

Exercise - Train and evaluate a classification model

8 minutes

This module requires a sandbox to complete. You have used 1 of 10 sandboxes for today. More sandboxes will be available tomorrow.

Activate sandbox





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Classification

Supervised machine learning techniques involve training a model to operate on a set of features and predict a label using a dataset that includes some already-known label values. You can think of this function like this, in which **y** represents the label we want to predict and **X** represents the vector of features the model uses to predict it.

$$y = f([x_1, x_2, x_3, ...])$$

Classification is a form of supervised machine learning in which you train a model to use the features (the *x* values in our function) to predict a label (*y*) that calculates the probability of the observed case belonging to each of a number of possible classes, and predicting an appropriate label. The simplest form of classification is *binary* classification, in which the label is 0 or 1, representing one of two classes; for example, "True" or "False"; "Internal" or "External"; "Profitable" or "Non-Profitable"; and so on.

Binary Classification

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```
import pandas as pd

# load the training dataset
!wget https://raw.githubusercontent.com/MicrosoftDocs/mslearn-introduction-to-machine-learning,
diabetes = pd.read_csv('diabetes.csv')
diabetes.head()
```

This data consists of diagnostic information about some patients who have been tested for diabetes. Scroll to the right if necessary, and note that the final column in the dataset (**Diabetic**) contains the value **0** for patients who tested negative for diabetes, and **1** for patients who tested positive. This is the label that we will train our model to predict; most of the other columns (**Pregnancies**, **PlasmaGlucose**, **DiastolicBloodPressure**, and so on) are the features we will use to predict the **Diabetic** label.

Let's separate the features from the labels - we'll call the features X and the labe 💥 🌊 😎 🥩



```
# Separate features and labels
features = ['Pregnancies','PlasmaGlucose','DiastolicBloodPressure','TricepsThickness','SerumIn:
label = 'Diabetic'
X, y = diabetes[features].values, diabetes[label].values

for n in range(0,4):
    print("Patient", str(n+1), "\n Features:",list(X[n]), "\n Label:", y[n])
```

Now let's compare the feature distributions for each label value.

```
from matplotlib import pyplot as plt
%matplotlib inline

features = ['Pregnancies','PlasmaGlucose','DiastolicBloodPressure','TricepsThickness','SerumIn:
    for col in features:
        diabetes.boxplot(column=col, by='Diabetic', figsize=(6,6))
        plt.title(col)
plt.show()
```



For some of the features, there's a noticeable difference in the distribution for each label value. In particular, **Pregnancies** and **Age** show markedly different distributions for diabetic patients than for non-diabetic patients. These features may help predict whether or not a patient is diabetic.

Split the data

Our dataset includes known values for the label, so we can use this to train a classifier so that it finds a statistical relationship between the features and the label value; but how will we know if our model is any good? How do we know it will predict correctly

naber values, use only some or it to train the moder, and hold back some to test the trained moder - enabling us to compare the predicted labels with the already known labels in the test set.

In Python, the **scikit-learn** package contains a large number of functions we can use to build a machine learning model - including a **train_lear_and** contains a large number of functions we can use to build a machine learning model - including a **train_lear_and** contains and hold back 30% for testing

70% for training and hold back 30% for testing. OK, now we're ready to train our model by fitting the training features (**X_train**) to the training labels (**y_train**). There are various algorithms we can use to train the model. In this example, we'll use *Logistic Regression*, which (despite its name) is a well-established algorithm for classification. In addition to the training features and labels, we'll need to set a *regularization* parameter. This is used to counteract any bias in the sample, and help the model generalize well by avoiding *overfitting* the model to the training data.

Note: Parameters for machine learning algorithms are generally referred to as *hyperparameters* (to a data scientist, *parameters* are values in the data itself - *hyperparameters* are defined externally from the data!)

Now we've trained the model using the training data, we can use the test data we held back to evaluate how well it predicts. Again, **scikit-learn** can help us do this. Let's start by using the model to predict labels for our test set, and compare the predicted labels to the known labels:



The arrays of labels are too long to be displayed in the notebook output, so we can only compare a few values. Even if we printed out all of the predicted and actual labels, there are too many of them to make this a sensible way to evaluate the model. Fortunately, **scikit-learn** has a few more tricks up its sleeve, and it provides some metrics that we can use to evaluate the model.

The most obvious thing you might want to do is to check the *accuracy* of the predictions - in simple terms, what proportion of the labels did the model predict correctly?

The accuracy is returned as a decimal value - a value of 1.0 would mean that the model got 100% of the predictions right; while an accuracy of 0.0 is, well, pretty useless!

Summary

Here we prepared our data by splitting it into test and train datasets, and applied logistic regression - a way of applying binary labels to our data. Our model was able to predict whether patients had diabetes with what appears like reasonable accuracy. But is this good enough? In the next notebook we will look at alternatives to accuracy that can be much more useful in machine learning.

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Next unit: Evaluate classification models

Continue >



