6주. Decision Tree, RF, SVM			
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PimalndiansDiabetes dataset을 가지고 Classification 을 하고자 한다. (마지막의 diabetes 컬럼 이 class label 임)

- Q1 (4점) scikit-learn에서 제공하는 DecisionTree, RandonForest, support vector machine 알고리즘를 이용하여 **PimalndiansDiabetes dataset**에 대한 분류 모델을 생성하고 accuracy를 비교하시오.
- 10-fold cross validation을 실시하여 mean accuracy를 비교한다
- 각 알고리즘의 hyper parameter 의 값은 default value를 이용한다.

Source code:

```
// source code 의 폰트는 Courier10 BT Bold으로 하시오
pip install pydot
import pandas as pd
import numpy as np
from sklearn.tree import DecisionTreeClassifier, export graphviz
from sklearn.ensemble import RandomForestClassifier
from sklearn import svm
from sklearn.model selection import train test split
from sklearn.metrics import confusion matrix
import pandas as pd
import pydot
from sklearn.model selection import KFold
#decision tree
df=pd.read csv('C:/Users/ATIV/Desktop/Deeplearning Cloud/dataset 09
14/PimaIndiansDiabetes.csv')
df X = df.loc[:, df.columns != 'diabetes']
df y = df['diabetes']
df X=df X.values
df_y=df_y.values
seed=32152339
n fold=10
kf = KFold(n_splits=n_fold, random_state=seed, shuffle=True)
10 fold
# Define learning model
DT model = DecisionTreeClassifier(random state=seed)
RF model = RandomForestClassifier(random state=seed)
```

```
SVM model = svm.SVC()
models = [DT model,RF model,SVM model]
for model in models:
   acc = np.zeros(n fold)
  i = 0
  print(model)
   for train_index, test_index in kf.split(df_X):
      print("fold:", i)
      train_X, test_X = df_X[train_index], df_X[test_index]
      train_y, test_y = df_y[train_index], df_y[test_index]
      # Train the model using the training sets
      model.fit(train X, train y)
      # Make predictions using the testing set
      pred_y = model.predict(test_X)
      #print(pred y)
      acc[i] = accuracy_score(test_y, pred_y)
     # print('Accuracy : {0:3f}'.format(acc[i]))
      i += 1
  print(model)
   #print(f"{n fold} fold:", acc)
   print("mean accuracy :", np.mean(acc))
   print('----')
```

실행화면 캡쳐:

```
DecisionTreeClassifier(random state=32152339)
fold: 0
fold: 1
fold: 2
fold: 3
fold: 4
fold: 5
fold: 6
fold: 7
fold: 8
fold: 9
DecisionTreeClassifier(random state=32152339)
mean accuracy : 0.6980177717019821
_____
RandomForestClassifier(random_state=32152339)
fold: 0
fold: 1
fold: 2
fold: 3
fold: 4
fold: 5
fold: 6
fold: 7
fold: 8
fold: 9
RandomForestClassifier(random_state=32152339)
mean accuracy: 0.7617737525632263
______
SVC()
fold: 0
fold: 1
fold: 2
fold: 3
fold: 4
fold: 5
fold: 6
fold: 7
fold: 8
fold: 9
SVC()
mean accuracy: 0.7564593301435407
```

-> randomforest가 0.76으로 가장 성능이 좋다.

Q2. (3점) 다음의 조건에 따라 support vector machine 알고리즘를 이용하여 **PimalndiansDiabetes dataset**에 대한 분류 모델을 생성하고 accuracy를 비교하시오.

- hyper parameter 중 kernel 에 대해 linear, poly, rbf, sigmoid, precomputed를 각 각 테스트하여 어떤 kernel 이 가장 높은 accuracy를 도출하는지 확인하시오.
- 10-fold cross validation을 실시하여 mean accuracy를 비교한다

Source code:

```
// source code 의 폰트는 Courier10 BT Bold으로 하시오
kernels=['linear', 'poly', 'rbf', 'sigmoid','precomputed']
for k in kernels:
   print(f'svc kernels : {k}')
   model = svm.SVC(kernel=k)
   acc = np.zeros(n_fold)
                          # 10 fold
                      # fold no
   for train index, test index in kf.split(df X):
      print("fold:", i)
      train X, test X = df X[train index], df X[test index]
      train_y, test_y = df_y[train_index], df_y[test_index]
      if k != 'precomputed':
          model.fit(train X, train y)
          pred_y = model.predict(test_X)
      else:
          gram_train = np.dot(train_X, train_X.T)
          model.fit(gram train, train y)
          gram test = np.dot(test X, train X.T)
          pred y = model.predict(gram test)
      acc[i] = accuracy score(test y, pred y)
      i += 1
   print(model)
  # print(f"{n fold} fold :", acc)
   print("mean accuracy :", np.mean(acc))
   print()
```

실행화면 캡쳐:

Deep Learning/Cloud

```
svc kernels : linear
fold: 0
fold: 1
fold: 2
fold: 3
fold: 4
fold: 5
fold: 6
fold: 7
fold: 8
fold: 9
SVC(kernel='linear')
mean accuracy: 0.7733937115516063
svc kernels : poly
fold: 0
fold: 1
fold: 2
fold: 3
fold: 4
fold: 5
fold: 6
fold: 7
fold: 8
fold: 9
SVC(kernel='poly')
mean accuracy: 0.7564422419685578
svc kernels : rbf
fold: 0
fold: 1
fold: 2
fold: 3
fold: 4
fold: 5
fold: 6
fold: 7
fold: 8
fold: 9
SVC()
mean accuracy: 0.7564593301435407
svc kernels : sigmoid
fold: 0
fold: 1
fold: 2
fold: 3
fold: 4
fold: 5
fold: 6
fold: 7
fold: 8
fold: 9
SVC(kernel='sigmoid')
mean accuracy: 0.49743677375256323
svc kernels : precomputed
fold: 0
fold: 1
fold: 2
fold: 3
fold: 4
fold: 5
fold: 6
fold: 7
fold: 8
SVC(kernel='precomputed')
mean accuracy: 0.7733937115516063
```

-> linear kernel과 precomputed kernel이 0.77로 accuracy 성능이 좋다.

```
Q3. (3점) 다음의 조건에 따라 Random Forest 알고리즘를 이용하여 PimalndiansDiabetes
dataset에 대한 분류 모델을 생성하고 accuracy를 비교하시오.
-다음의 hyper parameter를 테스트 하시오
. n_estimators : 100, 200, 300, 400, 500
. max_features : 1, 2, 3, 4, 5
어떤 조합이 가장 높은 accuracy를 도출하는지 확인하시오.
- 10-fold cross validation을 실시하여 mean accuracy를 비교한다
```

Source code:

```
// source code 의 폰트는 Courier10 BT Bold으로 하시오
n = [100, 200, 300, 400, 500]
\max \text{ features} = [1, 2, 3, 4, 5]
mean acc=[]
for n_estimator in n_estimators:
            print(f'n estimators : {n estimator}')
             for max feature in max features:
{\tt model=RandomForestClassifier(n\_estimators=n\_estimator, max\_features=n\_estimator, max\_featur
max feature)
                         acc = np.zeros(n fold)
                                                                                                                      # 10 fold
                         i = 0
                         for train index, test index in kf.split(df X):
                                     #print("fold:", i)
                                     train X, test X = df X[train index], df X[test index]
                                     train_y, test_y = df_y[train_index], df_y[test_index]
                                     model.fit(train X, train y)
                                     pred y = model.predict(test X)
                                     acc[i] = accuracy score(test y, pred y)
                                      i += 1
                        print(model)
                            print(f"{n fold} fold :", acc)
                        print("mean accuracy :", np.mean(acc))
mean_acc.append([n_estimator,max_feature,round(np.mean(acc),5)])
                         print()
```

```
mean acc
 mean acc=pd.DataFrame(mean acc)
 mean acc.columns=['n estimators','max features','mean accuracy']
 mean acc[mean acc.mean accuracy==np.max(mean acc.mean accuracy)]
실행화면 캡쳐:
n_estimators : 100
RandomForestClassifier(max_features=1)
mean accuracy: 0.756578947368421
RandomForestClassifier(max features=2)
mean accuracy: 0.7630382775119618
RandomForestClassifier(max features=3)
mean accuracy: 0.7617395762132604
RandomForestClassifier(max features=4)
mean accuracy: 0.7486671223513328
RandomForestClassifier(max features=5)
mean accuracy: 0.7435064935064937
n_estimators : 200
RandomForestClassifier(max_features=1, n_estimators=200)
mean accuracy: 0.7540157211209844
RandomForestClassifier(max_features=2, n_estimators=200)
mean accuracy: 0.7682330827067669
RandomForestClassifier(max_features=3, n_estimators=200)
mean accuracy: 0.7630895420369106
RandomForestClassifier(max_features=4, n_estimators=200)
mean accuracy: 0.7604408749145593
RandomForestClassifier(max_features=5, n_estimators=200)
mean accuracy: 0.7487183868762817
n_estimators : 300
RandomForestClassifier(max_features=1, n_estimators=300)
mean accuracy: 0.7669856459330144
RandomForestClassifier(max_features=2, n_estimators=300)
mean accuracy: 0.7643882433356117
RandomForestClassifier(max_features=3, n_estimators=300)
mean accuracy : 0.7643540669856461
RandomForestClassifier(max_features=4, n_estimators=300)
mean accuracy: 0.7526144907723855
```

```
RandomForestClassifier(max_features=5, n_estimators=300)
mean accuracy: 0.7448051948051949
n estimators: 400
RandomForestClassifier(max_features=1, n_estimators=400)
mean accuracy : 0.7591934381408065
RandomForestClassifier(max features=2, n estimators=400)
mean accuracy: 0.7630553656869445
RandomForestClassifier(max_features=3, n_estimators=400)
mean accuracy: 0.755228981544771
RandomForestClassifier(max_features=4, n_estimators=400)
mean accuracy: 0.7487012987012986
RandomForestClassifier(max_features=5, n_estimators=400)
mean accuracy: 0.7577922077922079
n estimators : 500
RandomForestClassifier(max features=1, n estimators=500)
mean accuracy: 0.7591934381408065
RandomForestClassifier(max_features=2, n_estimators=500)
mean accuracy : 0.760475051264525
RandomForestClassifier(max_features=3, n_estimators=500)
mean accuracy: 0.7604408749145593
RandomForestClassifier(max features=4, n estimators=500)
mean accuracy: 0.7473855092276145
RandomForestClassifier(max_features=5, n_estimators=500)
mean accuracy: 0.7473855092276145
Out[10]:
  n_estimators max_features mean_accuracy
           200
                           2
                                    0.76823
```

n_estimators: 200 max_features: 2

의 hyper parametr가 가장 성능이 좋다.