# **Problem Set 5: Probabilistic Models**

This assignment requires a working IPython Notebook installation, which you should already have. If not, please refer to the instructions in the previous problem sets.

Only PDF files are accepted for submission. To print this notebook to a pdf file, you can go to "File" -> "Download as" -> "PDF via LaTex(.pdf)" or simply use "print" in browser.

Total: 100 points (+20 points bonus).

# 1 Linear Discriminant Analysis (LDA) [50 pts + 10pts bonus]

In this exercise, we build a linear discriminant analysis (LDA) classifier for the problem of predicting whether a student gets admitted into a university.

LDA is a generative model for classification. Given a training dataset of positive and negative features (x,y) with  $y \in 0,1$ , LDA models the data x as generated from class-conditional Gaussians: P(x,y) = P(x|y)P(y), where  $P(y=1) = \pi$  and  $P(x|y=j) = N(x;\mu^j,\Sigma)$ . Here means  $\mu^j$  are class-dependent but the covariance matrix  $\Sigma$  is class-independent (the same for all classes).

A novel feature x is classified as a positive if P(y=1|x)>P(y=0|x), which is equivalent to a(x)>0, where the linear classifier  $a(x)=w^Tx+w_0$  has weights given by  $w=\Sigma^{-1}(\mu^1-\mu^0)$ .

In practice, and in this assignment, we use a(x) > some threshold, or equivalently,  $w^T x > T$  for some constant T. The specific threshold can be determined on a validation dataset or via cross-validation.

As we saw in lecture, LDA and logistic regression can be expressed in the same form

$$P(y=1|x)=rac{1}{1+e^{- heta^Tx}}.$$

However, they generally produce different solutions for the parameter theta.

## 1.1 Derivation of LDA [20 pts]

Show that the log-odds decision function a(x) for LDA

$$a(x) = \ln rac{p(x|y=1)p(y=1)}{p(x|y=0)p(y=0)}$$

is linear in x, that is, we can express  $a(x) = w^T x + w_0$  for some  $w, w_0$ . Show all your steps.

Answer: 
$$P(x|y=j)=N(x;\mu^j,\Sigma)=rac{1}{2\pi^{rac{p}{2}}|\Sigma|^{rac{1}{2}}}e^{-rac{1}{2}(x-\mu_j)^T\Sigma^{-1}(x-\mu_j)}$$

$$p(y=1)=\pi$$
 and  $p(y=0)=1-\pi$ . Therefore,  $\ln rac{p(y=1)}{p(y=0)}=const$  Thus we have:

$$\begin{split} \ln \frac{p(x|y=1)p(y=1)}{p(x|y=0)p(y=0)} &= \ln(p(x|y=1)) - \ln(p(x|y=0)) + \ln \frac{p(y=1)}{p(y=0)} \\ &= \ln(\frac{1}{2\pi^{\frac{p}{2}}|\Sigma|^{\frac{1}{2}}}e^{-\frac{1}{2}(x-\mu_1)^T\Sigma^{-1}(x-\mu_1)}) - \ln(\frac{1}{2\pi^{\frac{p}{2}}|\Sigma|^{\frac{1}{2}}}e^{-\frac{1}{2}(x-\mu_0)^T\Sigma^{-1}(x-\mu_0)}) + const \\ &= -\frac{1}{2}(x-\mu_1)^T\Sigma^{-1}(x-\mu_1) + \frac{1}{2}(x-\mu_0)^T\Sigma^{-1}(x-\mu_0) + const \\ &= -\frac{1}{2}(x^T\Sigma^{-1}x - 2x^T\Sigma^{-1}\mu_1 + \mu_1^T\Sigma^{-1}\mu_1) + \frac{1}{2}(x^T\Sigma^{-1}x - 2x^T\Sigma^{-1}\mu_0 + \mu_0^T\Sigma^{-1}\mu_0) + const \end{split}$$

$$= -\frac{1}{2}(x^T \Sigma^{-1} x - 2x^T \Sigma^{-1} \mu_1 + \mu_1^T \Sigma^{-1} \mu_1 - x^T \Sigma^{-1} x + 2x^T \Sigma^{-1} \mu_0 - \mu_0^T \Sigma^{-1} \mu_0) + const$$

$$= -\frac{1}{2}(-2x^T \Sigma^{-1} (\mu_1 - \mu_0) + \mu_1^T \Sigma^{-1} \mu_1 - \mu_0^T \Sigma^{-1} \mu_0) + const$$

We have  $w=\Sigma^{-1}(\mu^1-\mu^0)$ 

Let 
$$-2(w_0) = \mu_1^T \Sigma^{-1} \mu_1 - \mu_0^T \Sigma^{-1} \mu_0$$

Thus:

$$=-rac{1}{2}(-2x^T\Sigma^{-1}(\mu_1-\mu_0)+\mu_1^T\Sigma^{-1}\mu_1-\mu_0^T\Sigma^{-1}\mu_0)=w^Tx+w_0=a(x)$$

# **Implementation**

In this part, you can assume the prior probabilities for the two classes are the same (although the number of the positive and negative samples in the training data is not the same), and that the threshold T is zero (you do not need to find  $w_0$ ). As a bonus, you are encouraged to explore how the different prior probabilities shift the decision boundary.

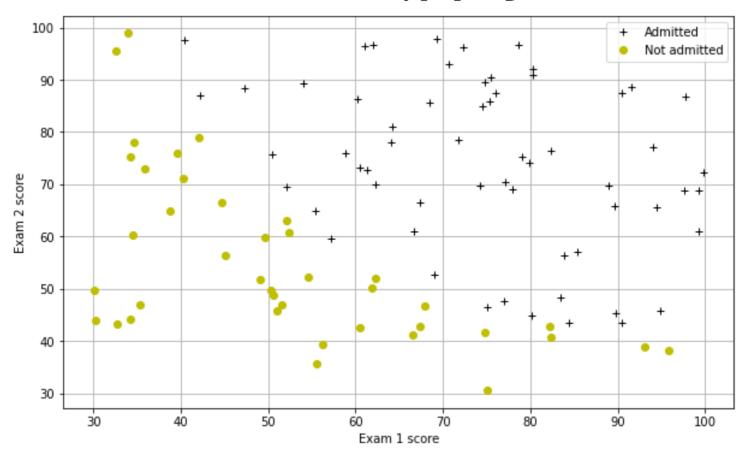
# 1.2 Loading the data

First, lets load and plot the data.

```
%matplotlib inline
import matplotlib.pyplot as plt
import numpy as np

datafile = 'data/ex2data1.txt'
#!head $datafile
```

```
cols = np.loadtxt(datafile, delimiter=',', usecols=(0, 1, 2), unpack=True) #Read in comma separat
##Form the usual "X" matrix and "v" vector
X = np.transpose(np.array(cols[:-1]))
v = np.transpose(np.arrav(cols[-1:]))
m = v.size # number of training examples
##Insert the usual column of 1's into the "X" matrix
X = np.insert(X, 0, 1, axis=1)
#Divide the sample into two: ones with positive classification, one with null classification
pos = np.array([X[i] for i in range(X.shape[0]) if v[i] == 1])
neg = np.array([X[i] for i in range(X.shape[0]) if y[i] == 0])
def plotData():
    plt.figure(figsize=(10, 6))
    plt.plot(pos[:, 1], pos[:, 2], 'k+', label='Admitted')
    plt.plot(neg[:, 1], neg[:, 2], 'vo', label='Not admitted')
    plt.xlabel('Exam 1 score')
    plt.ylabel('Exam 2 score')
    plt.legend()
    plt.grid(True)
plotData()
```



Implement the LDA classifier by completing the code below.

# 1.3 Centering the data [5 pts]

As an implementation detail, you should first center the positive and negative data separately, so that each set has a mean equal to [0, 0], before computing the covariance, as this tends to give a more accurate estimate of the covariance.

```
# IMPLEMENT THIS
pos_mean = np.mean(pos, axis = 0)
neg_mean = np.mean(neg, axis = 0)
```

```
#print(pos_mean)
#print(neg_mean)
pos_data = pos - (1/2)*(pos_mean+neg_mean)
neg_data = neg - (1/2)*(pos_mean+neg_mean)
```

## 1.4 Computing parameters and predictions [20 pts]

Implement the LDA algorithm here in vectorized form (avoid loops). First, compute the covariance matrix, then the classifier's weights, then use the classifier to make predictions on the training data. Note, you should center the whole training data set before applying the classifier. Namely, subtract the mean value of the two classes' means (1/2\*(pos\_mean+neg\_mean)), which is on the separating plane when their prior probabilities are the same and becomes the 'center' of the data

```
In [405...
# IMPLEMENT THIS
cov_all = np.cov(pos_data.T[1:]) + np.cov(neg_data.T[1:])# SHAPE: (2,2)
#print(cov_all.shape)
w = np.linalg.inv(cov_all)@((pos_mean - neg_mean)[1:])
#print(w)
X_centered = np.concatenate((neg_data , pos_data), axis = 0)
y_lda = np.dot(X_centered[:,1:],w) # SHAPE: (100,)
#print(y_lda.shape)
```

# 1.5 Training accuracy [5 pts]

Complete the code to compute the training set accuracy, which should be around 89%.

```
In [406...
# IMPLEMENT THIS
count = 0
y_sort = np.sort(y, axis = 0)

for id, i in enumerate(y_lda):
    if i < 0:</pre>
```

```
y_lda[id] = 0
if i > 0:
    y_lda[id] = 1
if y_lda[id] == y_sort[id]:
    count+=1

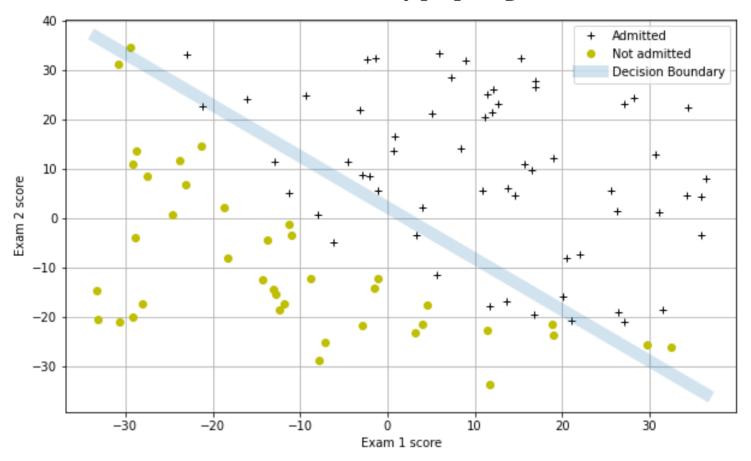
accuracy = count*100/m
print(accuracy)
```

89.0

## 1.6 Bonus [10 pts] (optional)

As an optional bonus exercise, try changing the prior probability of the positive class  $P(y=1)=\pi$  and show how the resulting decision boundary changes by plotting it. What is the effect of changing  $\pi$  on the boundary? The Boundary moves toward the other class respectively to alpha

```
In [407...
          # IMPLEMENT THIS
          #plt.plot(y lda)
          plt.figure(figsize=(10, 6))
          plt.plot(pos data[:, 1], pos data[:, 2], 'k+', label='Admitted')
          plt.plot(neg data[:, 1], neg data[:, 2], 'yo', label='Not admitted')
          plt.xlabel('Exam 1 score')
          plt.ylabel('Exam 2 score')
          plt.legend()
          plt.grid(True)
          linexs = np.array([np.max(X centered[:,1]),np.min(X centered[:,1])])
          alpha = 0.9 \# < 1
          lineys = np.log(alpha/(1-alpha))-linexs*w[0]/w[1]
          plt.plot(linexs, lineys, lw=10, alpha=0.2, label='Decision Boundary')
          plt.legend()
          plt.show()
```



# 2 Conjugate Prior [25 pts]

In this problem we will estimate the parameter  $\theta$  of the coin flipping problem using the Bayesian approach. In the Bayesian approach,  $\theta$  is considered a random variable sampled from some prior distribution  $p(\theta)$ . We represent the ith coin flip by a Bernoulli random variable  $x \in \{0,1\}$ , where  $x_i = 1$  if the ith flip lands on heads. We assume that the observed  $x_i$ 's are conditionally indendent given  $\theta$ . This means that the joint distribution of n coin flips and  $\theta$  can be factorized as

$$p(x_1,\dots,x_n, heta)=p( heta)\prod_{i=1}^n p(x_i| heta)$$

Suppose that our prior distribution on  $\theta$  is Beta(h,t), for some h,t>0. That is,

$$p(\theta) \propto \theta^{h-1} (1-\theta)^{t-1}$$

See Bishop Ch 2.1.1 or https://en.wikipedia.org/wiki/Beta\_distribution for more details on this distribution. There is a nice online graphing tool here http://eurekastatistics.com/beta-distribution-pdf-grapher/.

The Beta distribution is a conjugate to the Bernoulli distribution since the prior  $p(\theta)$  and posterior  $p(\theta|x_1,\ldots,x_n)$  are in the same family (the Beta family).

## 2.1 Posterior [10 pts]

Suppose our dataset of flips has H heads and T tails. Show that the posterior distribution fo  $\theta$  is Beta(h+H,t+T), i.e. show that

$$p(\theta|x_1,\ldots,x_n) \propto \theta^{h-1+H} (1-\theta)^{t-1+T}$$

#### **Answer:**

Based on the definition: Posterior = Prior x Likelihood

$$egin{aligned} p( heta|x_1,\ldots,x_n) &= p( heta)p(x_1,\ldots,x_n| heta) \ & p( heta|x_1,\ldots,x_n) &= Beta(h,t)*Binomial(n, heta) \ &= Beta(x_1+x_2+\ldots+x_n+h,n-x_1-x_2-\ldots-x_n+t) \end{aligned}$$

We have  $x_i=1$  when it is heads, and 0 when it is tails. Thus,  $x_1+x_2+\ldots+x_n=H$  and  $n-x_1-x_2-\ldots-x_n=T$ . Therefore, we have:

$$p( heta|x_1,\ldots,x_n) = Beta(h+H,t+T) = rac{\Gamma(h+H+t+T)}{\Gamma(h+H)\Gamma(t+T)} heta^{h-1+H}(1- heta)^{t-1+T}$$

$$p( heta|x_1,\dots,x_n) \propto heta^{h-1+H} (1- heta)^{t-1+T}$$

# 2.2 Parameter estimates [10 pts]

Give the expressions for three estimates of  $\theta$ : the Maximum Likelihood ( $\theta_{ML}$ ), the Maximum a Posteriori ( $\theta_{MAP}$ ) or the mode of the posterior over  $\theta$ , and the mean of the posterior ( $\theta_{MP}$ ).

Hint: You may use the fact that a Beta(h, t) distribution has mean h/(h+t) and has mode (h-1)/(h+t-2) for h,t>1.

**Answer:** We have maximum likelihood  $p(x_i|\theta) = \theta^H (1-\theta)^T$  and:

$$heta_{ML} = rac{H}{H+T}$$

We have posterior  $p(\theta|x_i) \propto \theta^{h-1+H}(1-\theta)^{t-1+T}$ , proven above. Thus, with the same way we solve for maximum likelihood,

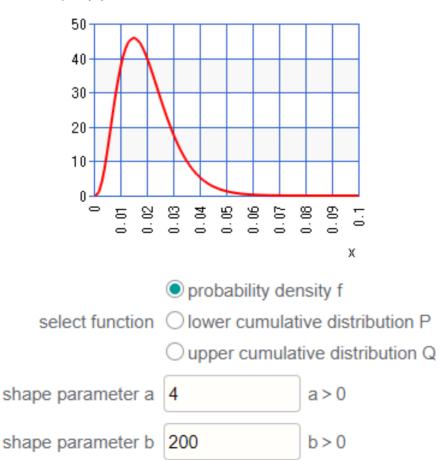
$$heta_{MAP}=rac{h-1+H}{h-1+H+t-1+T}=rac{H+h-1}{H+T+h+t-2}$$
  $heta_{MP}=rac{H+h}{H+T+h+t}$ 

## 2.3 Example [5 pts]

Supose you are modeling some event with a binary outcome (such as a user clicking on an ad after viewing it). You know that typical fractions of positive events observed in data are between 0.01 and 0.02. What are reasonable hyperparameters of the Beta prior to use, and why? Include a plot of the resulting prior PDF.

#### **Answer:**

Beta(4, 200) satisfy  $p(\theta)$  between 0.01 and 0.02



# 3 Bayesian Linear Regression [25 pts + 10pts bonus]

In this exercise, we implement Bayesian regression with linear basis function models. Recall that a linear regression model  $y(\mathbf{x}, \mathbf{w})$  can be defined more generally as

$$y(\mathbf{x},\mathbf{w}) = w_0 + \sum_{j=1}^{M-1} w_j \phi_j(\mathbf{x}) = \sum_{j=0}^{M-1} w_j \phi_j(\mathbf{x}) = \mathbf{w}^T \phi(\mathbf{x})$$

where  $\phi_j$  are basis functions and M is the total number of parameters  $w_j$  including the bias term  $w_0$ . The target variable t of an observation  $\mathbf{x}$  is given by a deterministic function  $y(\mathbf{x}, \mathbf{w})$  plus additive random noise  $\epsilon$ .

$$t = y(\mathbf{x}, \mathbf{w}) + \epsilon$$

We assume that the noise is normally distributed i.e. follows a Gaussian distribution with zero mean and precision (inverse variance)  $\beta$ . The corresponding probabilistic model i.e. the conditional distribution of t given  $\mathbf{x}$  can therefore be written as

$$p(t|\mathbf{x},\mathbf{w},eta) = \mathcal{N}(t|y(\mathbf{x},\mathbf{w}),eta^{-1}) = \sqrt{rac{eta}{2\pi}} \mathrm{exp}\left(-rac{eta}{2}(t-y(\mathbf{x},\mathbf{w}))^2
ight)$$

where the mean of this distribution is the regression function  $y(\mathbf{x}, \mathbf{w})$ .

## 3.1 Posterior and posterior predictive distribution [10 pts]

For a Bayesian treatment of linear regression, we need a prior probability distribution over the model parameters **w** with zero mean:

$$p(\mathbf{w}|lpha) = \mathcal{N}(\mathbf{w}|0,lpha^{-1}\mathbf{I})$$

where  $\alpha^{-1}I$  is the diagonal covariance matrix where all diagonal elements have the same variance  $\alpha^{-1}$ .

The posterior distribution has the following mean and inverse covariance matrix:

$$m_N = eta \mathbf{S}_N \Phi^T \mathbf{t}$$

$$\mathbf{S}_N^{-1} = \alpha \mathbf{I} + \beta \Phi^T \Phi$$

Hence, the posterior distribution can be written as

$$p(\mathbf{w}|\mathbf{t},lpha,eta) = \mathcal{N}(\mathbf{w}|\mathbf{m}_N,\mathbf{S}_N)$$

For making a prediction t at a new location x we use the posterior predictive distribution which is defined as

$$p(t|\mathbf{t},\mathbf{w},lpha,eta) = \int p(t|\mathbf{x},\mathbf{w},eta) p(\mathbf{w}|\mathbf{t},lpha,eta) d\mathbf{w}$$

Implement the posterior and posterior predictive distributions.

```
In [408...
          import numpv as np
          def posterior(Phi, t, alpha, beta, return inverse=False):
              """Computes mean and covariance matrix of the posterior distribution."""
              #raise NotImplementedError
              S N = np.linalg.inv(alpha * np.eye(Phi.shape[1]) + beta * Phi.T.dot(Phi) )
              m N = beta * S N.dot(Phi.T).dot(t)
              if return inverse:
                  return m N, S N, S N inv
              else:
                  return m N, S N
          def posterior predictive(Phi test, m N, S N, beta):
              """Computes mean and variances of the posterior predictive distribution."""
              y = Phi test.dot(m N)
              y var = 1 / beta + np.sum(Phi test.dot(S N) * Phi test, axis = 1)
              #raise NotImplementedError
              return y, y_var
```

# 3.2 Generating the data [5 pts]

Target values  $\mathbf{t}$  are generated from the design matrix  $\mathbf{X} \in \mathbb{R}^{N \times 1}$  with a linear function f which can also generate random noise of specified variance. Implement f below.

#### 3.3 Basis functions [5 pts]

For straight line fitting, we do not need to transform x with a basis function, which is equivalent to using an identity basis function. Other basis functions that are non-linear are necessary to model the non-linear relationship between input x and target t. Below is an example of non-linear basis functions: the polynomial. Implement the polynomial basis function. Note: You will not use it in the rest of the exercise but your implementation should be reasonable.

```
def identity_basis_function(x):
    return x

def polynomial_basis_function(x, power):
    #raise NotImplementedError
    return x**power

def expand(x, bf, bf_args=None):
    if bf_args is None:
        return np.concatenate([np.ones(x.shape), bf(x)], axis=1)
    else:
        return np.concatenate([np.ones(x.shape)] + [bf(x, bf_arg) for bf_arg in bf_args], axis=1)
```

## 3.3 Straight line fitting [5 pts]

For straight line fitting, we use a linear regression model of the form  $y(\mathbf{w}, \mathbf{x}) = w_0 + w_1 x$  and perform Bayesian inference for model parameters  $\mathbf{w}$ . Predictions are made with the posterior predictive distribution. Since this model has only two parameters,  $w_0$  and  $w_1$ , we can visualize the posterior density in 2D which is done in the first column of the following output. Rows use an increasing number of training data from a training dataset.

How does the dataset size affect the posterior density in the first column of the plots?

#### **Answer:**

Higher density (less spread out/less variance) with higher dataset size

```
In [411...
          from scipy import stats
          import matplotlib.pyplot as plt
          def plot data(x, t):
              plt.scatter(x, t, marker='o', c="k", s=20)
          def plot truth(x, y, label='Truth'):
              plt.plot(x, y, 'k--', label=label)
          def plot predictive(x, y, std, y label='Prediction', std label='Uncertainty', plot xy labels=True)
              y = y.ravel()
              std = std.ravel()
              plt.plot(x, y, label=y label)
              plt.fill between(x.ravel(), y + std, y - std, alpha = 0.5, label=std label)
              if plot xy labels:
                  plt.xlabel('x')
```

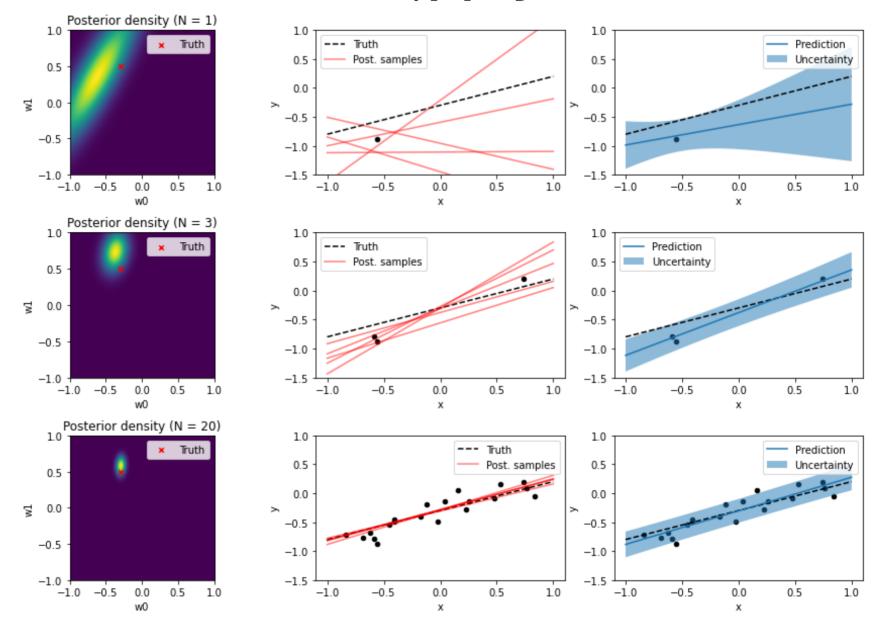
```
plt.vlabel('v')
def plot posterior samples(x, ys, plot xy labels=True):
    plt.plot(x, ys[:, 0], 'r-', alpha=0.5, label='Post. samples')
    for i in range(1, vs.shape[1]):
        plt.plot(x, vs[:, i], 'r-', alpha=0.5)
    if plot xv labels:
        plt.xlabel('x')
        plt.vlabel('v')
def plot posterior(mean, cov, w0, w1):
    resolution = 100
    grid x = grid v = np.linspace(-1, 1, resolution)
    grid flat = np.dstack(np.meshgrid(grid x, grid y)).reshape(-1, 2)
    densities = stats.multivariate normal.pdf(grid flat, mean=mean.ravel(), cov=cov).reshape(resol
    plt.imshow(densities, origin='lower', extent=(-1, 1, -1, 1))
    plt.scatter(w0, w1, marker='x', c="r", s=20, label='Truth')
    plt.xlabel('w0')
    plt.ylabel('w1')
def print comparison(title, a, b, a prefix='np', b prefix='br'):
    print(title)
    print('-' * len(title))
    print(f'{a prefix}:', a)
    print(f'{b prefix}:', b)
    print()
```

```
In [412...
```

import matplotlib.pyplot as plt
%matplotlib inline

```
# fix random seed so we always get the same data
np.random.seed(5)
# Trainina dataset sizes
N list = [1, 3, 20]
beta = 25.0
alpha = 2.0
# Training observations in [-1, 1)
X = np.random.rand(N list[-1], 1) * 2 - 1
# Trainina taraet values
t = f(X, noise variance=1/beta)
# Test observations
X \text{ test} = \text{np.linspace}(-1, 1, 100).\text{reshape}(-1, 1)
# Function values without noise
y true = f(X test, noise variance=0)
# Design matrix of test observations
Phi test = expand(X test, identity basis function)
plt.figure(figsize=(15, 10))
plt.subplots adjust(hspace=0.4)
for i, N in enumerate(N list):
    X N = X[:N]
    t N = t[:N]
    # Design matrix of training observations
    Phi N = expand(X N, identity basis function)
    # Mean and covariance matrix of posterior
    m N, S N = posterior(Phi N, t N, alpha, beta)
    # Mean and variances of posterior predictive
```

```
v, v var = posterior predictive(Phi test, m N, S N, beta)
# Draw 5 random weight samples from posterior and compute v values
w samples = np.random.multivariate normal(m N.ravel(), S N, 5).T
v samples = Phi test.dot(w samples)
plt.subplot(len(N list), 3, i * 3 + 1)
plot posterior(m N, S N, w0, w1)
plt.title(f'Posterior density (N = {N})')
plt.legend()
plt.subplot(len(N list), 3, i * 3 + 2)
plot data(X N, t N)
plot truth(X test, y true)
plot posterior samples(X test, y samples)
plt.vlim(-1.5, 1.0)
plt.legend()
plt.subplot(len(N list), 3, i * 3 + 3)
plot data(X N, t N)
plot truth(X test, y true, label=None)
plot predictive(X test, y, np.sqrt(y var))
plt.ylim(-1.5, 1.0)
plt.legend()
```



# 3.4 Bonus [10 pts] (optional)

We now fit a Gaussian basis function model to a noisy sinusoidal dataset for which we provide a sinusoidal function g with noise variance. Implement the gaussian basis function below.

The model uses 7 Gaussian basis functions with mean values equally distributed over [0,1] each having a standard deviation of 0.1. We then infer the model parameters **w** using the posterior predictive distribution.

How many parameters  $w_i$  are there? Run Bayesian inference and generate the plots. Do they follow the same trend as the ones generated by Bayesian linear regression?

#### **Answer:**

```
8 w_i from 0 - 7
```

Yes. Higher dataset size gives less variance.

```
def g(X, noise_variance):
    '''Sinusoidial function plus noise'''
    return 0.5 + np.sin(2 * np.pi * X) + noise(X.shape, noise_variance)

def gaussian_basis_function(x, mu, sigma=0.1):
    #raise NotImplementedError
    return np.exp(-(1/2) * (x-mu)*(x-mu)/(sigma*sigma))
```

```
import matplotlib.pyplot as plt
%matplotlib inline
from scipy.optimize import curve_fit

# fix random seed so we always get the same data
np.random.seed(5)

N_list = [3, 8, 20]

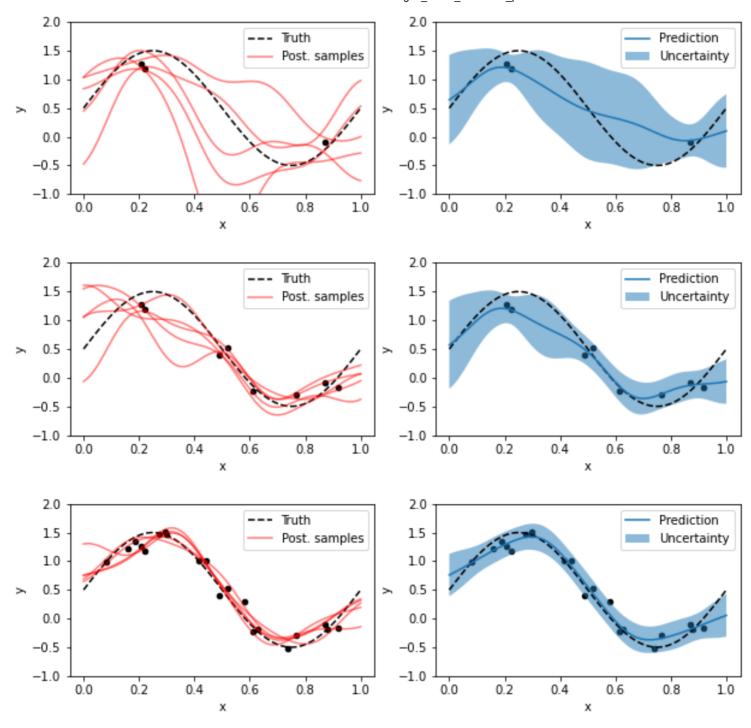
beta = 25.0
alpha = 2.0

# Training observations in [-1, 1)
```

```
X = np.random.rand(N list[-1], 1)
# Trainina taraet values
t = g(X, noise variance=1/beta)
# Test observations
X \text{ test} = \text{np.linspace}(0, 1, 100).\text{reshape}(-1, 1)
# Function values without noise
v true = g(X test, noise variance=0)
# Design matrix of test observations
Phi test = expand(X test, bf=gaussian basis function, bf args=np.linspace(0, 1, 7))
plt.figure(figsize=(10, 10))
plt.subplots adjust(hspace=0.4)
for i, N in enumerate(N list):
    X N = X[:N]
    t N = t[:N]
    # Design matrix of training observations
    Phi N = expand(X N, bf=gaussian basis function, bf args=np.linspace(0, 1, 7))
    # Mean and covariance matrix of posterior
    m N, S N = posterior(Phi N, t N, alpha, beta)
    # Mean and variances of posterior predictive
    y, y var = posterior predictive(Phi test, m N, S N, beta)
    # Draw 5 random weight samples from posterior and compute y values
    w samples = np.random.multivariate normal(m N.ravel(), S N, 5).T
    y samples = Phi test.dot(w samples)
    plt.subplot(len(N list), 2, i * 2 + 1)
    plot data(X N, t N)
    plot truth(X test, y true)
    plot posterior samples(X test, y samples)
```

```
plt.ylim(-1.0, 2.0)
plt.legend()

plt.subplot(len(N_list), 2, i * 2 + 2)
plot_data(X_N, t_N)
plot_truth(X_test, y_true, label=None)
plot_predictive(X_test, y, np.sqrt(y_var))
plt.ylim(-1.0, 2.0)
plt.legend()
```



In [ ]: