HW10

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1

```
trom sklearn.model selection import train_test_split
from sklearn.metrics import confusion matrix
from sklearn.preprocessing import StandardScaler
from sklearn.preprocessing import OneHotEncoder
from sklearn.ensemble import RandomForestClassifier
# reading from the file
name=["class","cap-shape","cap-surface","cap-color","bruises?","odor","gill-attachment","gill-spacing","gill-size","gill-color","stalk-shape"
,"stalk-root", "stalk-surface-above-ring", "stalk-surface-below-ring", "stalk-color-above-ring", "stalk-color-below-ring", "veil-type", "veil-color"
, "ring-number", "ring-type", "spore-print-color", "population", "habitat"]
data = pd.read_csv('agaricus-lepiota.data',names=name)
feature=data[["cap-shape","cap-surface","cap-color","bruises?","odor","gill-attachment","gill-spacing","gill-size","gill-color","stalk-shape"
,"stalk-root", "stalk-surface-above-ring", "stalk-surface-below-ring", "stalk-color-above-ring", "stalk-color-below-ring", "veil-type", "veil-color"
, "ring-number", "ring-type", "spore-print-color", "population", "habitat"]]
rf_classifier = RandomForestClassifier(
                      min_samples_leaf=50,
                      n estimators=150,
                      bootstrap=True,
                      oob score=True,
                      n jobs=-1,
                      random state=42,
                      max_features='auto')
b=data["class"].astype("category").cat.codes#since created dummy, each columns was seperated into more columns, we need original class column to m
hot=pd.get_dummies(data)
print(hot)
b=pd.concat([hot, b.rename("class")], axis=1)
b= b.drop(["class_e","class_p"], axis = 1)#drop the seperated class columns
x_train, x_test, y_train, y_test = train_test_split(
            b, b["class"], test_size=0.2, random_state=42)
print(y train)
print(x train)
rf_classifier.fit(x_train,y_train)
predictions = rf classifier.predict(x test)
print(confusion matrix(y test,predictions))
```

Figure 1: q1

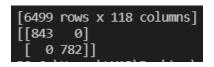


Figure 2: q1

The probability that people get poisoned is nearly 0.

 $\mathbf{2}$

2.1

```
mushroom.py
                 eeg.py 9+ X 💎 homework4.py
                                                     haberman.py
                                                                       🖶 haberman.data
🕏 eeg.py > ...
      11 OH 2 PYTEGI H. HIGGI TC2 THIBOLC COLLUSTOH HIGGI TY
      from sklearn.preprocessing import StandardScaler
      from sklearn.preprocessing import OneHotEncoder
      from sklearn.ensemble import RandomForestClassifier
      from scipy.io import arff
      from sklearn.tree import DecisionTreeClassifier
      from sklearn import metrics
      from sklearn.linear model import LogisticRegression
      from sklearn.model selection import cross val predict
      from sklearn.svm import LinearSVC
12
      data = arff.loadarff('EEG Eye State.arff')
      df = pd.DataFrame(data[0])
      b=df["eyeDetection"].astype("category").cat.codes
      b=pd.concat([df, b.rename("class")], axis=1)
      print(b)
      b= b.drop(["eyeDetection"], axis = 1)#drop the seperated class columns
      p=b.copy()
      p=p.drop(["class"], axis = 1)
      #b=pd.concat([df, b.rename("class")], axis=1)
      print(p)
      x_train, x_test, y_train, y_test = train test split(
                  p, b["class"], test size=0.2, random state=42)
      #random forest
      rf = RandomForestClassifier(
28
                            max depth=50,
                            n estimators=100)
      rf.fit(x train, y train)
      predictions = rf.predict(x test)
      print(confusion matrix(y test,predictions))
      print(metrics.accuracy_score(y_test, predictions))
      #decision tree
      clf = DecisionTreeClassifier(max depth=50)
      clf.fit(x train, y train)
      decision tree predictions = clf.predict(x test)
      print(confusion matrix(y test,decision tree predictions), "decision tree")
      print(metrics.accuracy_score(y_test, decision_tree_predictions))
```

Figure 3: Coding

```
[14980 rows x 14 columns]

[[1519 67]

[ 160 1250]]

0.9242323097463284

[[1353 233]

[ 258 1152]] decision tree

0.8361148197596796
```

Figure 4: Decision Tree confusion matrix is the second one. Random forest is the top one

2.2

Random forest uses bagging to randomly select subset features with replacement instead of choosing the specific features in fixed subsets. This bagging technique creates more diversity. Thus, it is less possible to become overfitting and genealize the dataset.

2.3

By decrease the max depth size of the random forest to 2, I am able to produce a better accuracy result for decision tree than random forest.

```
eeg.py 9+ X
                                                                       mushroom.py
                                 homework4.py
                                                    haberman.py
🕏 eeg.py > ...
      TLONE SKTEMINIMENTES THINDLE CONTRISTON MUCH TY
      from sklearn.preprocessing import StandardScaler
      from sklearn.preprocessing import OneHotEncoder
      from sklearn.ensemble import RandomForestClassifier
      from scipy.io import arff
      from sklearn.tree import DecisionTreeClassifier
      from sklearn import metrics
      from sklearn.linear model import LogisticRegression
      from sklearn.model selection import cross val predict
      from sklearn.svm import LinearSVC
      data = arff.loadarff('EEG Eye State.arff')
      df = pd.DataFrame(data[0])
      b=df["eyeDetection"].astype("category").cat.codes
      b=pd.concat([df, b.rename("class")], axis=1)
      print(b)
      b= b.drop(["eyeDetection"], axis = 1)#drop the seperated class columns
      p=b.copy()
      p=p.drop(["class"], axis = 1)
      #b=pd.concat([df, b.rename("class")], axis=1)
      print(p)
      x train, x test, y train, y test = train test split(
                   p, b["class"], test size=0.2, random state=42)
      #random forest
      rf = RandomForestClassifier(
 28
                            max depth=2,
                             n estimators=100)
      rf.fit(x train, y train)
      predictions = rf.predict(x test)
      print(confusion matrix(y test,predictions))
      print(metrics.accuracy score(y test, predictions))
      #decision tree
      clf = DecisionTreeClassifier(max depth=50)
      clf.fit(x_train, y_train)
      decision tree predictions = clf.predict(x test)
      print(confusion matrix(y test,decision tree predictions), "decision tree")
      print(metrics.accuracy score(y test, decision tree predictions))
```

Figure 5: Coding

```
[14980 rows x 14 columns]

[[1442 144]

[ 980 430]]

0.6248331108144193

[[1355 231]

[ 256 1154]] decision tree

0.8374499332443258
```

Figure 6: Decision Tree confusion matrix is the second one. Random forest is the top one

3

3.1

Both classifier failed to produce good result since the features provided might not have strong correlation with the final classification. Or perhaps the sample data does not represent the true population survival proportion. Or perhaps the sample size is too small since there are only 306 instances

3.2

It is impossible to say that one is always outperforming the other classifier. For example, decision tree performs better when dataset is small and there are not many features to select from. As a result, making decision based on all possible features is helpful than selecting features from sampled features. However, if the dataset is large, sampling the features before making the decision will prevent overfitting, which will boost the accuracy more than decision tree.

```
mushroom.py 7
                    eeg.py 9+
                                   homework4.py
                                                       haberman.py 8 X

    habern

haberman.py > ...
       import numpy as np
       import pandas as pd
       from sklearn.model selection import train test split
       from sklearn.metrics import confusion matrix
       from sklearn.preprocessing import StandardScaler
       from sklearn.preprocessing import OneHotEncoder
       from sklearn.ensemble import RandomForestClassifier
       from sklearn.tree import DecisionTreeClassifier
       name=["age","year","Number of positive axillary nodes", "class"]
 11
       data = pd.read_csv('haberman.data',names=name)
 12
       feature=data[["age", "year", "Number of positive axillary nodes"]]
       rf classifier = RandomForestClassifier(
                             n estimators=100,
                             max_depth=50)
 16
 17
       x train, x test, y train, y test = train test split(
                   feature, data["class"], test size=0.2, random state=42)
       print(y train)
       print(x train)
       rf classifier.fit(x train,y train)
       predictions = rf classifier.predict(x test)
       print(confusion matrix(y test,predictions),"rf")
       #decision tree
       clf = DecisionTreeClassifier(max depth=50, random state=42)
       clf.fit(x train, y train)
       decision tree predictions = clf.predict(x test)
       print(confusion matrix(y test,decision tree predictions), "decision tree")
 PROBLEMS 27
               OUTPUT
                       DEBUG CONSOLE
                                      TERMINAL
[[39 5]
 [14 4]] rf
 [[35 9]
  [13 5]] decision tree
```

Figure 7: Code

```
#linear logistic regression
logis=LogisticRegression(penalty='12',solver='lbfgs',multi class='multinomial')
logis prediction = cross_val_predict(logis, p, b["class"], cv=5)
print(confusion_matrix(b["class"],logis_prediction),"logistic regression_cv")
print(metrics.accuracy score(b["class"], logis prediction),"cv")
logis o=LogisticRegression(penalty='12',solver='lbfgs',multi class='multinomial')
logis o.fit(x train, y train)
logis pre=logis o.predict(x test)
print(confusion_matrix(y_test,logis_pre),"train_test_split")
print(metrics.accuracy_score(y_test, logis_pre),"train test split")
#linearSVC
linear=LinearSVC(max iter=2000)
1 prediction = cross val predict(linear, p, b["class"], cv=5)
print(confusion matrix(b["class"],l prediction),"linear svc cv")
print(metrics.accuracy_score(b["class"], l_prediction))
linear o=LinearSVC(max iter=2000)
linear o.fit(x train, y train)
1 pre=linear o.predict(x test)
print(confusion_matrix(y_test,l_pre),"train_test_split")
print(metrics.accuracy_score(y_test, l_pre), "train_test_split")
```

Figure 8: Coding

4

```
[[4410 3847]
[3937 2786]] linear svc_cv

0.4803738317757009

C:\Users\LXAB\anaconda3\envs\mltorch\l:
warnings.warn("Liblinear failed to co
[1242 344]
[770 640]] train_test_split
0.6281708945260347 train test split
0.6158210947930574 train test split
```

Based on this confusion matrix, train test split is a better solution. And train test split logistic regression is the best model. The method that I used will directly output the best confusion matrix and its accuracy, so no need to manually select the best one. The confusion matrix shown is the best solution for each classifier.