COMP5318 Assignment 1: Classification

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In [1]:

```
# Import all libraries
from sklearn.impute import SimpleImputer
from sklearn.preprocessing import MinMaxScaler
import pandas as pd
import numpy as np
from sklearn.model selection import StratifiedKFold
from sklearn.model selection import cross val score
from sklearn.neighbors import KNeighborsClassifier
from sklearn.linear model import LogisticRegression
from sklearn.naive bayes import GaussianNB
from sklearn.tree import DecisionTreeClassifier
from sklearn.ensemble import BaggingClassifier
from sklearn.ensemble import AdaBoostClassifier
from sklearn.ensemble import GradientBoostingClassifier
from sklearn.model selection import GridSearchCV
from sklearn.svm import SVC
from sklearn.ensemble import RandomForestClassifier
```

In [2]:

```
# Load Breast Cancer Wisconsin Dataset
# Make sure the file is in the same root directory as 'breast-cancer-wisconsin.csv'
db = pd.read_csv('breast-cancer-wisconsin.csv')
```

In [3]:

```
#Because SimpleImputer requires the missing and replacement values to be in the same
#But '? is a 'string', and the average of each column is a 'float'
#So need to replace '?' to np.nan at first
db.replace('?', np.nan, inplace=True)
#Changing the class values
db['class'].replace("class1",0,inplace=True)
db['class'].replace("class2",1,inplace=True)
#Using SimpleImputer to replace missing values
imp = SimpleImputer(missing_values=np.nan,strategy='mean')
db = imp.fit_transform(db)
#Using MinMaxScaler to Normalise the data between (0,1)
scaler = MinMaxScaler(feature range=(0, 1))
db = scaler.fit_transform(db)
#Setting X as features, y as classes
df = pd.DataFrame(db) #Convert db into dataframe form to facilitate data processing
X = df.drop(df.columns[-1], axis=1).values #Delete the last column to get the data s
y = df[df.columns[-1]].values #Get the last column as a target set
```

In [4]:

```
# Print first ten rows of pre-processed dataset to 4 decimal places
# Iterate through the first 10 items of the array.
# Use the Map and List functions to convert the array to a list.
# And then use the Join function to join the data and output.
for i in range(0,10):
    print(','.join(list(map('{:.4f}'.format,db[i]))))
```

Part 1: Cross validation without parameter tuning

In [5]:

```
## Setting the 10 fold stratified cross-validation
# The stratified folds from cvKFold should be provided to the classifiers
cvKFold=StratifiedKFold(n_splits=10, shuffle=True, random_state=0)
```

In [6]:

```
# K-Nearest Neighbour
def kNNClassifier(X, y, k):
    # Create the KNN classifier, and import the parameter k as the number of neighbor
knn = KNeighborsClassifier(n_neighbors=k)
    # Cross-validation is performed using the cross_val_score function.
    # Set the classifier parameters to KNN.
    # Set the dataset to the processed dataset X and the target set y.
    # Set the number of folds to cvKFold as required.
    scores = cross_val_score(knn, X, y, cv=cvKFold)
    return scores.mean()
```

In [7]:

```
# Logistic Regression
def logregClassifier(X, y):
    # Create the Logistic regression classifier
    logreg = LogisticRegression()
    # Cross-validation is performed using the cross_val_score function.
    # Set the classifier parameters to Logistic Regression.
    # Set the dataset to the processed dataset X and the target set y.
    # Set the number of folds to cvKFold as required.
    scores = cross_val_score(logreg, X, y, cv=cvKFold)
    return scores.mean()
```

In [8]:

```
#Naïve Bayes
def nbClassifier(X, y):
    # Create the Naïve Bayes classifier
    nb = GaussianNB()
    # Cross-validation is performed using the cross_val_score function.
    # Set the classifier parameters to Naïve Bayes.
    # Set the dataset to the processed dataset X and the target set y.
    # Set the number of folds to cvKFold as required.
    scores = cross_val_score(nb, X, y, cv=cvKFold)
    return scores.mean()
```

In [9]:

```
# Decision Tree
def dtClassifier(X, y):
    # Create the Decision Tree classifier
    dt = DecisionTreeClassifier()
    # Cross-validation is performed using the cross_val_score function.
    # Set the classifier parameters to Decision Tree.
    # Set the dataset to the processed dataset X and the target set y.
    # Set the number of folds to cvKFold as required.
    scores = cross_val_score(dt, X, y, cv=cvKFold)
    return scores.mean()
```

In [10]:

```
# Ensembles: Bagging, Ada Boost and Gradient Boosting
def bagDTClassifier(X, y, n estimators, max samples, max depth):
    # Create the Bagging classifier, and import parameters as required.
    bag clf = BaggingClassifier(
        DecisionTreeClassifier(max depth=max depth), n estimators=n estimators, max
    # Cross-validation is performed using the cross val score function.
    # Set the classifier parameters to Bagging.
    # Set the dataset to the processed dataset X and the target set y.
    # Set the number of folds to cvKFold as required.
    scores = cross val score(bag clf, X, y, cv=cvKFold)
    return scores.mean()
def adaDTClassifier(X, y, n_estimators, learning_rate, max_depth):
    # Create the Ada Boost classifier, and import parameters as required.
    ada clf = AdaBoostClassifier(
        DecisionTreeClassifier(max depth=max_depth), n_estimators=n_estimators, lear
    # Cross-validation is performed using the cross val score function.
    # Set the classifier parameters to Ada Boost.
    # Set the dataset to the processed dataset X and the target set y.
    # Set the number of folds to cvKFold as required.
    scores = cross val score(ada clf, X, y, cv=cvKFold)
    return scores.mean()
def gbClassifier(X, y, n_estimators, learning_rate):
    # Create the Gradient Boosting classifier, and import parameters as required.
    gb clf = GradientBoostingClassifier(n estimators=n estimators, learning rate=lea
    # Cross-validation is performed using the cross val score function.
    # Set the classifier parameters to Gradient Boosting.
    # Set the dataset to the processed dataset X and the target set y.
    # Set the number of folds to cvKFold as required.
    scores = cross val score(qb clf, X, y, cv=cvKFold)
    return scores.mean()
```

Part 1 Results

In [11]:

```
# Parameters for Part 1:
#KNN
k=3
#Bagging
bag n estimators = 50
bag max samples = 100
bag max depth = 5
#AdaBoost
ada n estimators = 50
ada_learning_rate = 0.5
ada bag max depth = 5
#GB
qb n estimators = 50
gb learning rate = 0.5
# Print results for each classifier in part 1 to 4 decimal places here:
print("kNN average cross-validation accuracy: {:.4f}".format(kNNClassifier(X, y, k))
print("LR average cross-validation accuracy: {:.4f}".format(logregClassifier(X, y)))
print("NB average cross-validation accuracy: {:.4f}".format(nbClassifier(X, y)))
print("DT average cross-validation accuracy: {:.4f}".format(dtClassifier(X, y)))
print("Bagging average cross-validation accuracy: {:.4f}".format(bagDTClassifier(X,
print("AdaBoost average cross-validation accuracy: {:.4f}".format(adaDTClassifier(X,
print("GB average cross-validation accuracy: {:.4f}".format(gbClassifier(X, y, gb n
```

```
kNN average cross-validation accuracy: 0.9642
LR average cross-validation accuracy: 0.9642
NB average cross-validation accuracy: 0.9585
DT average cross-validation accuracy: 0.9356
Bagging average cross-validation accuracy: 0.9642
AdaBoost average cross-validation accuracy: 0.9585
GB average cross-validation accuracy: 0.9614
```

Part 2: Cross validation with parameter tuning

In [12]:

```
# Linear SVM
# You should use SVC from sklearn.svm
C = \{0.001, 0.01, 0.1, 1, 10, 100\}
gamma = \{0.001, 0.01, 0.1, 1, 10, 100\}
def bestLinClassifier(X,y):
    # Iterate through the parameters C and gamma, converting them to the form of a I
    C list=[]
    for value in C:
        C list.append(value)
    gamma list=[]
    for value in gamma:
        gamma list.append(value)
    # Wrap the parameters C and gamma into a new dictionary.
    para grid = {'C':C list, 'gamma':gamma list}
    # Use the spilt function to get the training set and test set in cvKFold.
    for train index, test index in cvKFold.split(X, y):
        X train, X test = X[train index], X[test index]
        y train, y test = y[train index], y[test index]
    # Create the SVM classifier, and set the kernel method to linear.
    lin svm = SVC(kernel="linear")
    # Set estimator for grid search.
    grid_search = GridSearchCV(lin_svm, para_grid, cv=cvKFold, return train score=Tr
    # Import the training set for training.
    grid_search.fit(X_train, y_train)
    # Get the desired results.
    best C = grid search.best params ['C']
    best gamma = grid search.best params ['gamma']
    best val score = grid search.best score
    best set score = grid search.score(X test, y test)
    return best_C, best_gamma, best_val_score, best_set_score #(appropriate values s
```

In [13]:

```
# Random Forest
# You should use RandomForestClassifier from sklearn.ensemble with information gain
n = \{10, 20, 30, 50, 100\}
max leaf nodes = \{4, 10, 16, 20, 30\}
def bestRFClassifier(X,y):
    # Iterate through the parameters n estimators and max leaf nodes, converting the
    n estimators list=[]
    for value in n estimators:
        n estimators list.append(value)
    max leaf nodes list=[]
    for value in max leaf nodes:
        max leaf nodes list.append(value)
    # Wrap the parameters n estimators and max leaf nodes into a new dictionary.
    para_grid = {'n_estimators':n_estimators_list,'max leaf nodes':max leaf nodes li
    # Use the spilt function to get the training set and test set in cvKFold.
    for train index, test index in cvKFold.split(X, y):
        X train, X test = X[train index], X[test index]
        y train, y test = y[train index], y[test index]
    # Create the Random Forest classifier, and set the criterion as entropy and max
    rnd clf = RandomForestClassifier(criterion='entropy', max features='sqrt')
    # Set estimator for grid search.
    grid search = GridSearchCV(rnd clf, para grid, cv=cvKFold, return train score=Tx
    # Import the training set for training.
    grid search.fit(X train, y train)
    # Get the desired results.
    best n estimators = grid search.best params ['n estimators']
    best max leaf nodes = grid search.best params ['max leaf nodes']
    best val score = grid search.best score
    best set score = grid search.score(X test, y test)
    return best_n_estimators, best_max_leaf_nodes, best_val_score, best_set_score#(a
```

Part 2 Results

In [14]:

```
# Perform Grid Search with 10-fold Stratified Cross Validation (GridSearchCV in skle
# The stratified folds from cvKFold should be provided to GridSearchV

# This should include using train_test_split from sklearn.model_selection with strat
# Print results for each classifier here. All results should be printed to 4 decimal
# "n_estimators" and "max_leaf_nodes" which should be printed as integers.

best_C, best_gamma, best_val_score, best_set_score = bestLinClassifier(X,y)
print("SVM best C: {:.4f}".format(best_C))
print("SVM best gamma: {:.4f}".format(best_gamma))
print("SVM cross-validation accuracy: {:.4f}".format(best_val_score))

best_n_estimators, best_max_leaf_nodes, best_val_score, best_set_score = bestRFClass
print("RF best n_estimators: {:.4f}".format(best_n_estimators))
print("RF best max_leaf_nodes: {:.4f}".format(best_max_leaf_nodes))
print("RF cross-validation accuracy: {:.4f}".format(best_val_score))
print("RF test set accuracy: {:.4f}".format(best_val_score))
```

```
SVM best C: 1.0000

SVM best gamma: 0.1000

SVM cross-validation accuracy: 0.9683

SVM test set accuracy: 0.9710

RF best n_estimators: 20.0000

RF best max_leaf_nodes: 16.0000

RF cross-validation accuracy: 0.9714

RF test set accuracy: 0.9565
```