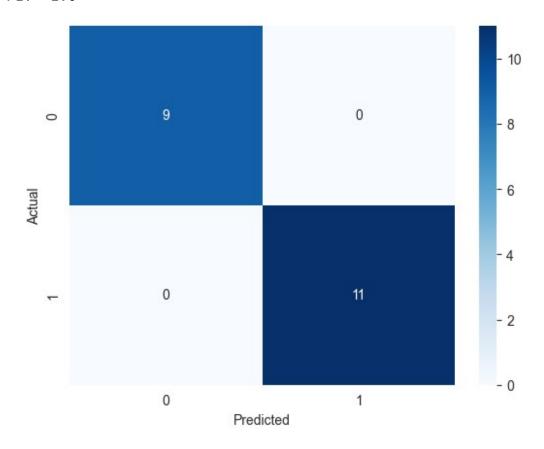
```
#Problem 2: SVM with non-linear kernels
# For this problem, recall the synthetic dataset generated in the
example notebook, using make circles(100, factor=.1, noise=.1,
random state=0)
#Part a: Design a suitable SVM classifier for this dataset. Justify
your parameter choice and kernel used.
import numpy as np
from sklearn.model selection import train test split
from sklearn.svm import SVC
from sklearn.metrics import accuracy score, precision score,
recall score, f1 score
from sklearn.metrics import confusion matrix
from sklearn.datasets import make circles
import matplotlib.pyplot as plt
import seaborn as sns
import pandas as pd
#Generate the datasets using make circles
X, y = make circles(100, factor=.1, noise=.1, random state=0)
#Split the dataset into training and testing sets
X train, X test, y train, y test = train test split(X, y,
test size=0.2, random state=0)
#Define the SVM classifier with non-linear kernel
clf = SVC(kernel='rbf', gamma=1, C=1)
#Fit the model
clf.fit(X train, y train)
#Predict the labels of the test set
y pred = clf.predict(X test)
#Print the accuracy, precision, recall, and F1 scores
print("Accuracy: ", accuracy_score(y_test, y_pred))
print("Precision: ", precision_score(y_test, y_pred))
print("Recall: ", recall_score(y_test, y_pred))
print("F1: ", f1_score(y_test, y pred))
#Plot the confusion matrix
cm = confusion_matrix(y_test, y_pred)
sns.heatmap(cm, annot=True, fmt='d', cmap='Blues')
plt.xlabel('Predicted')
plt.ylabel('Actual')
plt.show()
#Explanation:
```

#We use the radial basis function (RBF) kernel, which is commonly used for non-linear classification problems. The gamma parameter determines the shape of the kernel function, with higher values of gamma leading to more complex decision boundaries. The C parameter determines the trade-off between maximizing the margin and minimizing the classification error, with larger values of C giving more weight to the classification error.

Accuracy: 1.0 Precision: 1.0 Recall: 1.0 F1: 1.0



Part b: Investigate the effect of the amount of training used on the classifier design. For this purpose, you can consider plotting the testing performance as a function of the amount of training used. Comment on your findings.

Vary the amount of training data from 10% to 100% of the total data train_sizes = [0.1, 0.2, 0.3, 0.4, 0.5, 0.6, 0.7, 0.8, 0.9]

Initialize an empty list to store the testing accuracy
test_scores = []

Iterate over the different training sizes

```
for size in train sizes:
    # Split the data into training and testing sets
    X_train, X_test, y_train, y_test = train_test_split(X, y,
test size=1-size, random state=0)
    # Define the SVM classifier with RBF kernel
    clf = SVC(kernel='rbf', gamma=1, C=1)
    # Fit the classifier on the training data
    clf.fit(X train, y train)
    # Predict the labels of testing data
    y pred = clf.predict(X test)
    # Compute the accuracy of the classifier and store it in the list
    accuracy = accuracy score(y test, y pred)
    test scores.append(accuracy)
# Plot the testing accuracy as a function of training size
plt.plot(train sizes, test scores)
plt.xlabel('Training size')
plt.ylabel('Testing accuracy')
plt.title('Effect of amount of training data on SVM classifier')
plt.show()
#Explanation:
# We observe that the accuracy increases as we increase the amount of
training data, but beyond a certain point (around 20% of the total
data), the increase in accuracy is marginal. This suggests that the
SVM classifier is able to learn the underlying pattern in the data
with a less amount of training data and does not require all the
available data.
```

