1 Independence vs Correlation

(a) Answer: X and Y are uncorrelated, but they are not independent.

Prove uncorrelated:

Two random variables X, Y are uncorrelated if their covariance E(XY) - E(X)E(Y) is 0.

$$\begin{split} E(X) &= 0 \cdot P(X=0) + 1 \cdot P(X=1) + (-1) \cdot P(X=-1) \\ &= P(X=1) - P(X=-1) \\ &= P(X=1,Y=0) + P(X=1,Y=1) - P(X=-1,Y=0) - P(X=-1,Y=1) \\ &= P(X=1,Y=0) - P(X=-1,Y=0) \\ &= P(X=1|Y=0) \cdot P(Y=0) - P(X=-1|Y=0) \cdot P(Y=0) \\ &= \frac{1}{2} \cdot \frac{1}{2} - \frac{1}{2} \cdot \frac{1}{2} = 0 \\ \text{Similarly } E(Y) &= 0 \end{split}$$

Similarly, E(Y) = 0.

Because either X or Y is zero, X and Y are never nonzero at the same time, so E(XY) = 0.

Therefore, we have proved that E(XY) = E(X)E(Y) = 0

Prove dependent:

Two random variables X, Y are independent if P(X|Y) = P(X).

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

P(X = 1) = P(X = 1, Y = 0) = P(X = 1|Y = 0) · P(Y = 0) = $\frac{1}{2}$ · $\frac{1}{2}$ = $\frac{1}{4}$
Therefore, X, Y are not independent.

(b) Answer: X, Y, Z are pairwise independent but not mutually independent. Prove pairwise independent:

For P(X|Y) we are given Y which put some constraints on C and D, between which only C influences X. However, B randomly chooses between 0 and 1 and it is independent of C, so knowing Y won't influence P(X). Thus P(X|Y) = P(X).

Similar applies to other pairs.

Prove not mutually independent:

 $P(X|Y,Z) \neq P(X)$ because if given Y = 1 and Z = 1, we can deduce that B = C = D, so that X is definitely = 1 too. Thus Y, Z influence X.

2/20/2017 hw3_code

```
In [1]: import numpy as np
   import matplotlib.pyplot as plt
   import scipy.io
   from scipy.stats import multivariate_normal
   import math
   %matplotlib inline
```

2. Isocontours of Normal Distributions

Reference: https://docs.scipy.org/doc/scipy-

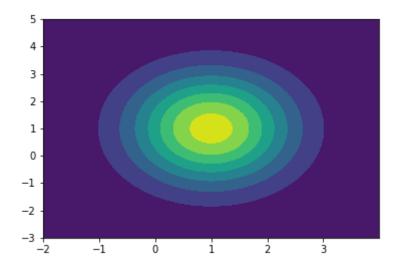
<u>0.14.0/reference/generated/scipy.stats.multivariate_normal.html (https://docs.scipy.org/doc/scipy-0.14.0/reference/generated/scipy.stats.multivariate_normal.html)</u>

```
In [2]: def isocontours(mu, sig, x, y):
            pos = np.empty(x.shape + (2,))
            pos[:, :, 0] = x
            pos[:, :, 1] = y
            rv = multivariate_normal(mu, sig)
            fig = plt.figure()
            plot = fig.add_subplot(1.2,1.7,1)
            return plot.contourf(x, y, rv.pdf(pos))
        def two_isocontours(mu1, mu2, sig1, sig2, x, y):
            pos = np.empty(x.shape + (2,))
            pos[:, :, 0] = x
            pos[:, :, 1] = y
            rv1 = multivariate normal(mu1, sig1)
            rv2 = multivariate normal(mu2, sig2)
            fig = plt.figure()
            plot = fig.add subplot(1.2,1.7,1)
            return plot.contourf(x, y, rv1.pdf(pos) - rv2.pdf(pos))
```

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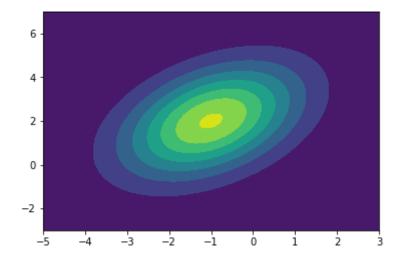
```
In [3]: # (a)
    mu = [1, 1]
    sig = [[1,0],[0,2]]
    x, y = np.mgrid[-2:4:.01, -3:5:0.01]
    a = isocontours(mu, sig, x, y)
    a
```

Out[3]: <matplotlib.contour.QuadContourSet at 0x107e885f8>



```
In [4]: # (b)
mu = [-1, 2]
sig = [[2,1],[1,3]]
x, y = np.mgrid[-5:3:.01, -3:7:.01]
b = isocontours(mu, sig, x, y)
b
```

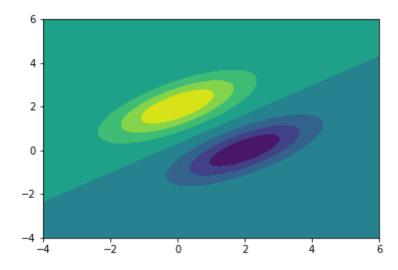
Out[4]: <matplotlib.contour.QuadContourSet at 0x10cae7e10>



2/20/2017 hw3_code

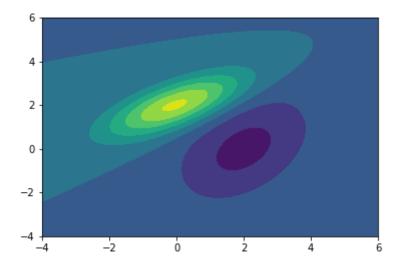
```
In [5]: # (c)
    mu1 = [0, 2]
    mu2 = [2, 0]
    sig = [[2,1],[1,1]]
    x, y = np.mgrid[-4:6:.01, -4:6:.01]
    c = two_isocontours(mu1, mu2, sig, sig, x, y)
    c
```

Out[5]: <matplotlib.contour.QuadContourSet at 0x1112f0438>



```
In [6]: # (d)
    mu1 = [0, 2]
    mu2 = [2, 0]
    sig1 = [[2,1],[1,1]]
    sig2 = [[2, 1], [1, 3]]
    x, y = np.mgrid[-4:6:.01, -4:6:.01]
    d = two_isocontours(mu1, mu2, sig1, sig2, x, y)
    d
```

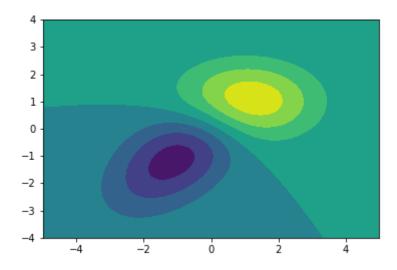
Out[6]: <matplotlib.contour.QuadContourSet at 0x1143b0588>



2/27/2017 hw3_code

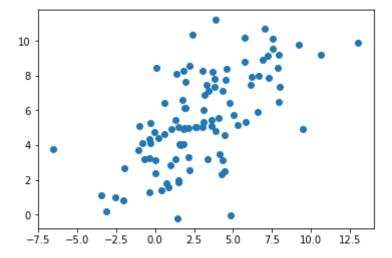
```
In [7]: # (3)
    mu1 = [1, 1]
    mu2 = [-1, -1]
    sig1 = [[2, 0], [0, 1]]
    sig2 = [[2, 1], [1, 2]]
    x, y = np.mgrid[-5:5:.01, -4:4:.01]
    e = two_isocontours(mu1, mu2, sig1, sig2, x, y)
    e
```

Out[7]: <matplotlib.contour.QuadContourSet at 0x11611ac50>



3. Eigenvectors of the Gaussian Covariance Matrix

```
In [18]: # randomly draw 100 sample points
n = 100
x1 = np.random.normal(3, 3, n)
x2 = 0.5 * x1 + np.random.normal(4, 2, n)
fig, axes = plt.subplots()
axes.scatter(x1, x2)
plt.show()
```



2/27/2017 hw3_code

```
In [19]: # (a)
         x1_{mean} = np.mean(x1)
         x2_{mean} = np.mean(x2)
         print("x1 mean = ", x1_mean)
         print("x2 mean = ", x2 mean)
         x1 mean = 2.95932073974
         x2 mean = 5.49994510123
In [20]: # (b)
         m = np.cov(x1, x2)
         print("Covariance Matrix = \n", m)
         Covariance Matrix =
          [[ 10.43613528
                          5.47600768]
             5.47600768
                           7.34461899]]
In [21]:
         # (C)
         evalue, evectors = np.linalg.eig(m)
         print("Eigenvalues: \n", evalue)
         print("Eigenvectors: \n", evectors)
         Eigenvalues:
          [ 14.58037084
                           3.200383441
         Eigenvectors:
          [[ 0.79739029 -0.60346394]
          [ 0.60346394  0.79739029]]
In [26]: # (d)
         arrow1 = evalue[0] * evectors[:,0]
         arrow2 = evectors[:, 1] * evalue[1]
         # Plot
         fig, axes = plt.subplots()
         axes.scatter(x1, x2)
         plt.quiver(x1_mean, x2_mean, arrow1[0], arrow1[1], angles='xy', scale_units=
         plt.quiver(x1_mean, x2_mean, arrow2[0], arrow2[1], angles='xy', scale_units=
         plt.axis([-15,15, -15,15])
         plt.show()
           15
           10
            5
            0
           -5
          -10
```

10

15

-15

-15

-10

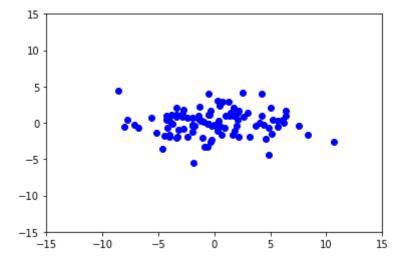
-5

2/27/2017 hw3_code

```
In [27]: # (e)
    x1_centered = np.reshape(np.subtract(x1, x1_mean), (len(x1), 1))
    x2_centered = np.reshape(np.subtract(x2, x2_mean), (len(x2), 1))
    x_centered = np.hstack([x1_centered, x2_centered])
    x_rotated = []
    for ptr in x_centered:
        new_ptr = np.matrix(evectors).T * np.reshape(ptr, (2, 1))
        x_rotated.append([new_ptr[0, 0], new_ptr[1, 0]])

x_rotated = np.array(x_rotated)
    x1 = x_rotated[:, 0]
    x2 = x_rotated[:, 1]

plt.axis([-15, 15, -15, 15])
    plt.plot(x1, x2, 'bo')
    plt.show()
```



In []:

4 Maximum Likelihood Estimation

(a) pdf for multivariate normal distribution $N(\mu, \Sigma)$:

$$p(x|\mu, \Sigma) = \frac{1}{(2\pi)^{n/2} \Sigma^{1/2}} e^{-\frac{1}{2}(x-\mu)^T \Sigma^{-1}(x-\mu)}$$

$$L(\mu, \Sigma|X) = \prod_{i=1}^n P(X = x_i) = \left(\frac{1}{(2\pi)^{n/2} \Sigma^{1/2}}\right)^n e^{\left(\frac{1}{2}\Sigma^{-1}\right)^n \sum_{i=1}^n (x_i - \mu)^T (x_i - \mu)}$$

$$\log(L(\mu, \Sigma|X)) = -n \log((2\pi)^{n/2} \Sigma^{1/2}) - \frac{1}{2} \sum_{i=1}^n (x_i - \mu)^T \Sigma^{-1}(x_i - \mu)$$

$$= -n \log((2\pi)^{n/2}) - \frac{n}{2} \log(\Sigma) - \frac{1}{2} \sum_{i=1}^n (x_i - \mu)^T \Sigma^{-1}(x_i - \mu)$$

Compute $\hat{\mu}$:

$$\frac{d \log L}{d\mu} = \frac{1}{2} [2c_1(x_1 - \mu) + 2c_2(x_2 - \mu)...2c_n(x_n - \mu)] = \sum_{i=1}^n x_i - n\mu = 0$$

$$\hat{\mu} = \frac{1}{n} \sum_{i=1}^n x_i$$

Compute $\hat{\Sigma}$:

According to the "trace trick" I found on CMU lecture slides (https://www.cs.cmu.edu/epxing/Class/08s/recitation/gaussian.pdf):

$$L(\mu, \Sigma | X) \propto -\frac{n}{2} \log(\Sigma) - \frac{1}{2} Trace(\Sigma^{-1} \sum_{i=1}^{n} (x_i - \mu)(x_i - \mu)^T)$$

Because $\frac{d}{dA} \log A = A^{-T}$ and $\frac{d}{dA} Tr(AB) = \frac{d}{dA} Tr(BA) = B^{T}$:

$$\frac{d\log L}{d\Sigma} = -\frac{n}{2}\Sigma + \frac{1}{2}\sum_{i=1}^{n} (x_i - \mu)(x_i - \mu)^T = 0$$
$$\hat{\Sigma} = \frac{1}{n}\sum_{i=1}^{n} (x_i - \hat{\mu})^T (x_i - \hat{\mu})$$

(b) MLE for $A\mu = \frac{1}{n} \sum_{i=1}^{n} x_i$ Because A is invertible

$$A^{-1}A\mu = \hat{\mu} = \frac{A^{-1}}{n} \sum_{i=1}^{n} x_i$$

5 Covariance Matirces and Decompositions

(a) Answer: Σ is not invertible when 1. there is one or more than one deterministic-value features; 2. there is one or more than one features that are dependent on others.

A matrix is invertible if all of its eigenvalues $\neq 0$

Case 1: ≥ 1 deterministic features

If a feature f_i always equals to certain value, $Cov(f_i, f_j)$ and $Cov(R_j, R_i)$ for j = 1, 2...n all equal 0, so there is a whole row of 0 and a whole column of 0 in the covariance matrix. This means one of the matrix's eigenvalue is 0. Geometrically, all sample points will have the same value for feature f_i , which means they are "squeezed" on a R^{n-1} -dimension space instead of R^n -dimension, and the eigenvector corresponding to f_i is 0. For example in a 2D space with features x and y, and y is constant, then all the sample points will spread on a horizontal line. If we plot eigenvector arrows on the plot, we won't see one that points in y-axis direction.

Case 2: > 1 features are dependent on other features

This means some eigenvectors might be dependent on others. Geometrically, if we assume there are 2 features x and y and y is dependent on x in a way that all y values are a certain multiple of x, we will see sample points only spread on a line. The eigenvector of y will overlap with that of x although might with a different magnitude.

(b) **Approach 1**: delete all dependent features which means delete the corresponding rows and columns that concern these features

Approach 2: to make the eigenvalue not 0, we can add a really small γ constant to it.

$$Ax = x$$

$$I\gamma x = \beta x$$

$$Ax + I\gamma x = \beta x + \gamma x$$

$$(A + I\gamma)x = (\beta + \gamma)x$$

So we just add $I\gamma$ to the covariance matrix, and we can use accuracy on validation dataset to tune this γ value.

(c) To maximize f(x), choose the eigenvector that corresponds to the maximum eigenvalue; to minimize f(x), choose the eigenvector that corresponds to the minimum eigenvalue.

$$f(x) = \frac{1}{(2\pi)^{n/2} \sum^{1/2}} e^{-\frac{1}{2}(x-\mu)^T \sum^{-1} (x-\mu)}$$

To maximize f(x), we need to minimize $\frac{1}{2}(x-\mu)^T\Sigma^{-1}(x-\mu)$. Because $\mu=0$, so we are minimizing $x^T\Sigma^{-1}x$

Becuase Σ is positive definite and invertible,

$$\Sigma = PDP^T$$

where P is a unitary matrix consist of orthonormal eigenvectors and D is diagonal matrix consist of eigenvalues λ_i . Therefore, $P^{-1} = P^T$ and D^{-1} is also diagonal matrix consist of σ eigenvalue's inverse.

$$\Sigma^{-1} = (PDP^T)^{-1} = (P^T)^{-1}D^{-1}P^{-1} = PD^{-1}P^T$$

Set S as a diagonal matrix consist of the square root of D^{-1} 's diagonal entries s.t. $D^{-1} = SS$

$$\Sigma^{-1} = PD^{-1}P^T = PSSP^T = PSS^TP^T$$

Set A = PS, then

$$\Sigma^{-1} = A^{T} A$$

$$x^{T} \Sigma^{-1} x = ||Ax||_{2}^{2}$$

$$||Ax||_{2}^{2} = x^{T} A^{T} A x = x^{T} P D^{-1} P^{T} x$$

Set v = Px. Because P is unitary and $||x||_2^2 = 1$, we choose any x vector that is in the Euclidean basis vector form (consist of one "1" entry and rest are all "0"). And x can be transformed by P to another Euclidean basis vector. In other words, different x can single out different eigenvectors in P.

$$x^{T} \Sigma^{-1} x = ||Ax||_{2}^{2} = v^{T} D^{-1} v$$

The above means we can choose a v vector to single out any diagonal entries on D^{-1} . As a reminder, D^{-1} 's diagonal consist of the inverse of Σ^{-1} 's eigenvalues, $\frac{1}{\lambda_i}$. Therefore, in order to minimize $||Ax||_2^2$, we need $\frac{1}{\lambda_{max}}$ So we need to choose x that can single out the eigenvector that corresponds to the maximal eigenvalue of Σ^{-1} .

Minimizing f(x) has opposite process, which is to find x that singles out the eigenvector that corresponds to the minimum eigenvalue.

MNIST

```
In [1]: import matplotlib.pyplot as plt
import scipy.io as sio
from scipy.stats import multivariate_normal
import numpy as np
import pandas as pd
from sklearn.preprocessing import normalize
%matplotlib inline
import pdb
```

Load Data

```
In [6]: trainMAT = sio.loadmat('./mnist/train.mat')
    testMAT = sio.loadmat('./mnist/test.mat')
    trainX = trainMAT["trainX"]
    trainY = trainX[:, -1:].reshape(1, len(trainX))[0]
    np.random.shuffle(trainX)
    validate_data = trainX[:10000, :]
    train_data = trainX[10000:, :]
    test_data = testMAT["testX"]
    CLASS_ = [0, 1, 2, 3, 4, 5, 6, 7, 8, 9]
    c = 0.0001
```

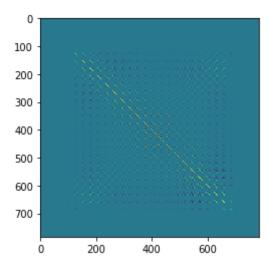
a) Fit Gaussian Distribution to each digit class using MLE

```
In [3]: def gaussian_mean_cov(data, label, if_print=False):
    df = pd.DataFrame(data)
    df = df[df[784]== label]
    features = normalize(df.values[:, :-1].astype(np.float32))
    mu = np.mean(features, axis=0)
    sigma = np.cov(features, rowvar=0)
    if if_print:
        print("Label ", label, " : ")
        print("Mu = ", mu)
        print("Sigma = ", sigma, "\n")
    return mu, sigma
```

b) Visualize covariance matrix

In [5]: plt.imshow(gaussian_mean_cov(train_data, 0)[1])

Out[5]: <matplotlib.image.AxesImage at 0x104163860>



Observation:

- 1. the values on one diagonal are comparably large. These values are the variances of the RVs, so it's reasonable that they are larger than any feature pair's covariance
- 2. There are parts of the values super small (the faded color on the graph), which means the features involved don't have much dependency on each other, or they might be always 0 (no color on this pixel)

c) LDA & QDA

```
In [15]: def lda_train(data):
             models = dict()
             x = normalize(data[:, :-1].astype(np.float32))
             sigma = np.cov(x, rowvar=0)
             for label in CLASS:
                 mu, _ = gaussian_mean_cov(data, label)
                 m = multivariate_normal(mu, sigma + c * np.identity(sigma.shape[0]))
                 models[label] = m
             return models
         def qda train(data):
             models = dict()
             for label in CLASS:
                 mu, sigma = gaussian mean cov(data, label)
                 m = multivariate_normal(mu, sigma + c * np.identity(sigma.shape[0]))
                 models[label] = m
             return models
         def test(x, models):
             y = list()
             x = normalize(x.astype(np.float32))
             for sample in x:
                 prob = [models[label].logpdf(sample) + compute prior(trainY, label)
                 y.append(np.argmax(prob))
             return y
         def batch train and evaluate(train data, validate data, categories, type="LI
             errors = []
             for size in categories:
                 models = qda train(train data[:size, :])
                 if type == "LDA":
                     models = lda_train(train_data[:size, :])
                 prediction = test(validate data[:, :-1], models)
                 y = validate data[:, -1:].reshape(1, len(validate data))[0]
                 err = 1 - np.mean(np.equal(prediction, y).astype(np.int32))
                 errors.append(err)
                 print("Error for training size {}: {}".format(size, err))
             plt.plot(categories, errors, 'ro')
             plt.axis([min(categories)-10, max(categories)+10, -0.1, max(errors)+0.1]
             plt.title("{} Error rate".format(type))
             return models
         def compute prior(data, label):
             return np.mean(np.equal(data, label).astype(np.int32))
```

Spam

```
In [1]: import matplotlib.pyplot as plt
   import scipy.io as sio
   from scipy.stats import multivariate_normal
   import numpy as np
   import pandas as pd
   import math
   from sklearn.preprocessing import normalize
   from sklearn.feature_extraction.text import TfidfVectorizer
%matplotlib inline
   import glob as g
   import re
   import pdb
```

Load Data

```
In [2]: spamfiles = g.glob('./spam/spam/*.txt')
    hamfiles = g.glob('./spam/ham/*.txt')
    testfiles = ['./spam/test/' + str(i) + '.txt' for i in range(10000)]
```

```
In [3]: def load_process_files(files):
    txts = list()
    for file in files:
        txt = open(file, "r", encoding='utf-8', errors='ignore').read()
        txt = txt.replace('\r\n', ' ')
        txts.append(txt)
    return txts
```

```
In [4]: spams = load_process_files(spamfiles)
hams = load_process_files(hamfiles)
tests = load_process_files(testfiles)
trains = spams + hams
all = trains + tests
```

```
In [5]: Ns = len(spams); Nh = len(hams); Nt = len(tests)
    validate_n = 10000
    vectorizer = TfidfVectorizer(min_df=0.05)

data = vectorizer.fit_transform(all).toarray()
    train_data = data[:-10000]
    test_data = data[-10000:]

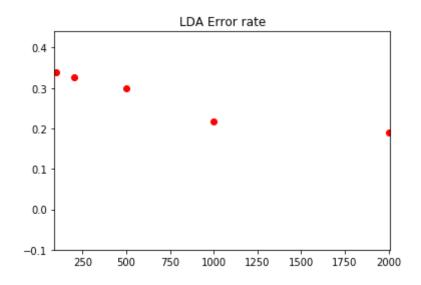
spam_data = train_data[:Ns]
    spam_validate = spam_data[validate_n:]

ham_data = train_data[Ns:]
    ham_validate = ham_data[validate_n:]

spam_prior = math.log(1.0 * Ns / (Ns + Nh))
    ham_prior = math.log(1.0 * Nh / (Ns + Nh))
```

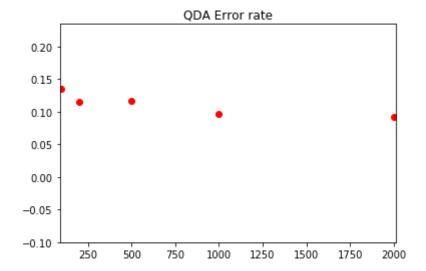
Model

```
In [15]: def gaussian_mean_cov(data):
             mu = np.mean(data, axis=0)
             sigma = np.cov(data, rowvar=0)
             return mu, sigma
         def lda_train(spam, ham, all):
             x = normalize(all.astype(np.float32))
             sigma = np.cov(x, rowvar=0)
             mu1, _ = gaussian_mean_cov(spam)
             mu2, _ = gaussian_mean_cov(ham)
             m1 = multivariate_normal(mu1, sigma + c * np.identity(sigma.shape[0]))
             m2 = multivariate_normal(mu2, sigma + c * np.identity(sigma.shape[0]))
             return m1, m2
         def qda_train(spam, ham):
             mul, sigmal = gaussian mean cov(spam)
             mu2, sigma2 = gaussian mean cov(ham)
             m1 = multivariate normal(mu1, sigma1 + c * np.identity(sigma1.shape[0]))
             m2 = multivariate normal(mu2, sigma2 + c * np.identity(sigma2.shape[0]))
             return m1, m2
         def test(x, spam m, ham m):
             y = list()
             for sample in x:
                 spam y = spam m.logpdf(sample) + spam prior
                 ham y = ham m.logpdf(sample) + ham prior
                 y.append(np.argmax([ham y, spam y]))
             return y
         def evaluate(spam data, ham data, spam m, ham m):
             spam_y = test(spam_data, spam_m, ham_m)
             ham y = test(ham data, spam m, ham m)
             correct = np.count_nonzero(spam_y) + (len(ham_y) - np.count_nonzero(ham_
             total = len(spam y) + len(ham y)
             return 1.0 * correct / total
         def batch train and evaluate(spam data, ham data, spam validate, ham validat
             errors = []
             for size in categories:
                 spam_m, ham_m = qda_train(spam_data[:size, :], ham_data[:size, :])
                 if type == "LDA":
                     all = np.concatenate((spam_data, ham_data), axis=0)
                     spam m, ham m = lda train(spam data[:size, :], ham data[:size,
                 err = 1 - evaluate(spam validate, ham validate, spam m, ham m)
                 errors.append(err)
                 print("Error for training size {}: {}".format(size, err))
             plt.plot(categories, errors, 'ro')
             plt.axis([min(categories)-10, max(categories)+10, -0.1, max(errors)+0.1]
             plt.title("{} Error rate".format(type))
             return spam m, ham m
```



In [17]: qda_models = batch_train_and_evaluate(spam_data, ham_data, spam_validate, spa

```
Error for training size 100: 0.13479200432198812
Error for training size 200: 0.11561318206374938
Error for training size 500: 0.11642355483522415
Error for training size 1000: 0.09751485683414374
Error for training size 2000: 0.09292274446245274
```



```
In [24]: # Test
lda_y = test(test_data, lda_models[0], lda_models[1])
qda_y = test(test_data, qda_models[0], qda_models[1])
```

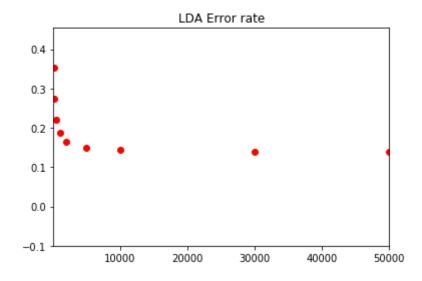
```
In [25]: df = pd.DataFrame(data = lda_y, columns=["Category"])
    df.index.name = "Id"
    df.to_csv("./spam.csv")
```

Kaggle: 0.95520

Feature Description:

Used bag of words approach plus normalization. More specifically, the features are the frequencies of each unique word in the text; after normalization, they become probabilities.

```
Error for training size 100: 0.35419999999999996
Error for training size 200: 0.2751
Error for training size 500: 0.22060000000000002
Error for training size 1000: 0.1877999999999997
Error for training size 2000: 0.1655999999999997
Error for training size 5000: 0.14970000000000006
Error for training size 10000: 0.14370000000000005
Error for training size 30000: 0.14
Error for training size 50000: 0.1392
```



```
Error for training size 200: 0.1490000000000002

Error for training size 500: 0.0950999999999996

Error for training size 1000: 0.0777999999999998

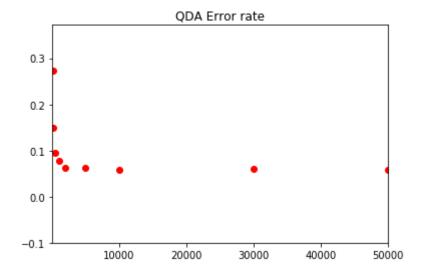
Error for training size 2000: 0.0634000000000001

Error for training size 5000: 0.06200000000000055

Error for training size 10000: 0.0590000000000005

Error for training size 30000: 0.05930000000000002

Error for training size 50000: 0.05810000000000000
```



```
In [9]: # Test
    lda_y = test(test_data, lda_models)
    qda_y = test(test_data, qda_models)
```

c) LDA vs QDA

QDA is better. Because we have 10 classes in this problem, but LDA uses the same variance for all their distributions, which isn't reasonable because different class random variable might have very different variance.

```
In [21]: df = pd.DataFrame(data = qda_y, columns=["Category"])
    df.index.name = "Id"
    df.to_csv("./mnist.csv")
```

Kaggle: 0.95640