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1a) X and Y are not independent, knowing the value of X or Y affects the probability of the other variable. For example, if we know that X has a value of 1, then the variable Y has to have a value of zero. If they were independent, they would satisfy the relationship:

$$P(x|y) = P(x)$$
 and  $P(y|x) = P(y)$ 

Since P(y = 0|x = 1) = 1 and  $P(y = 0) = \frac{1}{2}$ , we can see that the two variables are not independent. We can, however, show that X and Y are uncorrelated. Two random variables are uncorrelated if cov(X,Y) = 0.

$$cov(X,Y) = E(XY) - E(X)E(Y)$$

$$P(x=1) = \frac{1}{4}$$

$$P(x=0) = \frac{1}{2}$$

$$P(x = -1) = \frac{1}{4}$$

$$P(y=1) = \frac{1}{4}$$

$$P(y=0) = \frac{1}{2}$$

$$P(y = -1) = \frac{1}{4}$$

$$P(x = 1, y = 0) = \frac{1}{4}$$

$$P(x = -1, y = 0) = \frac{1}{4}$$

$$P(x = 0, y = 1) = \frac{1}{4}$$

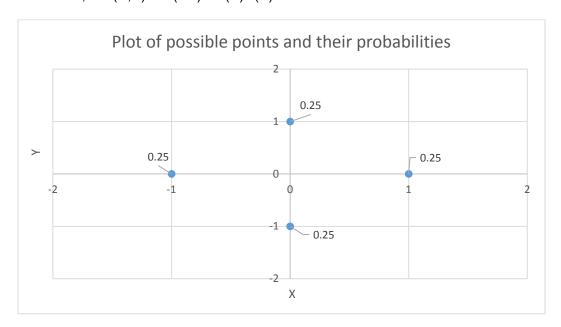
$$P(x = 0, y = -1) = \frac{1}{4}$$

$$E(X) = P(x = 1)*1 + P(x = 0)*0 + P(x = -1)*-1 = 0$$

Similarly, E(Y) = 0

E(XY) = 0, since either X or Y will be 0

Therefore, cov(X,Y) = E(XY) - E(X)E(Y) = 0



1b)

B1,B2,B3 = 0,0,0 -> X,Y,Z = 0,0,0  
B1,B2,B3 = 1,0,0 -> X,Y,Z = 1,0,1  
B1,B2,B3 = 0,1,0 -> X,Y,Z = 1,1,0  
B1,B2,B3 = 0,0,1 -> X,Y,Z = 0,1,1  
B1,B2,B3 = 1,1,0 -> X,Y,Z = 0,1,1  
B1,B2,B3 = 1,0,1 -> X,Y,Z = 1,1,0  
B1,B2,B3 = 0,1,1 -> X,Y,Z = 1,0,1  
B1,B2,B3 = 1,1,1 -> X,Y,Z = 0,0,0  

$$P(X = 1) = \frac{1}{2} P(X = 0) = \frac{1}{2}$$
  
 $P(Y = 1) = \frac{1}{2} P(Y = 0) = \frac{1}{2}$   
 $P(Z = 1) = \frac{1}{2} P(Z = 0) = \frac{1}{2}$ 

## Consider the pair X,Y:

$$P(X = 1, Y = 1) = \frac{1}{4} = P(X=1)P(Y=1)$$
  
 $P(X = 0, Y = 1) = \frac{1}{4} = P(X=0)P(Y=1)$   
 $P(X = 0, Y = 0) = \frac{1}{4} = P(X=0)P(Y=0)$   
 $P(X = 1, Y = 0) = \frac{1}{4} = P(X=1)P(Y=0)$ 

So P(X,Y) = P(X)P(Y), meaning X and Y are independent.

From symmetry of the problem, we can then see that this is true for the pairs Y,Z and X,Z as well, so X,Y,Z are pairwise independent, i.e. :

$$P(X,Y) = P(X)P(Y)$$

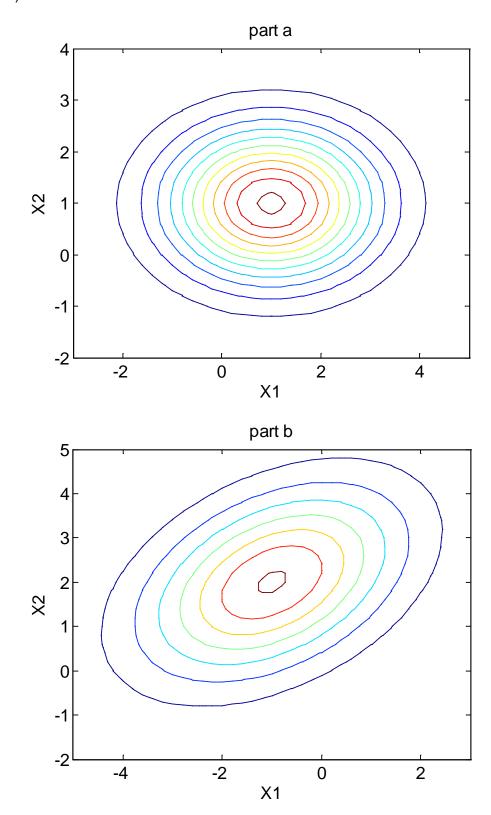
$$P(Y,Z) = P(Y)P(Z)$$

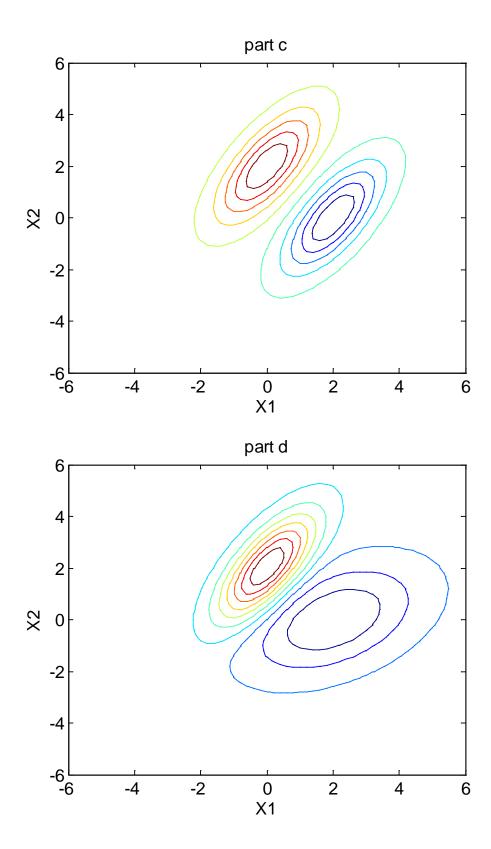
$$P(X,Z) = P(X)P(Z)$$

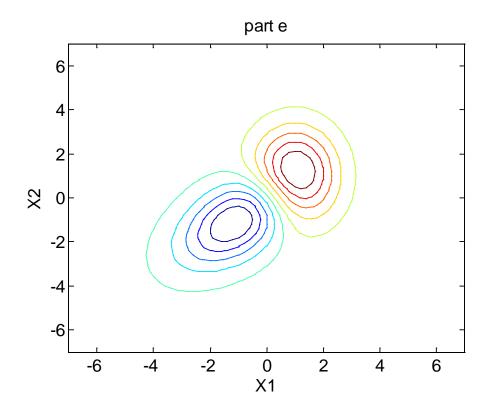
X,Y,Z are not mutually independent since P(X,Y,Z) = P(X)P(Y)P(Z).

## For example:

$$P(X = 1, Y = 1, Z = 1) = 0$$
  
 $P(X = 1)P(Y=1) P(Z=1) = 1/8$   
 $P(X=1, Y=1, Z=1) = P(X=1)P(Y=1)P(Z=1)$ 





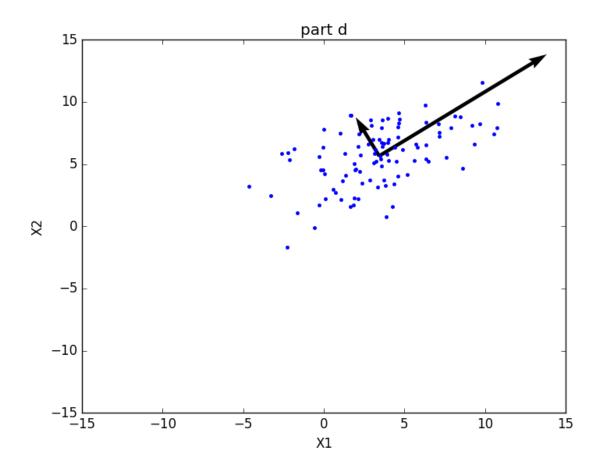


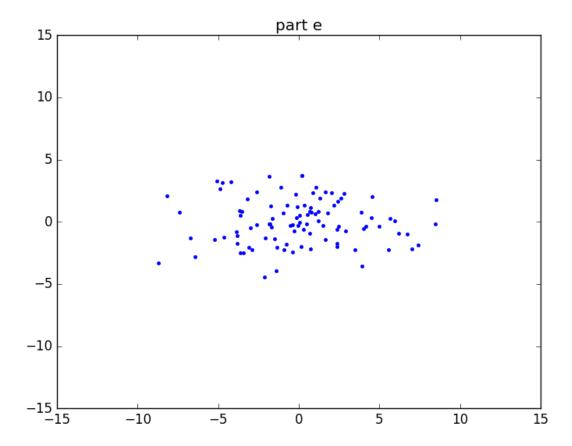
3) a) mean: [3.137, 5.698]

b) covariance matrix:  $\begin{bmatrix} 8.209 & 3.721 \\ 3.721 & 5.694 \end{bmatrix}$ 

c) eigenvalues: [10.879 3.024]

eigenvectors:  $\left[\begin{array}{c} 0.812 \\ -0.583 \end{array}\right]$  ,  $\left[\begin{array}{c} 0.583 \\ 0.812 \end{array}\right]$ 





4) a) 
$$\Sigma = E((\mathbf{x} - \mathbf{\mu})^{\mathrm{T}}(\mathbf{x} - \mathbf{\mu}))$$
  
For all  $a \in \mathbb{R}^n$ ,  $\mathbf{a}^{\mathrm{T}}\Sigma \mathbf{a} = E(\mathbf{a}^T(\mathbf{x} - \mathbf{\mu})^{\mathrm{T}}(\mathbf{x} - \mathbf{\mu})\mathbf{a}) \ge 0$ , so  $\mathbf{a}^{\mathrm{T}}\Sigma \mathbf{a} \ge 0$ .

Since  $a^T\Sigma a \geq 0$ , we know that  $\Sigma$  is positive semidefinite. In homework 2, we showed that all positive definite matrices are invertible. Therefore, we know that  $\Sigma^{-1}$  will not exist when  $\Sigma$  is positive semidefinite but not positive definite (one or more of the eigenvalues is 0 and one or more of the diagonal entries is 0). In order to ensure that  $\Sigma^{-1}$  exists, we must transform X into X' such that the diagonals of  $\Sigma$  are always positive. Essentially, we need to find X' such that  $cov(X_i, X_i) = E((x_i - \mu_i)^2) > 0$ . We can linearly transform X by adding a positive constant  $\gamma$  to each  $X_i$  which will give us an X' that ensures an invertible covariance matrix.

b) 
$$x^{T}\Sigma^{-1}x = ||Ax||_{2}^{2} = x^{T}A^{T}Ax \rightarrow A^{T}A = \Sigma^{-1}$$

From the spectral decomposition theorem, we know that we can express  $\Sigma^{-1}$  as  $U\Lambda^{-1}U^T$ , where U is an array with the eigenvectors of  $\Sigma^{-1}$  as columns and  $\Lambda$  is a diagonal matrix with the eigenvalues of  $\Sigma^{-1}$  as diagonal entries. Let B be a diagonal matrix with the square roots of the eigenvalues of  $\Sigma^{-1}$  as diagonal entries.

$$A^{T}A = \Sigma^{-1} = U\Lambda^{-1}U^{T} = U(B^{-1})^{T}B^{-1}U^{T} = (B^{-1}U^{T})^{T}B^{-1}U^{T}$$
  
Therefore,  $A = B^{-1}U^{T}$ .

- c) The intuitive meaning of  $x^T\Sigma^{-1}x$  is a measure of how far the point x is from the mean of the Gaussian.
- d) The maximum value of  $\|Ax\|_2^2$  is the largest eigenvalue of  $\Sigma^{\text{-1}}$  and the minimum value is the smallest eigenvalue. If  $X_i$  and  $X_j$  are orthogonal, then they are independent and the intuitive meaning for the maximum and minimum of  $\|Ax\|_2^2$  is a measure of the higher variance and lower variance corresponding to  $X_i$  and  $X_j$ . From the equation for f(x), we can see that minimizing  $\|Ax\|_2^2$  will maximize f(x), so we will choose x corresponding to the smallest eigenvalue of  $\Sigma^{\text{-1}}$ .

5) a) The maximum likelihood estimates for the mean and covariance matrix of a Gaussian distribution are simply the mean and covariance matrix of the sample data. For i.i.d observations  $X_1, ..., X_n$  from the same class, the mean will have the same dimensions as  $X_i$  and the elements will be the averages of each feature that is observed. The covariance matrix will have entries  $\Sigma_{ij}$  which correspond to the covariance between feature i and feature j across all n samples. For class k, the mean and covariance matrix are given by:

$$\hat{\mu}_k = \frac{1}{n_k} \sum_{j \mid Y_j = k} X_j \qquad \hat{\Sigma} = \frac{1}{n} \sum_k \sum_{j \mid Y_j = k} (X_j - \hat{\mu}_k) (X_j - \hat{\mu}_k)^T$$

b) The prior distribution for each class i is given by (number of training samples in class i)/(total samples). Using all 60,000 training samples, we find the priors to be:

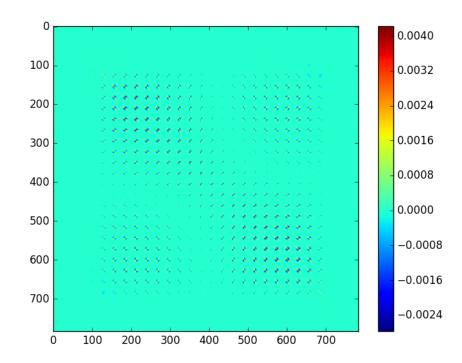
class 0: 0.0987 class 1: 0.1124 class 2: 0.0993 class 3: 0.1022

class 4: 0.0974 class 5: 0.0904

class 6: 0.0986 class 7: 0.1044 class 8: 0.0975

class 9: 0.0992

c) This is the covariance matrix for the digit 1.

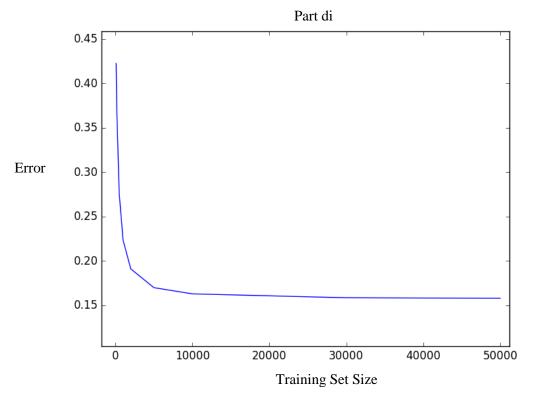


The diagonal entries of the matrix represent the variances of each feature. From this diagonal, we can see that the variance is highest where people would generally start writing and where they would finish writing. This intuitively makes sense, as people tend to draw their ones at different angles and different lengths. However, the midsection of a one usually passes through the center of the image regardless of the angle the one is drawn at. This can be seen in the structure of the covariance matrix, as the covariance matrix values decrease in the middle of the matrix.

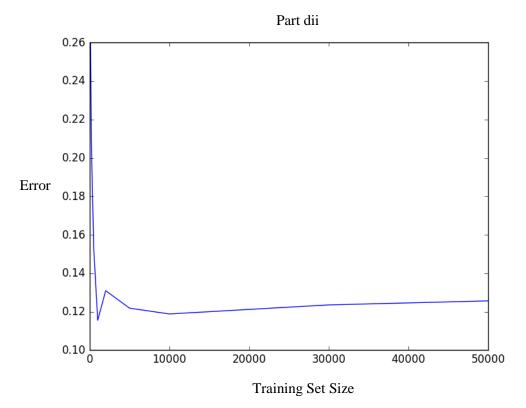
d) i) Here we are using linear discriminant analysis to classify our images. We find the likelihood of each class for each sample and then classify the sample with the class of highest likelihood. The decision boundary in this case separates our 784-dimensional space into 10 sections (one for each digit). The implementation is a linear discriminant analysis classifier, so the decision boundary takes a linear form. The boundary takes on this form because the same covariance matrix was used for each of the Gaussians associated with each digit.

Error rates: [0.4386, 0.3777, 0.2911, 0.2301, 0.1936, 0.1731, 0.1664, 0.1615, 0.1604] These error rates correspond to the training set sizes:

[100, 200, 500, 1000, 2000, 5000, 10000, 30000, 50000]



ii) In this part, we have implemented quadratic discriminant analysis. Our algorithm for classifying is essentially the same as that of LDA, but by allowing each class to have its own covariance matrix, we create quadratic decision boundary surfaces. Error rates: [0.2600, 0.2082, 0.1529, 0.1155, 0.1309, 0.1218, 0.1188, 0.1235, 0.1256] These error rates correspond to the training set sizes: [100, 200, 500, 1000, 2000, 5000, 10000, 30000, 50000]



iii) The error rate in part (ii) is much lower than that of part (i) for small training set sizes. The decision boundaries are linear in part (i) and quadratic in part (ii). As a result, the decision boundary chosen in part (ii) is more complex and results in better performance, especially for small training set sizes where underfitting is more significant. This increased decision boundary can also result in more overfitting, which can be seen by the slightly increasing error rate in part (ii).

iv) QDA validation error: 0.1296. QDA Kaggle score: 0.88320 LDA validation error: 0.1598 LDA Kaggle score: 0.85320

e) validation error: 0.2005 Kaggle score: 0.74240 6.1) Want to find w,  $\alpha$  that minimizes:  $\Sigma_n (X_i w + \alpha - y_i)^2$ . We can insert a column of ones to the end of X to absorb the  $\alpha$  term.

Find w that minimizes  $|Xw - y|^2$   $\rightarrow$  Find w that minimizes RSS =  $(Xw - y)^T(Xw - y)$ 

 $(Xw-y)^T(Xw-y)$  is minimized when  $X^TXw = X^Ty$   $\rightarrow$  If  $X^TX$  is invertible,  $w = (X^TX)^{-1}X^Ty$ 

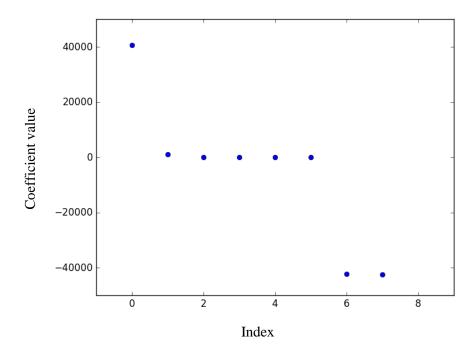
Using the training data, we find that w = [4.06e+04, 1.20e+03, -8.50, 1.18e+02, -3.79e+01, 4.31e+01, -4.22e+04, -4.25e+04, -3.57e+06]

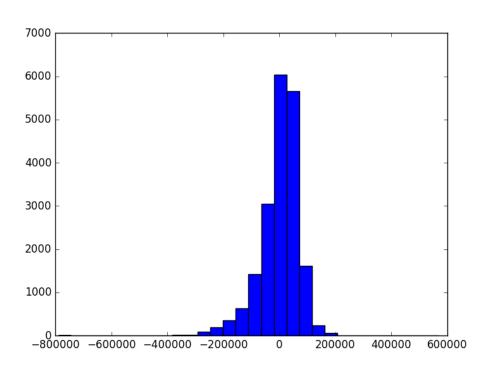
6.2) RSS: 5.795e+12

predicted min: -56562.827545 predicted max: 710798.838689

These numbers make sense. Dividing the RSS by the number of samples and then taking the square root gives us the average error per data point, which comes out to be \$69,492. This is a reasonable amount of error for the value of a home. The actual range for the validation set is \$28,300 to \$500,001. The actual range differs significantly from the predicted range, which suggests that the median home value is not linear at the high and low ends of the spectrum.

6.3)





The distribution resembles a Gaussian.

## Code for problem 2 (MATLAB)

```
% Part (a)
mu = [1 1];
Sigma = [2 0; 0 1];
x1 = -3:.2:5; x2 = -2:.2:4;
[X1,X2] = meshgrid(x1,x2);
F = mvnpdf([X1(:) X2(:)], mu, Sigma);
F = reshape(F,length(x2),length(x1));
figure()
contour(X1,X2,F)
set(gca,'FontSize', 14)
title('part a')
xlabel('X1')
ylabel('X2')
% Part (b)
mu = [-1 \ 2];
Sigma = [3 1;1 2];
x1 = -5:.2:3; x2 = -2:.2:5;
[X1,X2] = meshgrid(x1,x2);
F = mvnpdf([X1(:) X2(:)],mu,Sigma);
F = reshape(F,length(x2),length(x1));
figure()
contour(X1,X2,F)
set(gca,'FontSize', 14)
title('part b')
xlabel('X1')
ylabel('X2')
% Part (c)
mu1 = [0 2];
Sigmal = [1 1;1 2];
mu2 = [2 0];
Sigma2 = [1 1;1 2];
x1 = -6:.2:6; x2 = -6:.2:6;
[X1,X2] = meshgrid(x1,x2);
F1 = mvnpdf([X1(:) X2(:)],mu1,Sigma1);
F2 = mvnpdf([X1(:) X2(:)],mu2,Sigma2);
F3 = F1 - F2;
F3 = reshape(F3, length(x2), length(x1));
figure()
contour(X1,X2,F3,10)
set(gca,'FontSize', 14)
title('part c')
xlabel('X1')
ylabel('X2')
% Part (d)
mu1 = [0 2];
Sigma1 = [1 1;1 2];
mu2 = [2 0];
Sigma2 = [3 1;1 2];
x1 = -6:.2:6; x2 = -6:.2:6;
[X1,X2] = meshgrid(x1,x2);
F1 = mvnpdf([X1(:) X2(:)],mu1,Sigma1);
```

```
F2 = mvnpdf([X1(:) X2(:)],mu2,Sigma2);
F3 = F1 - F2;
F3 = reshape(F3, length(x2), length(x1));
figure()
contour(X1,X2,F3,10)
set(gca,'FontSize', 14)
title('part d')
xlabel('X1')
ylabel('X2')
% Part (e)
mu1 = [1 1];
Sigma1 = [1 0; 0 2];
mu2 = [-1 -1];
Sigma2 = [2 1;1 2];
x1 = -7:.2:7; x2 = -7:.2:7;
[X1,X2] = meshgrid(x1,x2);
F1 = mvnpdf([X1(:) X2(:)],mu1,Sigma1);
F2 = mvnpdf([X1(:) X2(:)],mu2,Sigma2);
F3 = F1 - F2;
F3 = reshape(F3, length(x2), length(x1));
figure()
contour(X1,X2,F3,10)
set(gca,'FontSize', 14)
title('part e')
xlabel('X1')
ylabel('X2')
```

## Code for problems 3, 5, 6 (Python)

import matplotlib.mlab as mlab import numpy as np import matplotlib.pyplot as plt import math, random import scipy.io

```
# Problem 3
"
X1 = np.random.normal(3,3,100)
X2 = (X1/2) + np.random.normal(4,2,100)
# Part (a)
mu = [np.sum(X1)/100, np.sum(X2)/100]
print('mean:\n', mu)
# Part (b)
sigma = np.cov(np.vstack((X1,X2)))
print('covariance matrix:\n',sigma)
```

```
# Part (c)
eigen = np.linalg.eig(sigma)
eigvalues = eigen[0]
eigvectors = eigen[1].T
print('eigenvalues:\n',eigvalues)
print('eigenvectors:\n',eigvectors)
# Part (d)
plt.figure()
plt.plot(X1,X2,'.')
plt.xlim(-15,15)
plt.ylim(-15,15)
U,V = zip(*eigvectors)
U = np.multiply(U,eigvalues)
V = np.multiply(V,eigvalues)
ax = plt.gca()
ax.quiver(mu[0], mu[1], U,V,scale_units='xy',scale=1)
plt.title('part d')
plt.xlabel("X1")
plt.ylabel("X2")
# Part (e)
if eigvalues[0] > eigvalues[1]:
  U = np.array(eigvectors).T
else:
  U = np.array([eigvectors[1],eigvalues[0]]).T
rotated = np.dot(U.T,(np.vstack([X1,X2]) - np.tile(mu,[100,1]).T))
plt.figure()
plt.plot(rotated[0],rotated[1],'.')
plt.xlim(-15,15)
plt.ylim(-15,15)
plt.title('part e')
plt.show()
# Problem 5
#benchmark.m, converted
def benchmark(pred_labels, true_labels):
  errors = pred_labels != true_labels
  err_rate = sum(errors) / float(len(true_labels))
  indices = errors.nonzero()
  return err_rate, indices
```

```
#montage images.m, converted
def montage_images(images):
  num_images=min(1000,np.size(images,2))
  numrows=math.floor(math.sqrt(num images))
  numcols=math.ceil(num images/numrows)
  img=np.zeros((numrows*28,numcols*28));
  for k in range(num_images):
     r = k % numrows
     c = k // numrows
     img[r*28:(r+1)*28,c*28:(c+1)*28]=images[:,:,k];
  return img
digit_train_data = scipy.io.loadmat("data/digit_dataset/train.mat")
digit train images= digit train data["train images"]
digit_train_labels= digit_train_data["train_labels"]
train_vectors=[]
for i in range(np.shape(digit_train_images)[2]):
  vector = np.float64(digit train images[:,:,i].flatten())
  train vectors.append(np.divide(vector,np.linalg.norm(vector, 2)))
normalized vectors= np.array(train vectors)
indices = list(range(np.shape(normalized vectors)[0]))
random.shuffle(indices)
train vectors = []
train_labels = []
val vectors = []
val labels = []
for i in indices[:50000]:
  train_vectors.append(normalized_vectors[i])
  train_labels.append(digit_train_labels[i])
for i in indices[50000:]:
  val vectors.append(normalized vectors[i])
  val labels.append(digit train labels[i])
val_labels = np.array(val_labels).ravel()
train_labels = np.array(train_labels).ravel()
train_vectors = np.array(train_vectors)
val_vectors = np.array(val_vectors)
# part (a)
mean list = []
cov_list = []
for i in range(10):
  indices = np.where(digit_train_labels==i)[0]
  subset = normalized_vectors[indices]
  mean list.append(np.sum(subset,axis=0)/subset.shape[0])
  cov list.append(np.cov(subset.T))
# mean list and cov list are the mean and covariance matrices for part (a)
```

```
# mean_list[i] and cov_list[i] correspond to digit i
# part (b)
priors = []
for i in range(10):
  indices = np.where(digit_train_labels==i)[0]
  priors.append(indices.shape[0]/digit_train_labels.shape[0])
print('Priors:\n', priors)
# part (c)
plt.imshow(cov_list[1])
plt.colorbar()
plt.show()
# part (di)
training_size = [100,200,500,1000,2000,5000,10000,30000,50000]
errors = []
alpha = np.eye(784)*0.0001
for N in training_size:
  subset_vectors = train_vectors[:N]
  subset_labels = train_labels[:N]
  mean_list = []
  cov_list = []
  for i in range(10):
     indices = np.where(subset labels==i)[0]
     subset = subset_vectors[indices]
     mean_list.append(np.sum(subset,axis=0)/subset.shape[0])
     cov_list.append(np.cov(subset.T))
  priors = []
  for i in range(10):
     indices = np.where(subset_labels==i)[0]
     priors.append(indices.shape[0]/subset_labels.shape[0])
  sigma_overall = np.sum(cov_list,axis=0)/len(cov_list) + alpha
  sig_inv = np.linalg.inv(sigma_overall)
  predictions=[]
  for sample in val_vectors:
     guess = 0
     best = -float('inf')
     for i in range(10):
```

```
f = (np.dot(np.dot(mean_list[i].T,sig_inv),sample) -
np.dot(np.dot(mean_list[i].T,sig_inv),mean_list[i])/2 + math.log(priors[i]))
       if f > best:
          quess = i
          best = f
     predictions.append(quess)
  errors.append(benchmark(np.array(predictions),val_labels)[0])
plt.figure()
plt.plot(training_size,errors,'-')
plt.show()
# part (dii)
training\_size = [100,200,500,1000,2000,5000,10000,30000,50000]
errors = []
alpha = np.eye(784)*0.00001
for N in training size:
  randomidx = random.randint(0,50000-N)
  subset vectors = train vectors[randomidx:randomidx+N]
  subset_labels = train_labels[randomidx:randomidx+N]
  mean_list = []
  cov_list = []
  for i in range(10):
     indices = np.where(subset_labels==i)[0]
     subset = subset_vectors[indices]
     mean_list.append(np.sum(subset,axis=0)/subset.shape[0])
     cov_list.append(np.cov(subset.T) + alpha)
  priors = []
  for i in range(10):
     indices = np.where(subset labels==i)[0]
     priors.append(indices.shape[0]/subset_labels.shape[0])
  sig_inv = [np.linalg.inv(sigma) for sigma in cov_list]
  logdet_list = [np.linalg.slogdet(sigma)[1] for sigma in cov_list]
  predictions=[]
  for sample in val vectors:
     guess = 0
     best = -float('inf')
     for i in range(10):
       f = np.dot(np.dot((sample - mean_list[i]).T,sig_inv[i]),(sample - mean_list[i]))/-2 +
math.log(priors[i]) - logdet_list[i]/2
       if f > best:
          guess = i
          best = f
     predictions.append(guess)
```

```
errors.append(benchmark(np.array(predictions),val_labels)[0])
plt.figure()
plt.plot(training_size,errors,'-')
plt.show()
# part (div)
# LDA for part (div)
digit_test_data = scipy.io.loadmat("data/digit_dataset/test.mat")
digit_test_images= digit_test_data["test_images"].transpose()
test vectors=[]
for i in range(np.shape(digit_test_images)[1]):
  vector = np.float64(digit_test_images[:,i])
  vector = vector.reshape((28,28)).T
  vector = vector.flatten()
  test_vectors.append(np.divide(vector,np.linalg.norm(vector, 2)))
test vectors= np.array(test vectors)
alpha = np.eye(784)*0.0001
subset vectors = train vectors
subset_labels = train_labels
mean list = []
cov_list = []
for i in range(10):
  indices = np.where(subset_labels==i)[0]
  subset = subset_vectors[indices]
  mean_list.append(np.sum(subset,axis=0)/subset.shape[0])
  cov list.append(np.cov(subset.T))
priors = []
for i in range(10):
  indices = np.where(subset_labels==i)[0]
  priors.append(indices.shape[0]/subset_labels.shape[0])
sigma overall = np.sum(cov list,axis=0)/len(cov list) + alpha
sig_inv = np.linalg.inv(sigma_overall)
predictions=[]
for sample in val_vectors:
  guess = 0
  best = -float('inf')
  for i in range(10):
     f = (np.dot(np.dot(mean list[i].T,sig inv),sample) -
np.dot(np.dot(mean_list[i].T,sig_inv),mean_list[i])/2 + math.log(priors[i]))
     if f > best:
       guess = i
```

```
best = f
  predictions.append(guess)
error = benchmark(np.array(predictions),val_labels)[0]
test predict = []
for sample in test_vectors:
  quess = 0
  best = -float('inf')
  for i in range(10):
     f = (np.dot(np.dot(mean list[i].T,sig inv),sample) -
np.dot(np.dot(mean_list[i].T,sig_inv),mean_list[i])/2 + math.log(priors[i]))
     if f > best:
       quess = i
       best = f
  test_predict.append(guess)
numbers = (np.arange(10000) + 1)
print(test_predict[:25])
test predict = np.vstack((numbers,test predict))
print(error)
np.savetxt("digitsLDA.csv", test_predict.transpose(), delimiter=",",fmt = '%u')
img = montage_images(test_vectors.T.reshape((28,28,test_vectors.shape[0]))[;,;;25])
plt.imshow(img)
plt.show()
# QDA for part (div)
digit_test_data = scipy.io.loadmat("data/digit_dataset/test.mat")
digit test images= digit test data["test images"].transpose()
test vectors=[]
for i in range(np.shape(digit_test_images)[1]):
  vector = np.float64(digit_test_images[:,i])
  vector = vector.reshape((28,28)).T
  vector = vector.flatten()
  test_vectors.append(np.divide(vector,np.linalg.norm(vector, 2)))
test_vectors= np.array(test_vectors)
alpha = np.eye(784)*0.00001
mean_list = []
cov_list = []
```

```
for i in range(10):
  indices = np.where(train_labels==i)[0]
  subset = train_vectors[indices]
  mean list.append(np.sum(subset,axis=0)/subset.shape[0])
  cov list.append(np.cov(subset.T) + alpha)
priors = []
for i in range(10):
  indices = np.where(train_labels==i)[0]
  priors.append(indices.shape[0]/train_labels.shape[0])
sig inv = [np.linalg.inv(sigma) for sigma in cov list]
logdet list = [np.linalg.slogdet(sigma)[1] for sigma in cov list]
validation predict = []
for sample in val_vectors:
  guess = 0
  best = -float('inf')
  for i in range(10):
     f = np.dot(np.dot((sample - mean_list[i]).T,sig_inv[i]),(sample - mean_list[i]))/-2 +
math.log(priors[i]) - logdet list[i]/2
     if f > best:
       guess = i
       best = f
  validation_predict.append(guess)
error = benchmark(np.array(validation_predict),val_labels)[0]
test_predict = []
for sample in test vectors:
  quess = 0
  best = -float('inf')
  for i in range(10):
     f = np.dot(np.dot((sample - mean_list[i]).T,sig_inv[i]),(sample - mean_list[i]))/-2 +
math.log(priors[i]) - logdet_list[i]/2
     if f > best:
       quess = i
       best = f
  test_predict.append(guess)
numbers = (np.arange(10000) + 1)
print(test predict[:25])
test_predict = np.vstack((numbers,test_predict))
print(benchmark(validation predict, val labels)[0])
np.savetxt("digitsQDA.csv", test_predict.transpose(), delimiter=",",fmt = '%u')
img = montage_images(test_vectors.T.reshape((28,28,test_vectors.shape[0]))[;,;;25])
plt.imshow(img)
plt.show()
```

```
# part (e)
spam data = scipy.io.loadmat("data/spam dataset/spam data.mat")
spam train data= spam data["training data"]
spam train labels= np.ravel(spam data["training labels"])
spam_test = spam_data["test_data"]
normalize = []
for vector in spam test:
  if not np.linalg.norm(vector, 2):
     normalize.append(vector)
  else:
     normalize.append(np.divide(vector,np.linalg.norm(vector, 2)))
spam_test = np.array(normalize)
normalize = []
for vector in spam train data:
  if not np.linalg.norm(vector, 2):
     normalize.append(vector)
  else:
     normalize.append(np.divide(vector,np.linalg.norm(vector, 2)))
spam train data = np.array(normalize)
alpha = np.eye(32)*0.0001
indices = list(range(np.shape(spam_train_data)[0]))
random.shuffle(indices)
subset_vectors = []
subset labels = []
val_vectors = []
val labels = []
for i in indices[:4000]:
  subset vectors.append(spam train data[i])
  subset_labels.append(spam_train_labels[i])
for i in indices[4000:]:
  val vectors.append(spam_train_data[i])
  val_labels.append(spam_train_labels[i])
val_labels = np.array(val_labels).ravel()
subset labels = np.array(subset labels).ravel()
subset_vectors = np.array(subset_vectors)
val_vectors = np.array(val_vectors)
mean_list = []
cov_list = []
priors = []
for i in range(2):
  indices = np.where(subset labels==i)[0]
  subset = subset vectors[indices]
  mean_list.append(np.sum(subset,axis=0)/subset.shape[0])
```

```
cov list.append(np.cov(subset.T) + alpha)
  priors.append(indices.shape[0]/subset_labels.shape[0])
sig inv = [np.linalg.inv(sigma) for sigma in cov list]
logdet_list = [np.linalg.slogdet(sigma)[1] for sigma in cov_list]
predictions=[]
for sample in val_vectors:
  quess = 0
  best = -float('inf')
  for i in range(2):
     f = np.dot(np.dot((sample - mean_list[i]).T,sig_inv[i]),(sample - mean_list[i]))/-2 +
math.log(priors[i]) - logdet_list[i]/2
     if f > best:
       quess = i
       best = f
  predictions.append(guess)
error = benchmark(np.array(predictions),val_labels)[0]
test predict = []
for sample in spam test:
  guess = 0
  best = -float('inf')
  for i in range(2):
     f = np.dot(np.dot((sample - mean_list[i]).T,sig_inv[i]),(sample - mean_list[i]))/-2 +
math.log(priors[i]) - logdet_list[i]/2
     if f > best:
       quess = i
       best = f
  test_predict.append(guess)
numbers = (np.arange(5857) + 1)
print('val error: ',error)
test_predict = np.vstack((numbers,test_predict))
np.savetxt("spam.csv", test_predict.transpose(), delimiter=",",fmt = '%u')
# problem 6
housing data = scipy.io.loadmat("data/housing dataset/housing data.mat")
X train= housing data['Xtrain']
X = np.hstack((X_train,np.ones((X_train.shape[0],1))))
Y_train= housing_data['Ytrain']
y = np.float_(Y_train)[:,0]
w = np.dot(np.linalg.inv(np.dot(X.T,X)),np.dot(X.T,y))
X validate = housing data['Xvalidate']
X_{validate} = np.hstack((X_{validate}, np.ones((X_{validate}.shape[0], 1))))
Y_validate= housing_data['Yvalidate'][:,0]
```

```
 \begin{tabular}{ll} validation\_predict = np.dot(X\_validate,w) \\ RSS = np.dot((validation\_predict - Y\_validate).T,(validation\_predict - Y\_validate)) \\ print('RSS: ',RSS) \\ print('predicted min: ',min(validation\_predict)) \\ print('predicted max: ',max(validation\_predict)) \\ plt.figure() \\ plt.plot(np.arange(8),w[:8],'o') \\ plt.xlim((-1,9)) \\ plt.figure() \\ plt.hist(np.dot(X,w) - y,30) \\ plt.show() \\ \end{tabular}
```