/or k?

Answer: They increase as the values of parameters increase

3. What would the effect be of changing the threshold (see code) in terms of false positives (cars we identify that aren't really there) and false negatives (real cars that we fail to identify)?

Answer: When the threshold increases, loss decreases.

## 5 Very-Short Answer Questions

Rubric: {reasoning:11}

1. Assuming we want to use the original features (no change of basis) in a linear model, what is an advantage of the "other" normal equations over the original normal equations?

Answer: Have less error and run faster

2. In class we argued that it's possible to make a kernel version of k-means clustering. What would an advantage of kernels be in this context?

Answer: kernel k-means partitions the points by linear separators in the new space

3. In the language of loss functions and regularization, what is the difference between MLE and MAP?

Answer: MLE starts only with the probability of observation given the parameter, while MAP also take prior distribution into consideration.

4. What is the difference between a generative model and a discriminative model?

Answer: Generative model explicitly models the actual distribution of each class, but discriminative model learns the conditional probability distribution

5. With PCA, is it possible for the loss to increase if k is increased? Briefly justify your answer.

Answer: Yes, when k increases, there are more dimensions and loss could increase.

6. What does "label switching" mean in the context of PCA?

Answer: Get same model if we permute rows of W

7. Why doesn't it make sense to do PCA with k > d?

Answer: when k > d, the least squares estimates tend to also have high variance, and hence the model will perform badly.

8. In terms of the matrices associated with PCA  $(X, W, Z, \hat{X})$ , where would an "eigenface" be stored? Answer: In X, each face/image is a element

9. What is an advantage and a disadvantage of using stochatic gradient over SVD when doing PCA?

Answer: Advanrage: stochatic gradient doesn't force any orthogonality constraints. Disadvatange: can be expensive when haing too many features.

- 10. Which of the following step-size sequences lead to convergence of stochastic gradient to a stationary point?
  - (a)  $\alpha^t = 1/t^2$ .
  - (b)  $\alpha^t = 1/t$ .
  - (c)  $\alpha^t = 1/\sqrt{t}$ .
  - (d)  $\alpha^t = 1$ .

Answer: a

11. We discussed "global" vs. "local" features for e-mail classification. What is an advantge of using global features, and what is advantage of using local features?

Answer: Global features can make prediction to larger population. Local feature can make precise prediction on a specific group.