Sheet 5

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In [1]: import numpy as np
 from matplotlib import pyplot as plt

1 Bayes: Signal or Noise?

$$p(\text{gamma ray} \mid \text{target direction}) = \frac{p(\text{target direction} \mid \text{gamma ray}) \cdot p(\text{gamma ray})}{p(\text{background})} = \frac{0.95 \cdot 0.1}{0.9} \approx 0.11$$

$$p(\text{target direction})$$

2 Bayes Classifier

a)

$$L(y,\hat{y}) = \left[egin{matrix} 0 & 1 \ 10 & 0 \end{array}
ight]$$

where:

- y=0 is the true class, and $\hat{y}=0$ or $\hat{y}=1$ are the predicted classes.
- y=1 is the true class, and $\hat{y}=0$ or $\hat{y}=1$ are the predicted classes.

With P(y=1|x)=p(x) and P(y=0|x)=1-p(x) we calculate the conditional risk $R(\hat{y}|x)$ for each possible predicted label \hat{y} :

1. **For** $\hat{y} = 0$:

$$R(0|x) = L(0,0) \cdot P(y=0|x) + L(1,0) \cdot P(y=1|x) = 0 \cdot (1-p(x)) + 10 \cdot p(x) = 10 \cdot p(x)$$

2. **For** $\hat{y} = 1$:

$$R(1|x) = L(0,1) \cdot P(y=0|x) + L(1,1) \cdot P(y=1|x) = 1 \cdot (1-p(x)) + 0 \cdot p(x) = 1 - p(x)$$

To minimize the conditional risk, choose $\hat{y} = 0$ if R(0|x) < R(1|x), and $\hat{y} = 1$ otherwise.

Thus, we need:

$$10 \cdot p(x) < 1 - p(x)$$

Rearrange this inequality:

$$11 \cdot p(x) < 1 \quad \Rightarrow \quad p(x) < \frac{1}{11}$$

The optimal Bayes classifier is:

$$\hat{y} = egin{cases} 0 & ext{if } p(x) < rac{1}{11} \ 1 & ext{if } p(x) \geq rac{1}{11} \end{cases}$$

This asymmetric loss matrix is appropriate in situations where the consequences of a false negative (predicting y=0 when y=1) are much more severe than a false positive (predicting y=1 when y=0).

b)

The Bayes classifier minimizes the **expected loss** by choosing the prediction \hat{y} that minimizes the **conditional risk** $R(\hat{y}|x)$ given x. For each potential prediction $\hat{y} \in \{0, 1, \dots, k\}$, we calculate $R(\hat{y}|x)$.

1. Conditional Risk for $\hat{y}=j$, where $j\in\{1,\ldots,k\}$: If we predict class j, the conditional risk is:

$$R(j|x) = \sum_{i=1}^k L(i,j) \cdot P(y=i|x)$$

$$R(j|x) = \sum_{\substack{i=1\\i \neq j}}^{k} P(y = i|x) = 1 - P(y = j|x)$$

2. Conditional Risk for the "Reject" Option $\hat{y}=0$: If we choose to reject, the conditional risk is:

$$R(0|x) = \sum_{i=1}^{k} L(i,0) \cdot P(y=i|x) = \alpha \sum_{i=1}^{k} P(y=i|x) = \alpha$$

since the probabilities P(y=i|x) sum to 1.

To minimize the conditional risk, we compare R(j|x) for $j \in \{1, \ldots, k\}$ with $R(0|x) = \alpha$.

• For each class j, we will choose $\hat{y} = j$ if:

$$R(j|x) < \alpha \quad \Rightarrow \quad 1 - P(y = j|x) < \alpha \quad \Rightarrow \quad P(y = j|x) > 1 - \alpha$$

• If no class j satisfies P(y=j|x)>1-lpha, we choose the "reject" option, $\hat{y}=0$.

Thus, the optimal Bayes classifier is:

$$\hat{y} = \left\{ egin{aligned} j & ext{if } P(y=j|x) > 1-lpha ext{ for some } j \in \{1,\ldots,k\} \ 0 & ext{if } P(y=j|x) \leq 1-lpha ext{ for all } j \in \{1,\ldots,k\} \end{aligned}
ight.$$

The parameter α controls the threshold for rejection:

- A lower α means that we are less willing to reject; we require a higher confidence in our prediction before choosing to classify. As α → 0,
 the classifier rarely chooses the "reject" option.
- A **higher** α increases the likelihood of rejection, as the threshold $1-\alpha$ becomes lower, requiring less certainty to avoid rejection.

This approach adds flexibility, allowing for rejection when the classifier's confidence is insufficient, thus potentially reducing overall decision-making risk.

3 QDA

(a)

```
In [24]: pts = np.load('data/dataid.npy')
labels = np.load('data/labelsid.npy')

x0, x1 = pts[labels == 0], pts[labels == 1]
plt.hist(x0, bins=10, alpha=0.5, density=True, color='r')
plt.hist(x1, bins=10, alpha=0.5, density=True, color='b')

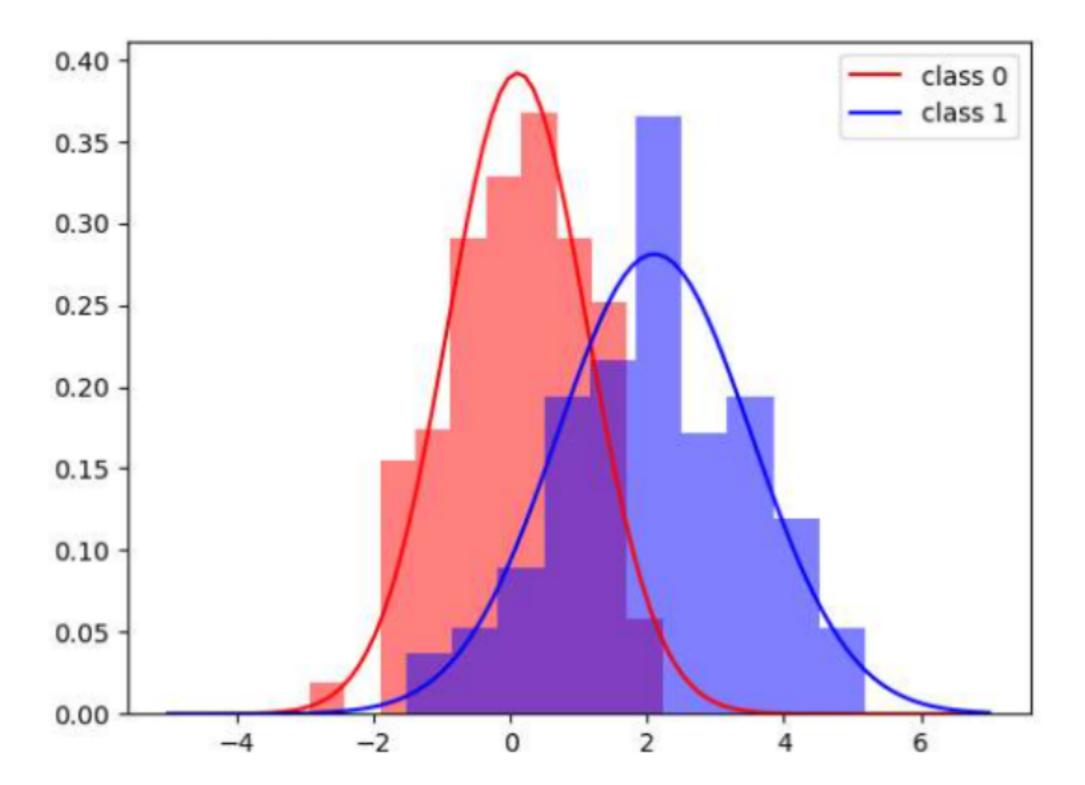
mean0, mean1 = np.mean(x0), np.mean(x1)
var0, var1 = np.var(x0), np.var(x1)

def gaussian(x, mean, var):
    return np.exp(-0.5*(x-mean)**2/var)/np.sqrt(2*np.pi*var)

x = np.linspace(-5, 7, 100)

plt.plot(x, gaussian(x, mean0, var0), label='class 0', color='r')
plt.plot(x, gaussian(x, mean1, var1), label='class 1', color='b')
plt.legend()
```

Out[24]: <matplotlib.legend.Legend at 0x23470f65d00>



(b)

```
In [26]: x = np.linspace(-10, 10, 100)

like0 = gaussian(x, mean0, var0)
like1 = gaussian(x, mean1, var1)

prior0, prior1 = 0.5, 0.5

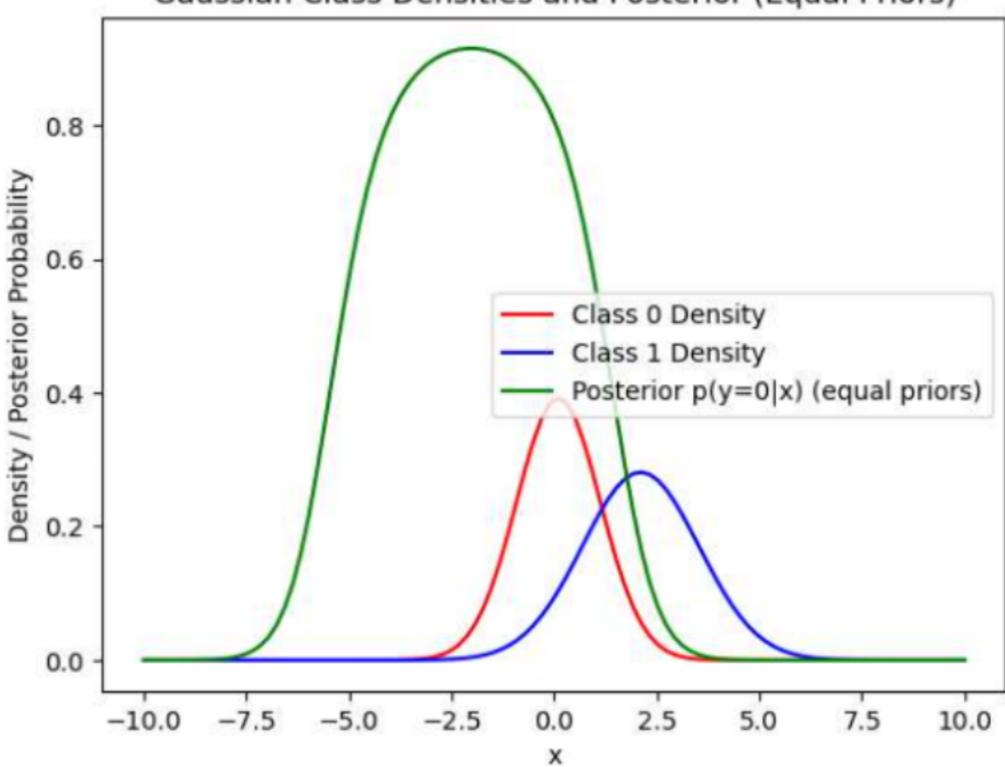
post0 = like0*prior0/(like0*prior0 + like1*prior1)

plt.plot(x, like0, label="Class 0 Density", color="r")
plt.plot(x, like1, label="Class 1 Density", color="b")
plt.plot(x, post0, label="Posterior p(y=0|x) (equal priors)", color="green")
```

```
plt.title("Gaussian Class Densities and Posterior (Equal Priors)")
plt.xlabel("x")
plt.ylabel("Density / Posterior Probability")
plt.legend()
```

Out[26]: <matplotlib.legend.Legend at 0x23470f833a0>

Gaussian Class Densities and Posterior (Equal Priors)



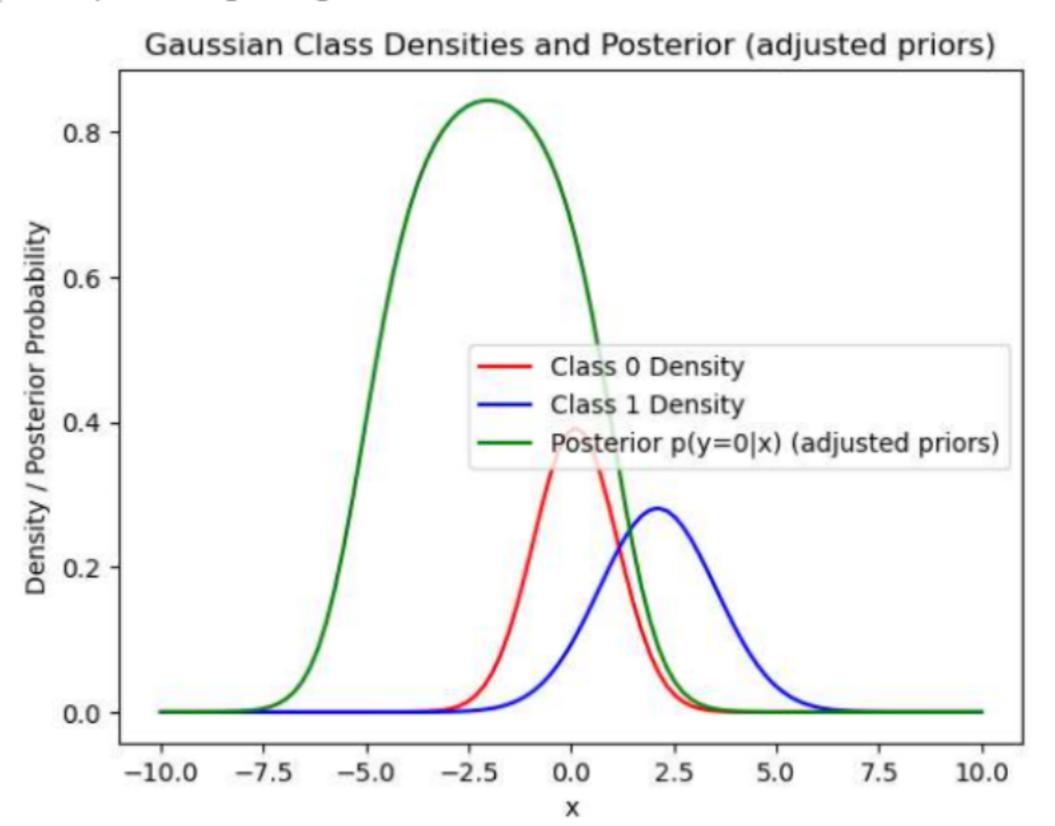
```
In [28]: prior0, prior1 = 1/3, 2/3

post0 = like0*prior0/(like0*prior0 + like1*prior1)

plt.plot(x, like0, label="Class 0 Density", color="r")
plt.plot(x, like1, label="Class 1 Density", color="b")
```

```
plt.plot(x, post0, label="Posterior p(y=0|x) (adjusted priors)", color="green")
plt.title("Gaussian Class Densities and Posterior (adjusted priors)")
plt.xlabel("x")
plt.ylabel("Density / Posterior Probability")
plt.legend()
```

Out[28]: <matplotlib.legend.Legend at 0x2346f84fe80>



4 Trees and Random Forests

a)

Misclassification Rate =
$$1 - \max(p(y = c))$$

$$\mathrm{Gini} = 1 - \sum_{c=1}^2 p(y=c)^2$$

$$\text{Entropy} = -\sum_{c=1}^{2} p(y=c) \log_2 p(y=c)$$

1. Node 1 (300, 100):

$$\begin{aligned} \text{Misclassification Rate} &= 1 - 0.75 = 0.25\\ \text{Gini Impurity} &= 1 - (0.75^2 + 0.25^2) = 1 - (0.5625 + 0.0625) = 0.375\\ \text{Entropy} &= -(0.75\log_2 0.75 + 0.25\log_2 0.25) \approx 0.811 \end{aligned}$$

2. Node 2 (100, 300):

$$\begin{aligned} \text{Misclassification Rate} &= 1 - 0.75 = 0.25 \\ \text{Gini Impurity} &= 1 - (0.25^2 + 0.75^2) = 0.375 \\ \text{Entropy} &= -(0.25\log_2 0.25 + 0.75\log_2 0.75) \approx 0.811 \end{aligned}$$

Weighted Averages for Split A:

$$\begin{split} \text{Misclassification Rate} &= 0.25 \times \frac{400}{800} + 0.25 \times \frac{400}{800} = 0.25 \\ \text{Gini Impurity} &= 0.375 \times \frac{400}{800} + 0.375 \times \frac{400}{800} = 0.375 \\ \text{Entropy} &= 0.811 \times \frac{400}{800} + 0.811 \times \frac{400}{800} \approx 0.811 \end{split}$$

4. Node 1 (200, 0):

$$\begin{aligned} \text{Misclassification Rate} &= 1 - 1 = 0 \\ \text{Gini Impurity} &= 1 - (1^2 + 0^2) = 0 \\ \text{Entropy} &= -(1 \cdot \log_2 1 + 0 \cdot \log_2 0) = 0 \end{aligned}$$

5. Node 2 (200, 400):

Misclassification Rate =
$$1-0.667 \approx 0.333$$

Gini Impurity = $1-(0.333^2+0.667^2)=1-(0.111+0.444)=0.444$
Entropy = $-(0.333\log_2 0.333+0.667\log_2 0.667)\approx 0.918$

Weighted Averages for Split B:

$$\begin{aligned} \text{Misclassification Rate} &= 0 \times \frac{200}{800} + 0.333 \times \frac{600}{800} \approx 0.25 \\ \text{Gini Impurity} &= 0 \times \frac{200}{800} + 0.444 \times \frac{600}{800} = 0.333 \\ \text{Entropy} &= 0 \times \frac{200}{800} + 0.918 \times \frac{600}{800} \approx 0.688 \end{aligned}$$

- Misclassification Rate: Both splits have the same misclassification rate of 0.25.
- Gini Impurity: Split B has a lower Gini impurity (0.333 vs. 0.375), so it would be preferred by the Gini criterion.
- Entropy: Split B also has a lower entropy (0.688 vs. 0.811), so it would be preferred by the entropy criterion.

Overall, Split B would be the preferred split based on both Gini impurity and entropy.

(b)

```
In []: # Load the data
pts = np.load('data/datald.npy')
labels = np.load('data/labelsld.npy')

# TODO: Sort the points to easily split them

sorted_pts = np.sort(pts)
sorted_labels = labels[np.argsort(pts)]

# TODO: Implement or find implementation for Gini impurity, entropy and misclassifcation rate

def probabilities(partition):
    # divide counts by size of dataset to get cluster probabilites
    return np.unique(partition, return_counts=True)[1] / len(partition)
```

```
def gini_impurity(partition):
    p = probabilities(partition)
    return 1 - np.sum(p**2)
def entropy(partition):
    p = probabilities(partition)
    return -np.sum(p * np.log2(p))
def misclassification_rate(partition):
    p = probabilities(partition)
    return 1 - np.max(p)
def compute_split_measure(1, 10, 11, method):
    p0 = probabilities(10)
    p1 = probabilities(11)
    p = probabilities(1)
    return method(p) - (len(10) * method(p0) + len(11) * method(p1)) / (len(1))
# TODO: Iterate over the possible splits, evaulating and saving the three criteria for each one
# TODO: Then, Compute the split that each criterion favours and visualize them
        (e.g. with a histogram for each class and vertical lines to show the splits)
```

(c)

In []:

```
In [7]: # Load the dijet data
features = np.load('data/dijet_features_normalized.npy')
labels = np.load('data/dijet_labels.npy')

# TODO: define train, val and test splits as specified (make sure to shuffle the data before splitting it!)

In [8]: from sklearn.ensemble import RandomForestClassifier

# TODO: train a random forest classifier for each combination of specified hyperparameters
# and evaluate the performances on the validation set.
```

In [9]: # TODO: for your preferred configuration, evaluate the performance of the best configuration on the test set