Lecture 9 - Conjugate Priors & The Naive Bayes Classifier

Review of Gaussian Distribution

(Read section 2.3 "The Gaussian Distribution" from the Bishop textbook)

Univariate Gaussian

The Gaussian distribution is a widely used probabilistic model for the probability density function (pdf) of continuous random variables.

The Gaussian distribution can model both univariate (1-D) or multivariate (multi-dimensional) samples.

In the **univariate** case, the pdf of a Gaussian distribution for a random variable $X \in \mathbb{R}$ can be written as

$$f_X(x) = rac{1}{\sqrt{2\pi\sigma^2}} \mathrm{exp}igg(-rac{(x-\mu)^2}{2\sigma^2}igg)$$

In this case, we say that X follows a Gaussian distribution with mean μ and variance σ^2 , or, $X \sim N(\mu, \sigma^2)$.

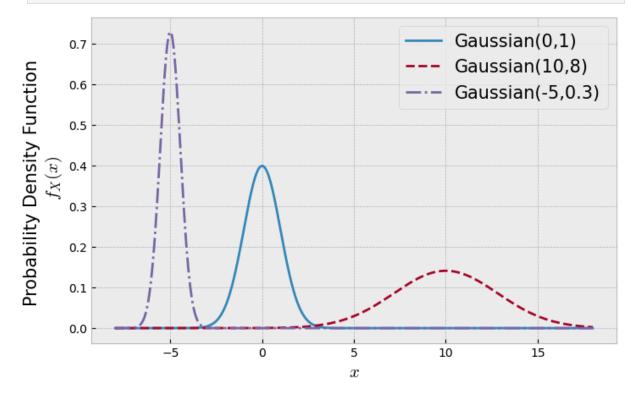
- We can define the **precision** parameter β as the inverse of the variance, that is, $\beta=\frac{1}{\sigma^2}.$
- A Gaussian distribution is called **Normal** when the mean is $\mu=0$ and variance is $\sigma^2=1$, $X\sim N(0,1)$.

```
import numpy as np
import matplotlib.pyplot as plt
%matplotlib inline
plt.style.use('bmh')
import scipy.stats as stats
```

```
In [2]: G1=stats.norm(0,1) # G(mean=0, variance=1^2)
G2=stats.norm(10,np.sqrt(8)) #G(mean=10, variance=8)
G3=stats.norm(-5,np.sqrt(0.3)) #G(mean=-5, variance=0.3^2)
```

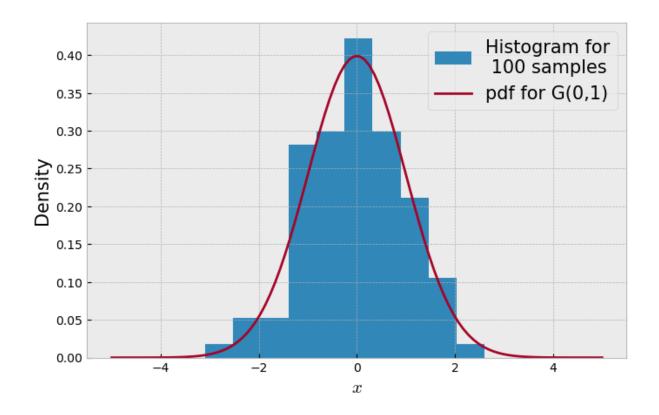
```
In [3]: plt.figure(figsize=(8,5))

x=np.linspace(-8,18,1000)
plt.plot(x,G1.pdf(x),label='Gaussian(0,1)')
plt.plot(x,G2.pdf(x),'--',label='Gaussian(10,8)')
plt.plot(x,G3.pdf(x),'--',label='Gaussian(-5,0.3)')
plt.legend(fontsize=15)
plt.xlabel('$x$',size=15)
plt.ylabel('Probability Density Function \n $f_X(x)$',size=15);
```



```
In [4]: samples=G1.rvs(size=100)

plt.figure(figsize=(8,5))
x=np.linspace(-5,5,100)
plt.hist(samples,density=True, label='Histogram for\n 100 samples')
plt.plot(x, G1.pdf(x), label='pdf for G(0,1)')
plt.legend(fontsize=15)
plt.xlabel('$x$',size=15)
plt.ylabel('Density',size=15);
```



Multivariate Gaussian

In the **multivariate** case, the pdf of a Gaussian distribution for a random variable $X \in \mathbb{R}^D$ can be written as

$$f_X(x) = rac{1}{\sqrt{(2\pi)^d \left|\Sigma
ight|}} \mathrm{exp}igg(-rac{1}{2}(\mathbf{x}-\mu)^T \Sigma^{-1} \left(\mathbf{x}-\mu
ight)igg)$$

In this case, we say that X follows a Gaussian distribution with mean μ and covariance Σ , or, $X \sim N(\mu, \Sigma)$.

- μ is a D-dimensional mean vector
- Σ is a D imes D covariance matrix
- $|\Sigma|$ denotes the determinant of Σ
- The precision parameter in a d-dimensional space is equal to $\beta=\Sigma^{-1}$

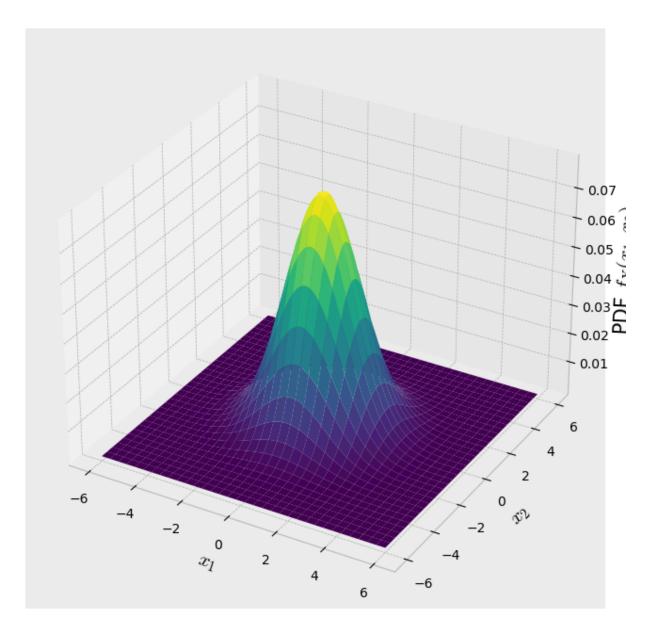
Let $X=[X_1,X_2]$. The **covariance** Σ measures the amount of variance is each individual dimension, X_1 and X_2 , as well as the amount of covariance between the two. We can write the covariance as

$$egin{aligned} ext{cov}(X_1, X_2) &= Eig[(X_1 - E\,[X_1])\,(X_2 - E\,[X_2])ig] \ &= egin{bmatrix} ext{var}(X_1) & ext{cov}(X_1, X_2) \ ext{cov}(X_1, X_2) & ext{var}(X_2) \end{bmatrix} \ &= egin{bmatrix} \sigma_{X_1}^2 & ext{cov}(X_1, X_2) \ ext{cov}(X_1, X_2) & \sigma_{X_2}^2 \end{bmatrix} \end{aligned}$$

• The **Pearson's correlation coefficient** between random variables X_1 and X_2 is defined as:

$$r=rac{ ext{cov}(X_1,X_2)}{\sqrt{ ext{var}(X_1)}\sqrt{ ext{var}(X_2)}}=rac{ ext{cov}(X_1,X_2)}{\sigma_{X_1}\sigma_{X_2}}$$

```
In [5]: mu = np.array([0,0]) # mean vector
        cov = np.array([[2,0],[0,2]]) # covariance matrix
        print(cov)
        G = stats.multivariate normal(mu, cov)
       [[2 0]
        [0 2]]
Out[5]: <scipy.stats._multivariate.multivariate_normal_frozen at 0x12e7afb10>
In [6]: x, y = np.mgrid[-6:6:100j, -6:6:100j]
        xy = np.column_stack([x.flat, y.flat])
        z = stats.multivariate_normal.pdf(xy, mean=mu, cov=cov)
        z = z.reshape(x.shape)
        fig = plt.figure(figsize=(8,8))
        ax = fig.add_subplot(111, projection='3d')
        ax.plot_surface(x,y,z, rstride=3, cstride=3, linewidth=1, antialiased=True,
                        cmap=plt.cm.viridis)
        ax.set_xlabel('$x_1$',size=15)
        ax.set_ylabel('$x_2$',size=15)
        ax.set_zlabel('PDF $f_X(x_1,x_2)$',size=15);
```



Let's use the Mathematica's demonstration "Joint Density of Bivariate Gaussian Random Variables" to better understand the role of the covariance matrix.

Example: Multivariate Gaussian-Gaussian Conjugate Prior Relationship

For a D-dimensional Gaussian data likelihood with mean μ and covariance βI and a prior distribution with mean μ_0 and covariance Σ_0

$$P(\mathbf{t}|\mathbf{w}) \sim \mathcal{N}(\mu, \beta \mathbf{I})$$

 $P(\mathbf{w}) \sim \mathcal{N}(\mu_0, \Sigma_0)$

The posterior distribution

$$egin{aligned} P(\mathbf{w}|\mathbf{t}) &\sim \mathcal{N}\left(\mu_N, \Sigma_N
ight) \ \mu_N &= \Sigma_N \left(\Sigma_0^{-1} \mu_0 + eta \mathbf{X}^T \mathbf{t}
ight) \ \Sigma_N^{-1} &= \Sigma_0^{-1} + eta \mathbf{X}^T \mathbf{X} \end{aligned}$$

where \mathbf{X} is the feature matrix of size $N \times M$.

• What happens with different values of β and Σ_0 ?

To simplify, let's assume the covariance of the prior to be **isotropic**, that is, it is a diagonal matrix with the same value along the diagonal, $\Sigma_0=\alpha^{-1}\mathbf{I}$. And, let $\mu_0=[0,0]$, thus

$$\mu_N = eta \Sigma_N \mathbf{X}^T \mathbf{t}$$

and

$$\Sigma_N^{-1} = \alpha \mathbf{I} + \beta \mathbf{X}^T \mathbf{X}$$

Let's consider the example presented in the Bishop textbook (Figure 3.7 in page 155).

Consider a single input variable \mathbf{x} , a single target variable \mathbf{t} and a linear model of the form $y(\mathbf{x}, \mathbf{w}) = w_0 + w_1 \mathbf{x}$. Because this has just two parameters coefficients, $w = [w_0, w_1]^T$, we can plot the prior and posterior distributions directly in parameter space (2-dimensional parameter space).

Let's generate some synthetic data from the function $f(x,a)=w_0+w_1x$ with parameter values $w_0=-0.3$ and $w_1=0.5$ by first choosing values of x_n from the uniform distribution $U(x_n|-1,1)$, then evaluating $f(x_n,\mathbf{w})$, and finally adding Gaussian noise with standard deviation of $\sigma=0.2$ to obtain the target values t_n .

$$t_n = f(x_n, \mathbf{w}) + \epsilon = -0.3 + 0.5x_n + \epsilon$$

where $\epsilon \sim \mathcal{N}(0, \beta \mathbf{I})$.

• Our goal is to recover the values of w_0 and w_1 from such data, and we will explore the dependence on the size of the data set.

For some data, $\{x_n,t_n\}_{n=1}^N$, we can pose this problem in terms of **Regularized Least Squares**:

$$egin{align} J(\mathbf{w}) &= rac{1}{2} \sum_{n=1}^{N} \left(t_n - y_n
ight)^2 + rac{\lambda}{2} \sum_{i=0}^{1} w_i^2 \ &= rac{1}{N} \sum_{n=1}^{2} \left(t_n - y_n
ight)^2 + rac{\lambda}{2} \left(w_0^2 + w_1^2
ight) \ &\Rightarrow rg_{\mathbf{w}} \min J(\mathbf{w}) \ \end{aligned}$$

Using MAP, we can rewrite our objective using the Bayesian interpretation:

$$\arg_{\mathbf{w}} \max P(\epsilon|\mathbf{w}) P(\mathbf{w})$$

Let's consider the data likelihood, $P(\epsilon|\mathbf{w})$, to be a Gaussian distribution with mean $\mu=0$ and variance $\sigma^2=\beta {\bf I}$. And let's also consider the prior distribution, $P({\bf w})$, to be a Gaussian distribution with mean μ_0 and variance $\sigma_0^2=lpha^{-1}{f I}$. Then, using the derivations from above, we can rewrite our optimization as:

$$rg_{\mathbf{w}} \max \mathcal{N}(\epsilon|0, eta \mathbf{I}) \mathcal{N}(\mathbf{w}|\mu_{\mathbf{0}}, lpha^{-1} \mathbf{I}) \ \propto rg_{\mathbf{w}} \max \mathcal{N}\left(eta \Sigma_{N}^{-1} \mathbf{X}^{T} \mathbf{t}, \Sigma_{N}\right)$$

where $\mu_0=[0,0]$, ${f X}$ is the polynomial feature matrix, and $\Sigma_N=\left(lpha^{-1}{f I}+eta{f X}^T{f X}
ight)^{-1}$ is the covariance matrix of the posterior distribution.

Note that we **do not known** the parameters of the prior distribution (μ_0 and σ_0 are unknown). The parameters of the prior distribution will have to be chosen by the user. And they will essentially encode any behavior or a priori knowledge we may have about the weights.

 Both our data likelihood and prior distributions are in a 2-dimensional space (this is because our *model order* is M=2 -- we have 2 parameters!).

We are going to generate data from $t=-0.3+0.5x+\epsilon$ where ϵ is drawn from a zeromean Gaussin distribution.

- The goal is to estimate the values $w_0=-0.3$ and $w_1=0.5$
- The feature matrix X can be computed using the polynomial basis functions
- Parameters to choose: β and α

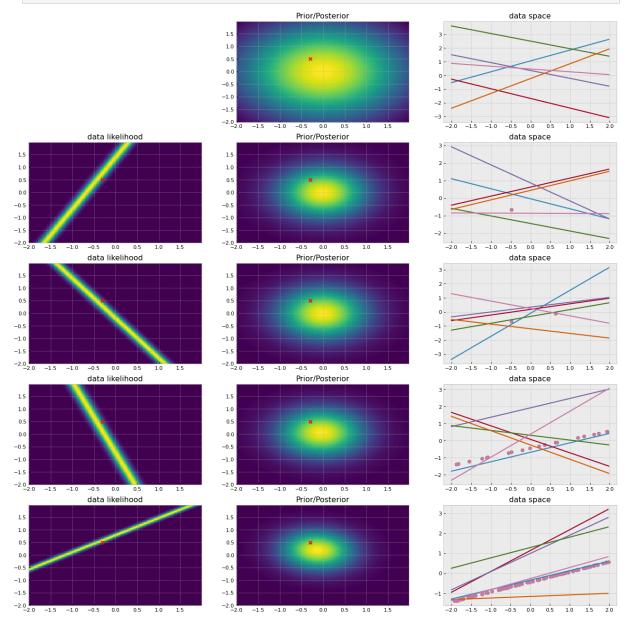
We want to implement this scenario for a case that we are getting more data every minute. As we get more and more data, we want to update our prior distribution using our posterior distribution (informative prior), because they take the have the same distribution form. This is only possible because because Gaussian-Gaussian have a conjugate prior relationship. That is, the posterior distribution is also a Gaussian.

```
import warnings
warnings.filterwarnings('ignore')
%matplotlib inline
# Fixing the random seed generator
np.random.seed(0)
def likelihood_prior_func(beta = 2, alpha = 1, draw_num=(0,1,10,20,50,100)):
    '''Online Update of the Posterior distribution for a Gaussian-Gaussian of
   Parameter:
   beta - variance of the data likelihood (of the additive noise)
   alpha - precision value or 1/variance of the prior distribution
   draw_num - number of points collected at each instance.
   This function will update the prior distribution as new data points are
   The prior distribution will be the posterior distribution from the last
   fig = plt.figure(figsize=(20, 20), dpi= 80, facecolor='w', edgecolor='k'
   # true (unknown) weights
   a = -0.3 \# w0
   b = 0.5 \# w1
   # set up input space
   rangeX = [-2, 2] # range of values for the input
   step = 0.025 # distance between points
   X = np.mgrid[rangeX[0]:rangeX[1]:step] # creates a grid of values for in
   #initialize prior/posterior and sample data
   S0 = (1/alpha)*np.eye(2) # prior covariance matrix
   sigma = S0 # copying it so we can update it later
   mean = [0,0] # mean for prior
   # Draws samples from Uniform(-1,1) distribution
   draws = np.random.uniform(rangeX[0],rangeX[1],size=draw_num[-1])
   # Generate the noisy target samples
   T = a + b*draws + np.random.normal(loc=0, scale=np.sqrt(beta))
   for i in range(len(draw_num)):
        if draw_num[i]>0: #skip first image
            # INPUT DATA
            #Feature Matrix (Polynomial features with M=2)
            FeatureMatrix = np.array([draws[:draw num[i]]**m for m in range(
            #Target Values
            t = T[0:draw_num[i]]
            # POSTERIOR PROBABILITY
            # Covariance matrix
            sigma = np.linalq.inv(S0 + beta*FeatureMatrix.T@FeatureMatrix)
            # Mean vector
           mean = beta*sigma@FeatureMatrix.T@t
            # PARAMETER SPACE
            # create a meshgrid of possible values for w's
            w0, w1 = np.mgrid[rangeX[0]:rangeX[1]:step, rangeX[0]:rangeX[1]:
```

```
# Define the Gaussian distribution for data likelihood
    p = multivariate normal(mean=t[draw num[i]-1], cov=beta)
    # Initialize the PDF for data likelihood
    out = np.empty(w0.shape)
    # For each value (w0,w1), compute the PDF for all data samples
    for j in range(len(w0)):
        out[j] = p.pdf(w0[j]+w1[j]*draws[draw_num[i]-1])
    # Plot the data likelihood
    ax = fig.add_subplot(*[len(draw_num),3,(i)*3+1])
    ax.pcolor(w0, w1, out)
    # Add the current value for parameters w=(w0,w1)
    ax.scatter(a,b, c='r',marker='x')
    myTitle = 'data likelihood'
    ax.set_title("\n".join(textwrap.wrap(myTitle, 100)))
# PARAMETER SPACE
# create a meshgrid of possible values for w's
w0, w1 = np.mgrid[rangeX[0]:rangeX[1]:step, rangeX[0]:rangeX[1]:step
# POSTERIOR PROBABILITY
# initialize the matrix with posterior PDF values
pos = np.empty(w1.shape + (2,))
# for w0
pos[:, :, 0] = w0
# and for w1
pos[:, :, 1] = w1
# compute the PDF
p = multivariate_normal(mean=mean, cov=sigma)
#Show prior/posterior
ax = fig.add_subplot(*[len(draw_num),3,(i)*3+2])
ax.pcolor(w0, w1, p.pdf(pos))
# Add the value for parameters w=(w0,w1) that MAXIMIZE THE POSTERIOF
ax.scatter(a,b, c='r',marker='x')
myTitle = 'Prior/Posterior'
ax.set_title("\n".join(textwrap.wrap(myTitle, 100)))
# DATA SPACE
ax = fig.add_subplot(*[len(draw_num),3,(i)*3+3])
for j in range(6):
    # draw sample from the prior probability to generate possible var{\epsilon}
    w0, w1 = np.random.multivariate normal(mean=mean, cov=sigma)
    # Estimated labels
    t = w0 + w1*X
    # Show data space
    ax.plot(X,t)
    if draw num[i] > 0:
        ax.scatter(FeatureMatrix[:,1], T[0:draw num[i]])
    myTitle = 'data space'
    ax.set_title("\n".join(textwrap.wrap(myTitle, 100)))
```

```
In [21]: # Parameters
# beta - variance of the data likelihood (for the additive noise)
# alpha - precision value or 1/variance for the prior distribution
```

draw_num - number of points collected at each instance
likelihood_prior_func(beta = 0.01, alpha = 0.5, draw_num=(0,1,2,20,100))



Introduction to Classification & Probabilistic Generative Models

So far we have focused on regression. We will begin to discuss classification.

Suppose we have training data from two classes, C_1 and C_2 , and we would like to train a classifier to assign a label to incoming test points as C_1 or C_2 .

There are *many* classifiers in the machine learning literature. We will cover a few in this course. Today we will focus on probabilistic generative approaches for classification.

There are two types of classification algorithms: **discriminative** or **generative**.

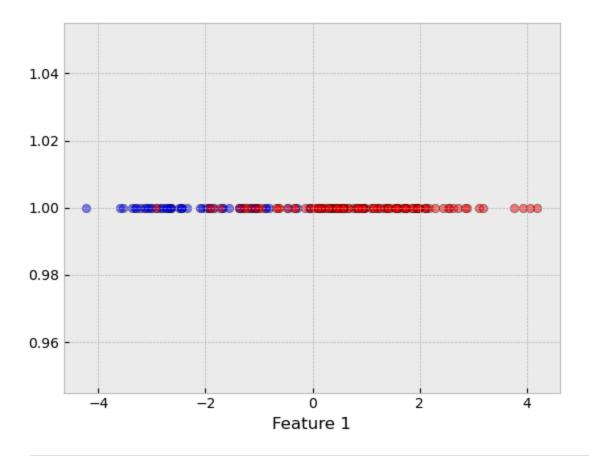
- A discriminative approach for classification is one in which we partition the feature space into regions for each class. Then, when we have a test point, we evaluate in which region it landed on and classify it accordingly.
- A **generative** approach for classification is one in which we estimate the parameters for distributions that generate the data for each class. Then, when we have a test point, we can compute the posterior probability of that point belonging to each class and assign the point to the class with the highest posterior probability.

```
In [9]: def generateData(mean1, mean2, cov1, cov2, N1, N2):
            # We are generating data from two Gaussians to represent two classes
            # In practice, we would not do this — we would just have data from the p
            data_C1 = stats.multivariate_normal(mean1, cov1).rvs(size=N1)
            data_C2 = stats.multivariate_normal(mean2, cov2).rvs(size=N2)
            # Entire Training Dataset
            data = np.concatenate((data C1, data C2))
            labels = np.concatenate((np.ones(N1),2*np.ones(N2)))
            if len(mean1)>1:
                plt.scatter(data[labels==1,0], data[labels==1,1], c='b', alpha=0.5,
                plt.scatter(data[labels==2,0], data[labels==2,1], c='r', alpha=0.5,
                plt.xlabel('Feature 1'); plt.ylabel('Feature 2');
            else:
                plt.scatter(data[labels==1], np.ones(N1), c='b', alpha=0.5, edgecold
                plt.scatter(data[labels==2], np.ones(N2), c='r', alpha=0.5, edgecold
                plt.xlabel('Feature 1');
            return data, labels
```

Case 1: Univariate Data Likelihood

```
In [10]: mean1 = [-2]
    mean2 = [1]
    var1 = [1]
    var2 = [2]
    N1 = 50 # C1 - blue
    N2 = 100 # C2 - red

data, labels = generateData(mean1, mean2, var1, var2, N1, N2)
```



```
In [11]: def drawMAP(data, labels):
             #### Estimate parameters (MLE solution)
             # Means
             mu1 = np.mean(data[labels==1])
             mu2 = np.mean(data[labels==2])
             # Variances
             var1 = np.cov(data[labels==1])
             var2 = np.cov(data[labels==2])
             #### Estimate Prior Probabilities - relative frequency
             N = len(data)
             N1 = np.sum(labels==1)
             N2 = N - N1
             p1 = N1/N # prior probability for C1
             p2 = N2/N # prior probability for C2
             #### Define data likelihoods
             G1=stats.norm(loc=mu1, scale=np.sqrt(var1)) # P(x|C1)
             G2=stats.norm(loc=mu2, scale=np.sqrt(var2)) # P(x|C2)
             x=np.linspace(-6,6,1001)
             #### Plot the weighted densities
             # these are proportional to the posteriors
             plt.figure(figsize=(10,5))
             plt.plot(x,p1*G1.pdf(x),label='$f_X(x|C_1)P(C_1)$')
             plt.plot(x,p2*G2.pdf(x),label='$f_X(x|C_2)P(C_2)$')
             #### Determine the regions where the posterior for deciding C1
```

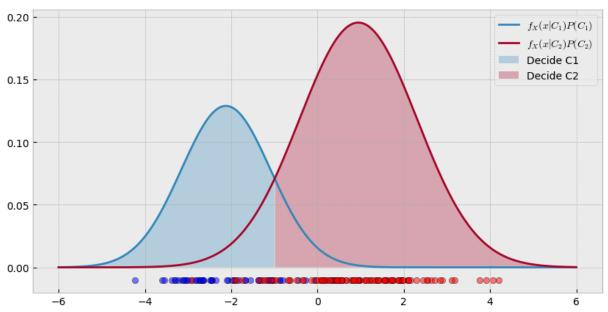
```
# and the posterior for deciding C2
R1=x[np.where(p1*G1.pdf(x)>= p2*G2.pdf(x))]
R2=x[np.where(p1*G1.pdf(x)< p2*G2.pdf(x))]

# Fill under the regions found above
plt.fill_between(R1,p1*G1.pdf(R1),alpha=0.3,label='Decide C1')
plt.fill_between(R2,p2*G2.pdf(R2),alpha=0.3,label='Decide C2')
plt.scatter(data[labels==1], -0.01*np.ones(N1), c='b', alpha=0.5, edgecc plt.scatter(data[labels==2], -0.01*np.ones(N2), c='r', alpha=0.5, edgecc plt.legend()

# Print the MAP threshold
print('MAP decision threshold to decide C2 is >',round(R2[0],2))
```

In [12]: drawMAP(data, labels)

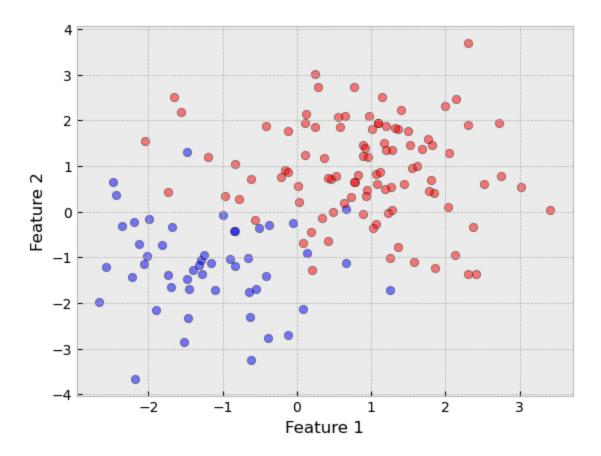
MAP decision threshold to decide C2 is > -0.98



Case 2: Multivariate Data Likelihood

```
In [13]: mean1 = [-1, -1]
    mean2 = [1, 1]
    cov1 = [[1,0],[0,1]]
    cov2 = [[1,0],[0,1]]
    N1 = 50
    N2 = 100

data, labels = generateData(mean1, mean2, cov1, cov2, N1, N2)
```



In the data we generated above, we have a "red" class and a "blue" class. When we are given a test sample, we will want to assign the label of red or blue.

We can compute the **posterior probability** for class C_1 as follows:

$$P(C_1|x)=rac{P(x|C_1)P(C_1)}{P(x)}$$

Understanding that the two classes, red and blue, form a partition of all possible classes, then we can utilize the *Law of Total Probability*, and obtain:

$$P(C_1|x) = rac{P(x|C_1)P(C_1)}{P(x|C_1)P(C_1) + P(x|C_2)P(C_2)}$$

Similarly, we can compute the posterior probability for class C_2 :

$$P(C_2|x) = rac{P(x|C_2)P(C_2)}{P(x|C_1)P(C_1) + P(x|C_2)P(C_2)}$$

Note that $P(C_1|x) + P(C_2|x) = 1$.

The Naive Bayes Classifier

Therefore, for a given test point \mathbf{x}^* , our decision rule is:

$$P(C_1|\mathbf{x}^*) \mathop{\gtrless}\limits_{C_2}^{C_1} P(C_2|\mathbf{x}^*)$$

Using the Bayes' rule, we can further rewrite it as:

$$\frac{P(\mathbf{x}^*|C_1)P(C_1)}{P(\mathbf{x}^*)} \underset{C_2}{\overset{C_1}{\geqslant}} \frac{P(\mathbf{x}^*|C_2)P(C_2)}{P(\mathbf{x}^*)}$$

$$\iff P(\mathbf{x}^*|C_1)P(C_1) \underset{C_2}{\overset{C_1}{\geqslant}} P(\mathbf{x}^*|C_2)P(C_2)$$

We assign \mathbf{x}^* as class 1 if $p(\mathbf{x}^*|C_1)p(C_1) > p(\mathbf{x}^*|C_2)p(C_2)$, or assign \mathbf{x}^* to class 2 if $p(\mathbf{x}^*|C_1)p(C_1) < p(\mathbf{x}^*|C_2)p(C_2)$.

This defines the Naive Bayes Classifier.

Training a Generative Classifier

• So, to train the classifier, what we need to do is to determine the parametric forms and the associated parameters for $P(x|C_1)$, $P(x|C_2)$, $P(C_1)$ and $P(C_2)$.

For example, we can assume that the data samples coming from either ${\cal C}_1$ and ${\cal C}_2$ are distributed according to Gaussian distributions. In this case,

$$P(x|C_k) = rac{1}{(2\pi)^{1/2} |\Sigma_k|^{1/2}} \mathrm{exp}igg\{ -rac{1}{2} (\mathbf{x} - \mu_k)^T \Sigma_k^{-1} (\mathbf{x} - \mu_k) igg\}, orall k = \{1, 2\}$$

• We can consider any distributional form we want.

What about the $P(C_1)$ and $P(C_2)$?

• We can consider the relative frequency of each class, that is, $P(C_i)=\frac{N_i}{N}$, where N_i is the number of points in class C_i and N is the total number of samples.

MLE Parameter Estimation Steps

For simplification, let's consider the covariance matrix Σ_k for k=1,2 to be **isotropic** matrices, that is, the covariance matrix is diagonal and the element along the diagonal is the same, or: $\Sigma_k = \sigma_k^2 \mathbf{I}$.

 What are the parameters? The mean and covariance of the Gaussian distribution for both classes. Given the assumption of the Gaussian form, how would you estimate the parameters for $p(x|C_1)$ and $p(x|C_2)$? We can use **maximum likelihood estimate** for the mean and covariance, because we are looking for the parameters of the distributions that *maximize* the data likelihood!

Assumption: Assuming the classes follow a (bivariate or 2-D) Gaussian distribution and, for simplicity, let's assume the covariance matrices are **isotropic**, that is, $\Sigma_k = \sigma_k^2 \mathbf{I}$.

The MLE steps for parameter estimation are:

1. Write down the observed data likelihood, \mathcal{L}^0

$$egin{aligned} \mathcal{L}^0 &= P(x_1, x_2, \dots, x_N | C_k) \ &= \prod_{n=1}^N P(x_n | C_k), ext{ data samples are i.i.d.} \ &= \prod_{n=1}^N rac{1}{(2\pi)^{d/2} |\Sigma|^{1/2}} \mathrm{exp} iggl\{ -rac{1}{2} (x_n - \mu_k)^T \Sigma_k^{-1} (x_n - \mu_k) iggr\} \ &= \prod_{n=1}^N rac{1}{(2\pi)^{d/2} |\sigma_k^2 \mathbf{I}|^{1/2}} \mathrm{exp} iggl\{ -rac{1}{2\sigma_k^2} (x_n - \mu_k)^T \mathbf{I} (x_n - \mu_k) iggr\} \ &= \prod_{n=1}^N rac{1}{(2\pi)^{d/2} (\sigma_k^2)^{d/2}} \mathrm{exp} iggl\{ -rac{1}{2\sigma_k^2} (x_n - \mu_k)^T (x_n - \mu_k) iggr\} \end{aligned}$$

where d is the dimensionality of the data space.

2. Take the log-likelihood, \mathbf{L} . This *trick* helps in taking derivatives.

$$egin{aligned} \mathcal{L} &= \ln\!\left(\mathcal{L}^0
ight) \ &= \sum_{n=1}^N \left(-rac{d}{2}\!\ln 2\pi - rac{d}{2}\!\ln \sigma_k^2 - rac{1}{2\sigma_k^2}(x_n - \mu_k)^T(x_n - \mu_k)
ight) \end{aligned}$$

3. Take the derivative of the log-likelihood function with respect to the parameters of interest. For Gaussian distribution they are the mean and covariance.

$$egin{aligned} rac{\partial \mathcal{L}}{\partial \mu_k} &= 0 \ \sum_{n \in C_k} rac{1}{\sigma_k^2} (x_n - \mu_k) &= 0 \ \sum_{n \in C_k} (x_n - \mu_k) &= 0 \ \sum_{n \in C_k} x_n - \sum_{n \in C_k} \mu_k &= 0 \ \sum_{n \in C_k} x_n - N_k \mu_k &= 0 \end{aligned}$$

This is the sample mean for each class. And,

$$egin{aligned} rac{\partial \mathcal{L}}{\partial \sigma_k^2} &= 0 \ \sum_{n \in C_k} -rac{d}{2\sigma_k^2} + rac{2(x_n - \mu_k)^T(x_n - \mu_k)}{(2\sigma_k^2)^2} &= 0 \ \sum_{n \in C_k} -d + rac{(x_n - \mu_k)^T(x_n - \mu_k)}{\sigma_k^2} &= 0 \ rac{\sum_{n \in C_k} (x_n - \mu_k)^T(x_n - \mu_k)}{\sigma_k^2} &= dN_k \ \sigma_k^2 &= rac{\sum_{n \in C_k} (x_n - \mu_k)^T(x_n - \mu_k)}{dN_k} \end{aligned}$$

This is the sample variance for each class. Then we can create $\Sigma_k = \sigma_k^2 \mathbf{I}$, which is the (biased) sample covariance for each class.

In practice, if we want to estimate an entire covariance matrix, we would have to take the derivative of the log-likelihood function with respect to every entry in the covariance matrix. Covariance matrices are symmetric, so we only need to determine the upper (or lower) half of the covariance matrix.

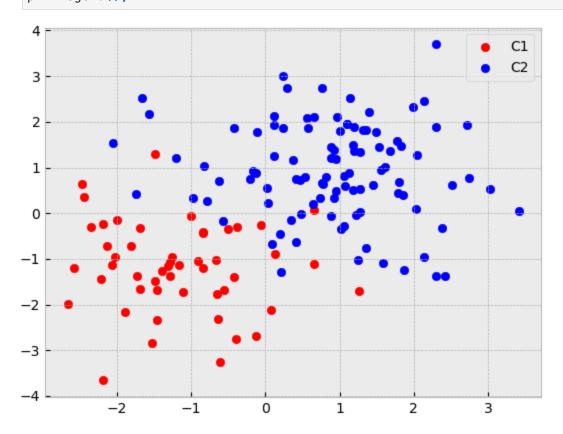
We can determine the values for $p(C_1)$ and $p(C_2)$ from the number of data points in each class:

$$p(C_k) = rac{N_k}{N}$$

where N is the total number of data points.

```
In [14]: data.shape
Out[14]: (150, 2)

In [15]: plt.scatter(data[labels==1,0], data[labels==1,1], color='red', label='C1')
    plt.scatter(data[labels==2,0], data[labels==2,1], color='blue', label='C2')
    plt.legend();
```



```
In [16]: #### Estimate parameters (with MLE solutions)
         # Means
         mu1 = np.mean(data[labels==1], axis=0)
         print('Mean of Class 1: ', mu1)
         mu2 = np.mean(data[labels==2], axis=0)
         print('Mean of Class 2: ', mu2)
         # Covariances — in this example we are showing the case where we estimate th
         cov1 = np.cov(data[labels==1,:].T) # np.cov expects input to be D-by-N
         print('Covariance of Class 1: ',cov1)
         cov2 = np.cov(data[labels==2, :].T)
         print('Covariance of Class 2: ',cov2)
         #### Estimate Prior Probabilities
         N = N1+N2
         p1 = N1/N
         print('Probability of Class 1: ',p1)
         p2 = N2/N
         print('Probability of Class 2: ',p2)
```

Next class.. we will define the data likelihood using these parameters and draw the decision boundary using the Naive Bayes Classifier.