Naive Bayes

Naive Bayes is a very simple but powerful classification method. For a given object x, Naive Bayes calculates x's probability to belong to each class y_i ($i = 1, \dots, k$), using the Bayes' theorem:

$$P\left(y_i \mid x\right) = \frac{P(y_i) \; P\left(x_1, \; \cdots, \; x_m \mid y_i\right)}{P\left(x_1, \; \cdots, \; x_m\right)}.$$

Additionally, it assumes that the features are independent from each other (which is the reason why it is called naive):

$$P(x_1, \dots, x_m \mid y_i) = \prod_{i=1}^m P(x_i \mid y_i).$$

So, we obtain:

$$P(y_i \mid x) = \frac{P(y_i) \prod_{j=1}^m P(x_j \mid y_i)}{P(x_1, \dots, x_m)}.$$

For a given object x, Naive Bayes will output the the Maximum a Posteriori (MAP) estimate:

$$\hat{\mathbf{y}} = \operatorname{arg\ max}_{\mathbf{y}} P\left(\mathbf{y} \mid \mathbf{x}\right) = \operatorname{arg\ max}_{\mathbf{y}} \frac{P\left(\mathbf{y}\right) \prod_{j=1}^{m} P\left(x_{j} \mid \mathbf{y}\right)}{P\left(x_{1}, \ \cdots, \ x_{m}\right)}.$$

Note that $P(x_1, \dots, x_m)$ is constant. Thus, we can drop it to obtain:

$$\hat{y} = \arg\max_{y} P(y_i) \prod_{i=1}^{m} P(x_i \mid y).$$

Practical example

First of all, we do all the necessary imports and load the Mushroom dataset.

In [1]:

```
import pandas as pd
import matplotlib.pyplot as plt

from sklearn.preprocessing import LabelEncoder
from sklearn.model_selection import train_test_split
from sklearn.naive_bayes import BernoulliNB
from sklearn.metrics import accuracy_score, roc_curve, auc

# setting random seed
seed = 10

data = pd.read_csv('data/mushroom.csv')

# We drop the 'stalk-root' feature because it is the only one containing missing values.
data = data.drop('stalk-root', axis=1)
data.head()
```

Out[1]:

| | cap- shape | cap- surface | cap- color | bruises? | odor | gill- attachment | _ | _ | | | | color- below- | | |
|---|---------------|-----------------|---------------|----------|------|---------------------|---|---|---|---|-------|------------------|---|---|
| 0 | х | s | n | t | р | f | С | n | k | е | w | w | p | w |
| 1 | х | s | у | t | a | f | С | b | k | е | w | W | р | w |
| 2 | b | s | w | t | I | f | С | b | n | е | W | W | р | w |
| 3 | х | у | w | t | р | f | С | n | n | е | w | w | р | w |
| 4 | х | S | g | f | n | f | w | b | k | t | w | W | р | w |

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Unfortunately, scikit-learn does not implement the classical Naive Bayes algorithm which calculates the conditional probabilities $P(x_j | y_i)$ as the proportion of objects from class y_i that assume each particular categorical value for feature j. However, scikit-learn contains the BernoulliNB class which assumes that data is distributed according to multivariate Bernoulli distributions.

So, for the Mushroom dataset, we can transform each categorical feature to dummy variables. **Note that such a conversion clearly violates the indepence assumption between features.** However, Naive Bayes has been proven to achieve good performance in several applications where indepence is violated (for example, in text classication).

In [2]:

```
# Creating a new DataFrame representation for each feature as dummy variables.
dummies = [pd.get dummies(data[c]) for c in data.drop('label', axis=1).columns]
# Concatenating all DataFrames containing dummy variables.
binary data = pd.concat(dummies, axis=1)
# Getting binary_data as a numpy.array.
X = binary_data.values
# Getting the labels.
le = LabelEncoder()
y = le.fit transform(data['label'].values)
# Splitting the binary dataset into train and test sets.
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.34, stratify=y, random_state=seed)
# Creates a BernoulliNB. binarize=None indicates that there is no need to binarize the input data.
nb = BernoulliNB(binarize=None)
nb.fit(X_train, y_train)
y pred = nb.predict(X test)
accuracy = accuracy_score(y_test, y_pred)
print('BernoulliNB accuracy score: {}'.format(accuracy))
```

BernoulliNB accuracy score: 0.9464350343829171

In [3]:

```
# Getting the probabilities for each class.
y prob = nb.predict proba(X test)
# Calculating ROC curve and ROC AUC.
false positive rate, true positive rate, thresholds = roc curve(y test, y prob[:, 1])
roc auc = auc(false positive rate, true positive rate)
# Plotting ROC curve.
lw = 2
plt.plot(false_positive_rate, true_positive_rate, color='blue', lw=lw, label='ROC curve (area = {:.4f})'.format(roc_auc)
plt.plot([0, 1], [0, 1], color='red', lw=lw, linestyle='--', label='Random classifier')
plt.xlim([0.0, 1.0])
plt.ylim([0.0, 1.05])
plt.xlabel('False Positive Rate')
plt.ylabel('True Positive Rate')
plt.title('ROC')
plt.legend(loc="lower right")
plt.show()
```

