FinalProject

December 15, 2024

```
[1]: # All necessary packages needed
     import pandas as pd
     import numpy as np
     import matplotlib
     import seaborn as sns
     from matplotlib import pyplot as plt
     from sklearn.model_selection import train_test_split, KFold
     from sklearn.preprocessing import OneHotEncoder, OrdinalEncoder, MinMaxScaler,

→StandardScaler

     from xgboost import XGBClassifier
     from sklearn.metrics import accuracy_score, confusion_matrix,_
      →ConfusionMatrixDisplay
     from sklearn.model_selection import ParameterGrid
     from sklearn.experimental import enable_iterative_imputer
     from sklearn.impute import IterativeImputer
     from sklearn.ensemble import RandomForestRegressor, RandomForestClassifier
     from sklearn.svm import SVC
     from sklearn.linear_model import LogisticRegression
     from sklearn.neighbors import KNeighborsClassifier
     import warnings
     from sklearn.exceptions import ConvergenceWarning
     import shap
[2]: # import data
     df = pd.read csv('data/kidney disease.csv')
     df.head()
     # EDA - print rows and columns
     print("Rows:", df.shape[0], "Columns:", df.shape[1])
```

Rows: 400 Columns: 26

0.1 Exporatory Data Analysis (EDA)

0.1.1 (i) About the whole data structure

```
Non-Null Count
         Column
                                         Dtype
                         -----
     0
         id
                         400 non-null
                                         int64
     1
         age
                         391 non-null
                                         float64
     2
                         388 non-null
                                         float64
         bр
     3
                         353 non-null
                                         float64
         sg
     4
         al
                         354 non-null
                                         float64
     5
                         351 non-null
                                         float64
         su
     6
         rbc
                         248 non-null
                                         object
     7
                         335 non-null
                                         object
         рс
     8
                         396 non-null
         рсс
                                         object
     9
                         396 non-null
                                         object
         ba
                         356 non-null
                                         float64
     10
         bgr
     11
         bu
                         381 non-null
                                         float64
     12
                         383 non-null
                                         float64
         sc
     13
         sod
                         313 non-null
                                         float64
     14
                         312 non-null
                                         float64
        pot
                         348 non-null
                                         float64
     15
         hemo
     16
                         330 non-null
                                         object
         pcv
     17
         WC
                         295 non-null
                                         object
                         270 non-null
     18
        rc
                                         object
     19
        htn
                         398 non-null
                                         object
     20
                         398 non-null
         dm
                                         object
     21
        cad
                         398 non-null
                                         object
     22
                         399 non-null
                                         object
         appet
                         399 non-null
     23
         ре
                                         object
     24
         ane
                         399 non-null
                                         object
     25 classification 400 non-null
                                         object
    dtypes: float64(11), int64(1), object(14)
    memory usage: 81.4+ KB
    None
[4]: # EDA - target variable type
     print("The column Y is discrete")
     print(df['classification'].value_counts())
    The column Y is discrete
    classification
    ckd
              248
              150
    notckd
    ckd\t
    Name: count, dtype: int64
[5]: df['classification'] = df['classification'].replace(to_replace={'ckd\t':'ckd',__
      print(df['classification'].value_counts())
    classification
```

Data columns (total 26 columns):

```
ckd 250
not ckd 150
Name: count, dtype: int64
```

Hence we don't need stratified spliting. We will use basic split for train, val and test sets.

```
rbc has [nan 'normal' 'abnormal']

pc has ['normal' 'abnormal' nan]

pcc has ['notpresent' 'present' nan]

ba has ['notpresent' 'present' nan]

htn has ['yes' 'no' nan]

dm has ['yes' 'no' 'yes' '\tno' '\tyes' nan]

cad has ['no' 'yes' nan]

ane has ['no' 'yes' nan]

appet has ['good' 'poor' nan]

sg has [1.02 1.01 1.005 1.015 nan 1.025]

al has [ 1. 4. 2. 3. 0. nan 5.]

su has [ 0. 3. 4. 1. nan 2. 5.]
```

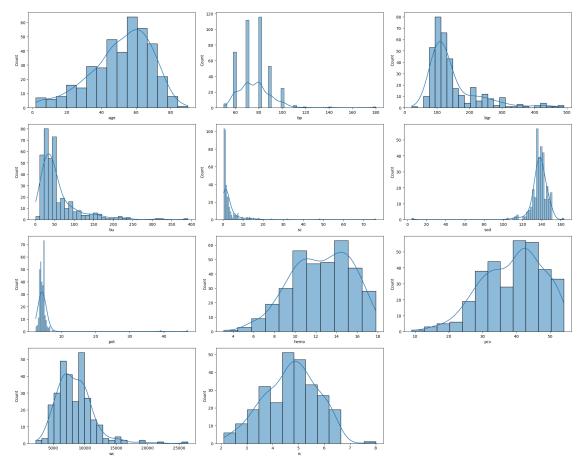
0.1.2 (ii) Then we first take a look at the continuous variables

```
[7]: # for the column pcv, wc, rc, there are some strings ? and some values with tabuahead
for column in ['pcv', 'wc', 'rc']:
    df[column] = df[column].str.strip() # Remove any leading/trailinguahhitespace
    df[column] = df[column].replace('?',np.nan)

plt.figure(figsize=(20, 16))
plotnumber = 1

for column in num_cols:
    # Convert to numeric
    df[column] = pd.to_numeric(df[column])

ax = plt.subplot(4, 3, plotnumber)
```

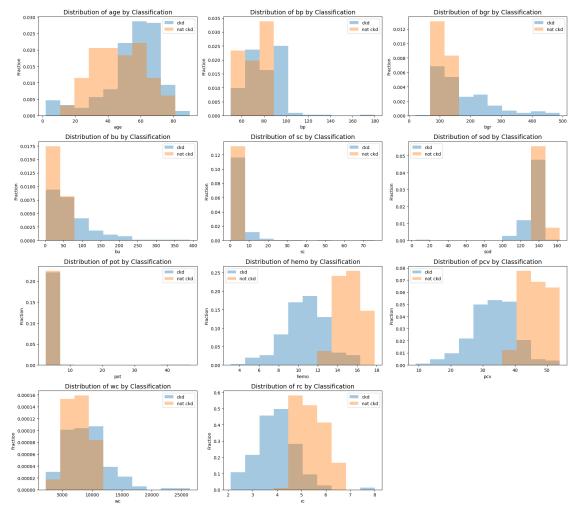


```
[8]: # compare the classification of continous variables
plt.figure(figsize=(18, 16))
categories = df['classification'].unique()

for i, column in enumerate(num_cols):
    plt.subplot(4,3,i+1) # Create subplots
    bin_range = (df[column].min(), df[column].max()) # Set the bin range forusthe current column
```

```
for c in categories:
    plt.hist(df[df['classification'] == c][column], alpha=0.4, label=c,_u
    range = bin_range, density=True)
    plt.legend()
    plt.ylabel('Fraction')
    plt.xlabel(column)
    plt.title(f'Distribution of {column} by Classification', fontsize = 14)

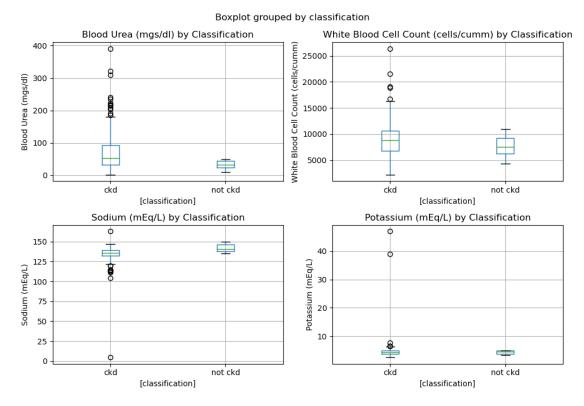
plt.tight_layout()
plt.savefig('/Users/ericdavis/Desktop/data1030/FinalProject1030/cont_hist.png',_u
    dpi = 350)
plt.show()
```



Now we find that age, bp, bgr, hemo, pcv, wc, rc may have large influence on classification(has disease or not). And we are interested in using bar plot to see if any columns have too many

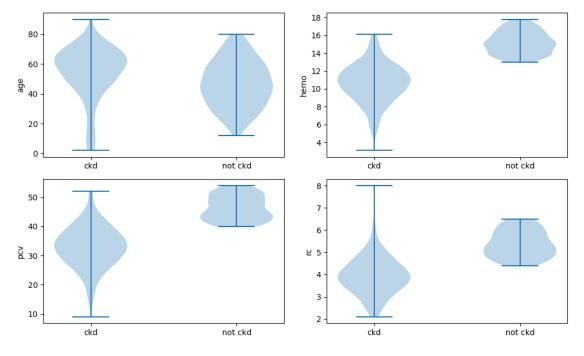
extreme outliers. From above plot, I will choose bgr, bu, we to investigate.

```
[9]: plt.figure(figsize = (10,7))
  bar_cols = ['bu', 'wc', 'sod', 'pot']
  bar_cols_full = ['Blood Urea (mgs/dl)', 'White Blood Cell Count (cells/cumm)', \( \text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\text{\tex{
```



Next, I am going to visualize some distribution of some attributes across different classidication by violin plot. Here I choose age, hemo, pcv and rc since they have large gap between two distributions from

```
[10]: # violin plot
plt.figure(figsize = (10,6))
```



```
[11]: # plot the same attributes using histogram

plt.figure(figsize=(12, 9))

categories = df['classification'].unique()

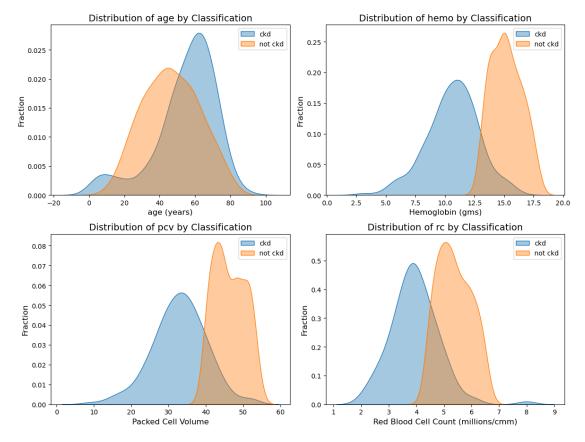
full_name_cols = ['age (years)', 'Hemoglobin (gms)', 'Packed Cell Volume', 'Red_

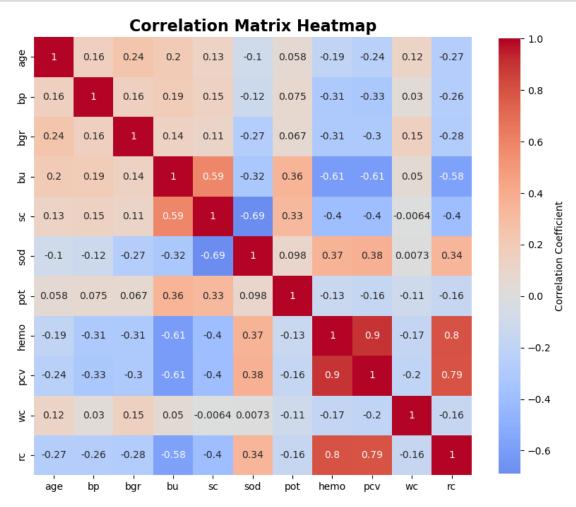
Blood Cell Count (millions/cmm)']

for i, column in enumerate(vio_cols):

plt.subplot(2,2,i+1) # Create subplots

for c in categories:
```





we find hemo and pcv have strong correlation relationship, hence we choose to drop one of these attributes – hemo.

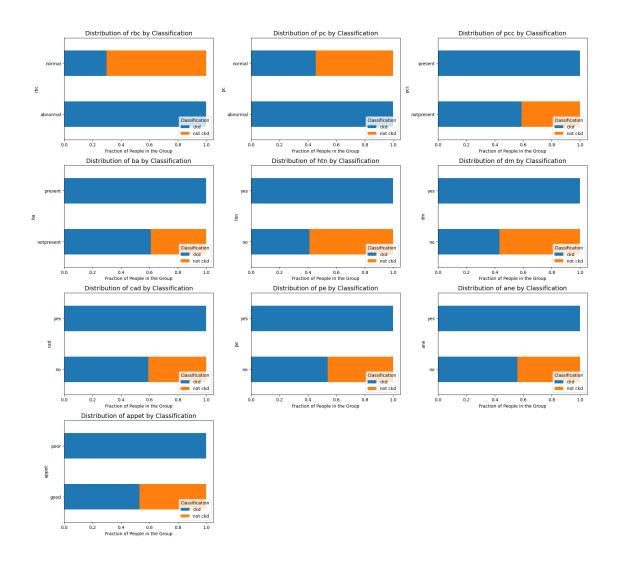
```
[13]: df = df.drop(columns = 'hemo')
num_cols.remove('hemo')
```

0.1.3 (iii) Next we take a look at the categorical variables with no order

```
[14]: plt.figure(figsize=(18, 20))
    categories = df['classification'].unique()
    cat_cols = ['rbc', 'pc', 'pcc', 'ba', 'htn', 'dm', 'cad', 'pe', 'ane', 'appet']
    for column in cat_cols:
```

```
df[column] = df[column].str.strip() # Remove any leading/trailing_
 \hookrightarrow whitespace
for i, column in enumerate(cat_cols):
    plt.subplot(5, 3, i + 1) # Use dynamic column count
    count_matrix = df.groupby([df[column], df['classification']]).size().

unstack()
    count_matrix_norm = count_matrix.div(count_matrix.sum(axis=1), axis=0)
    count_matrix_norm.plot(kind='barh', stacked=True, ax = plt.gca()) # Using_
 ⇔the current axis
    plt.legend(title='Classification', fontsize=10, loc='lower right')
    plt.xlabel('Fraction of People in the Group')
    plt.title(f'Distribution of {column} by Classification', fontsize = 14)
plt.tight_layout()
plt.savefig('/Users/ericdavis/Desktop/data1030/FinalProject1030/cat.png', dpi =__
 ⇒350)
plt.show()
```



We can find that all categorical attributes are binary. Also from the above stacked bar plot, we wonder if all not ckd happens in only one of the situation in all the binary attributes.

```
pd.DataFrame(results, index = ['notckdS1', 'ckdS1', 'notckdS2', 'ckdS2'])
```

[15]: appet rbc рс рсс ba htn \mathtt{dm} cad ре ane 0.47 notckdS1 0.701 0.544 0.412 0.39 0.00 0.000 0.407 0.461 0.44 ckdS1 0.299 0.456 0.588 0.61 1.00 1.000 0.593 0.539 0.56 0.53 notckdS2 0.000 0.000 0.000 0.00 0.59 0.567 0.000 0.000 0.00 0.00 ckdS2 1.000 1.000 1.000 1.00 0.41 0.433 1.000 1.000 1.00 1.00

0.1.4 (iv) Do some summary statistics

[16]: df.shape

[16]: (400, 25)

Before preprocessing, we have 25 features and 400 data points(One feature is dropped because of colinearity)

[17]: df.describe()

[17]:		id	age	bp	sg	al	su	\
	count	400.000000	391.000000	388.000000	353.000000	354.000000	351.000000	
	mean	199.500000	51.483376	76.469072	1.017408	1.016949	0.450142	
	std	115.614301	17.169714	13.683637	0.005717	1.352679	1.099191	
	min	0.000000	2.000000	50.000000	1.005000	0.000000	0.000000	
	25%	99.750000	42.000000	70.000000	1.010000	0.000000	0.000000	
	50%	199.500000	55.000000	80.000000	1.020000	0.000000	0.000000	
	75%	299.250000	64.500000	80.000000	1.020000	2.000000	0.000000	
	max	399.000000	90.000000	180.000000	1.025000	5.000000	5.000000	
		bgr	bu	sc	sod	pot	pcv	\
	count	356.000000	381.000000	383.000000	313.000000	312.000000	329.000000	
	mean	148.036517	57.425722	3.072454	137.528754	4.627244	38.884498	
	std	79.281714	50.503006	5.741126	10.408752	3.193904	8.990105	
	min	22.000000	1.500000	0.400000	4.500000	2.500000	9.000000	
	25%	99.000000	27.000000	0.900000	135.000000	3.800000	32.000000	
	50%	121.000000	42.000000	1.300000	138.000000	4.400000	40.000000	
	75%	163.000000	66.000000	2.800000	142.000000	4.900000	45.000000	
	max	490.000000	391.000000	76.000000	163.000000	47.000000	54.000000	

rc WC count 294.000000 269.000000 8406.122449 4.707435 meanstd 2944.474190 1.025323 2.100000 min 2200.000000 25% 6500.000000 3.900000 50% 8000.00000 4.800000 75% 9800.000000 5.400000 max 26400.000000 8.000000

0.2 Data Cleaning & Preparation

```
[18]: for i in cat_cols:
          print(df[i].unique())
     [nan 'normal' 'abnormal']
     ['normal' 'abnormal' nan]
     ['notpresent' 'present' nan]
     ['notpresent' 'present' nan]
     ['yes' 'no' nan]
     ['yes' 'no' nan]
     ['no' 'yes' nan]
     ['no' 'yes' nan]
     ['no' 'yes' nan]
     ['good' 'poor' nan]
[19]: # replace nan with string (another category)
      for i in cat_cols:
          df.loc[:,i] = df[i].replace(to_replace = {np.nan: 'missing'})
      for i in cat_cols:
          print(df[i].unique())
     ['missing' 'normal' 'abnormal']
     ['normal' 'abnormal' 'missing']
     ['notpresent' 'present' 'missing']
     ['notpresent' 'present' 'missing']
     ['yes' 'no' 'missing']
     ['yes' 'no' 'missing']
     ['no' 'yes' 'missing']
     ['no' 'yes' 'missing']
     ['no' 'yes' 'missing']
     ['good' 'poor' 'missing']
[20]: # transfer the target variable from ['ckd', 'not ckd'] into 0s and 1s (easy for_
       ⇔future analysis)
      warnings.filterwarnings("ignore")
      df['classification'] = pd.DataFrame(df['classification'].
       →replace(to_replace={'ckd': 1, 'not ckd': 0}).reset_index(drop=True))
[21]: # Check the missing values percentage for each attribute
      df.isnull().sum()/len(df)
[21]: id
                        0.0000
      age
                        0.0225
                        0.0300
      bp
```

```
0.1175
sg
                   0.1150
al
su
                   0.1225
                   0.0000
rbc
                   0.0000
рс
                   0.0000
рсс
                   0.0000
ba
                   0.1100
bgr
                   0.0475
bu
                   0.0425
sc
sod
                   0.2175
                   0.2200
pot
                   0.1775
pcv
                   0.2650
WC
                   0.3275
rc
htn
                   0.0000
                   0.0000
dm
                   0.0000
cad
                   0.0000
appet
                   0.0000
ре
                   0.0000
ane
                   0.0000
classification
dtype: float64
```

We can find the target variable doesn't contain missing values. And all attributes contain some missing values. And for the category variables, we can treat the missing values as another category.

0.3 Split and Preprocess the Dataset

```
[22]: # keep random state the same for spliting and training the model np.random.seed(42) # for reproducibility random_states = list(np.random.randint(1, 10000, size = 10))
```

```
[23]: X_train = []
X_val = []
X_test = []
y_train = []
y_val = []
y_test = []
X_test_withoutstd = []

X = df.loc[:, df.columns != 'classification']
y = df['classification']
# Define categorical and numerical features
cat_cols_noorder = ['rbc', 'pc', 'pcc', 'ba', 'htn', 'dm', 'cad', 'pe', 'ane', \underset{\underset}
\u
```

```
ordinal_cats = [[1.005,1.010,1.015,1.020,1.025,np.nan], [0,1,2,3,4,5,np.nan],
 \rightarrow [0,1,2,3,4,5,np.nan]]
for i in range(len(random states)):
    # use basic split to seperate out the test set
    X train unit, X other unit, y train unit, y other unit = 1
 otrain_test_split(X, y, train_size = 0.6, random_state = random_states[i])
    X_val_unit, X_test_unit, y_val_unit, y_test_unit =
 →train_test_split(X_other_unit, y_other_unit, train_size = 0.5, random_state_
 →= random_states[i])
    y_train.append(y_train_unit)
    y_val.append(y_val_unit)
    y_test.append(y_test_unit)
    # Preprocessing for the categorical data with no order -- OneHotEncoder
    # get the column names
    colnames = []
    for k in cat_cols_noorder:
        temp = [f'{k}_{value}' for value in X_train_unit[k].unique()]
        colnames.extend(temp)
    enc = OneHotEncoder(sparse_output = False, handle_unknown = 'ignore')
    X_train_pre_cat1 = pd.DataFrame(enc.
 afit_transform(X_train_unit[cat_cols_noorder]), columns = colnames).
 →reset_index(drop=True)
    X_test_pre_cat1 = pd.DataFrame(enc.
 stransform(X_test_unit[cat_cols_noorder]), columns = colnames).
 →reset_index(drop=True)
    X_val_pre_cat1 = pd.DataFrame(enc.transform(X_val_unit[cat_cols_noorder]),_

¬columns = colnames).reset_index(drop=True)

    # Preprocessing for the categorical data with order -- OrdinalEncoder
    ordinal = OrdinalEncoder(categories = ordinal_cats)
    X_train_pre_cat2 = pd.DataFrame(ordinal.

→fit_transform(X_train_unit[cat_cols_withorder]), columns =
□
 →cat_cols_withorder).reset_index(drop=True)
    X test pre cat2 = pd.DataFrame(ordinal.

¬transform(X_test_unit[cat_cols_withorder]), columns = cat_cols_withorder).

 →reset_index(drop=True)
    X_val_pre_cat2 = pd.DataFrame(ordinal.
 otransform(X_val_unit[cat_cols_withorder]), columns = cat_cols_withorder).
 →reset_index(drop=True)
    # Preprocessing for the continous variables -- StandardScalar
    scaler = StandardScaler()
```

```
X_train_pre_num = pd.DataFrame(scaler.
fit_transform(X_train_unit[num_cols]), columns = num_cols).
→reset_index(drop=True)
  X_test_pre_num = pd.DataFrame(scaler.transform(X_test_unit[num_cols]),_

¬columns = num_cols).reset_index(drop=True)

  X val_pre_num = pd.DataFrame(scaler.transform(X_val_unit[num_cols]),_

¬columns = num_cols).reset_index(drop=True)
  # merge the table
  X_train_merge = pd.concat([X_train_pre_num, X_train_pre_cat2,_
→X_train_pre_cat1], axis = 1)
  X_val_merge = pd.concat([X_val_pre_num, X_val_pre_cat2, X_val_pre_cat1],_
\Rightarrowaxis = 1)
  X_test_merge = pd.concat([X_test_pre_num, X_test_pre_cat2,_
→X_test_pre_cat1], axis = 1)
  # Apply StandardScalar to all the features
  X_train.append(pd.DataFrame(scaler.fit_transform(X_train_merge), columns =

¬X_train_merge.columns))
  X_val.append(pd.DataFrame(scaler.transform(X_val_merge), columns =_
→X val merge.columns))
  X_test.append(pd.DataFrame(scaler.transform(X_test_merge), columns =__
→X_test_merge.columns))
  # concat tables
  X_test_withoutstd.append(X_test_merge)
```

0.4 Choose ML model, Tune hyperparameter and Evaluate

0.4.1 (i) XGBoosting

```
[24]: # Initialze empty variables
best_model_xgb = []
best_valscore_xgb = np.zeros(len(random_states))
test_score_xgb = np.zeros(len(random_states))
y_pred_xgb = []

for i in range(len(random_states)):
    # Define hyperparameter grid for tuning
    param_grid = {
        'max_depth': [1, 3, 5, 8],
        'n_estimators': [10000],
        'learning_rate': [0.03],
        'reg_alpha': [1e-4, 1e-3, 1e-2, 1e-1, 0e0],
        'reg_lambda': [1e-4, 1e-3, 1e-2, 1e-1, 0e0]
}
```

```
val_score = np.zeros(len(ParameterGrid(param_grid)))
    models = []
    # train the model XGBoosting
    for p in range(len(ParameterGrid(param_grid))):
        params = ParameterGrid(param_grid)[p]
        clf = XGBClassifier(**params, early_stopping_rounds = 30, n_jobs = -1,__
  →random_state = random_states[i])
        clf.fit(X_train[i], y_train[i], eval_set = [(X_val[i], y_val[i])],_u
  ⇔verbose = False)
        models.append(clf)
        y_val_pred = clf.predict(X_val[i])
        val_score[p] = accuracy_score(y_val[i], y_val_pred)
    # get the best validation score
    best_valscore_xgb[i] = np.max(val_score)
    best_model_xgb.append(models[np.argmax(val_score)])
    # evaluate the model using test set (test score)
    y_test_pred = best_model_xgb[i].predict(X_test[i])
    test_score_xgb[i] = accuracy_score(y_test[i], y_test_pred)
    y_pred_xgb.append(y_test_pred)
    print('Random state: ', random_states[i])
    print('The best model parameters are: ', ParameterGrid(param_grid)[np.
 →argmax(val_score)])
    print('Best validation score: ', best valscore xgb[i])
    print('Test accuracy: ', test_score_xgb[i])
    print(' ')
Random state: 7271
The best model parameters are: {'reg_lambda': 0.0001, 'reg_alpha': 0.0001,
'n estimators': 10000, 'max depth': 1, 'learning rate': 0.03}
Best validation score: 1.0
Test accuracy: 1.0
Random state: 861
The best model parameters are: {'reg_lambda': 0.0001, 'reg_alpha': 0.0001,
'n_estimators': 10000, 'max_depth': 1, 'learning_rate': 0.03}
Best validation score: 0.9875
Test accuracy: 1.0
Random state: 5391
The best model parameters are: {'reg_lambda': 0.0001, 'reg_alpha': 0.0001,
'n_estimators': 10000, 'max_depth': 1, 'learning_rate': 0.03}
Best validation score: 0.9875
Test accuracy: 0.9875
```

```
Random state: 5192
     The best model parameters are: {'reg_lambda': 0.0001, 'reg_alpha': 0.0001,
     'n_estimators': 10000, 'max_depth': 1, 'learning_rate': 0.03}
     Best validation score: 1.0
     Test accuracy: 1.0
     Random state: 5735
     The best model parameters are: {'reg_lambda': 0.0001, 'reg_alpha': 0.0001,
     'n_estimators': 10000, 'max_depth': 1, 'learning_rate': 0.03}
     Best validation score: 1.0
     Test accuracy: 1.0
     Random state: 6266
     The best model parameters are: {'reg_lambda': 0.0001, 'reg_alpha': 0.0001,
     'n_estimators': 10000, 'max_depth': 1, 'learning_rate': 0.03}
     Best validation score: 1.0
     Test accuracy: 0.9875
     Random state: 467
     The best model parameters are: {'reg_lambda': 0.0001, 'reg_alpha': 0.0001,
     'n_estimators': 10000, 'max_depth': 1, 'learning_rate': 0.03}
     Best validation score: 1.0
     Test accuracy: 0.975
     Random state: 4427
     The best model parameters are: {'reg_lambda': 0.0001, 'reg_alpha': 0.0001,
     'n_estimators': 10000, 'max_depth': 1, 'learning_rate': 0.03}
     Best validation score: 1.0
     Test accuracy: 1.0
     Random state: 5579
     The best model parameters are: {'reg_lambda': 0.0001, 'reg_alpha': 0.0001,
     'n_estimators': 10000, 'max_depth': 1, 'learning_rate': 0.03}
     Best validation score: 1.0
     Test accuracy: 1.0
     Random state: 8323
     The best model parameters are: {'reg_lambda': 0.0001, 'reg_alpha': 0.0001,
     'n_estimators': 10000, 'max_depth': 1, 'learning_rate': 0.03}
     Best validation score: 1.0
     Test accuracy: 1.0
[25]: print('The average test accuracy for xgb is', test_score_xgb.mean())
      print('The test accuracy standard deviation for xgb is', test_score_xgb.std())
      y_test_flatten = [item for sublist in y_test for item in sublist]
```

```
y_pred_flatten = [item for sublist in y_pred_xgb for item in sublist] # get the_

-y_true and y_pred

overall_accuracy_xgb = accuracy_score(y_test_flatten, y_pred_flatten)

print('The OVERALL test accuracy for xgb is', overall_accuracy_xgb)
```

When training with validation data, set a large n_estimators (e.g., 10000) and use early stopping to terminate the training once performance on the validation set stops improving. Here the test score stablized at 0.99, because we set large n_estimators and use early stopping to avoid overfitting. And here other parameters become less important with those two parameters settled down.

0.4.2 (ii) Random Forest

```
[27]: # Initialze empty variables
best_model_rf = []
best_valscore_rf = np.zeros(len(random_states))
test_score_rf = np.zeros(len(random_states))
y_pred_rf = []

for i in range(len(random_states)):
    # Define hyperparameter grid for tuning
    param_grid = {
        'n_estimators': [100],
        'max_depth': [1, 3, 5, 8, 13],
        'max_features': [1, 3, 5, 8, 13, 18]
}

val_score = np.zeros(len(ParameterGrid(param_grid)))
models = []
```

```
# train the model random forest
    for p in range(len(ParameterGrid(param_grid))):
        params = ParameterGrid(param_grid)[p]
        clf = RandomForestClassifier(**params, n_jobs = -1, random_state =__
  →random_states[i])
        clf.fit(X train[i], y train[i])
        models.append(clf)
        y_val_pred = clf.predict(X_val[i])
        val_score[p] = accuracy_score(y_val[i], y_val_pred)
    # get the best validation score
    best_valscore_rf[i] = np.max(val_score)
    best_model_rf.append(models[np.argmax(val_score)])
    # evaluate the model using test set (test score)
    y_test_pred = best_model_rf[i].predict(X_test[i])
    test_score_rf[i] = accuracy_score(y_test[i], y_test_pred)
    y_pred_rf.append(y_test_pred)
    print('Random state: ', random_states[i])
    print('The best model parameters are: ', ParameterGrid(param_grid)[np.
 →argmax(val_score)])
    print('Best validation score: ', best_valscore_rf[i])
    print('Test accuracy: ', test_score_rf[i])
    print(' ')
Random state: 7271
The best model parameters are: {'n_estimators': 100, 'max_features': 3,
'max depth': 1}
Best validation score: 1.0
Test accuracy: 0.95
Random state: 861
The best model parameters are: {'n_estimators': 100, 'max_features': 3,
'max depth': 3}
Best validation score: 1.0
Test accuracy: 0.9875
Random state: 5391
The best model parameters are: {'n_estimators': 100, 'max_features': 13,
'max_depth': 3}
Best validation score: 1.0
Test accuracy: 0.9875
Random state: 5192
The best model parameters are: {'n_estimators': 100, 'max_features': 5,
'max_depth': 1}
```

```
Best validation score: 1.0
     Test accuracy: 0.9375
     Random state: 5735
     The best model parameters are: {'n_estimators': 100, 'max_features': 3,
     'max depth': 3}
     Best validation score: 1.0
     Test accuracy: 1.0
     Random state: 6266
     The best model parameters are: {'n_estimators': 100, 'max_features': 3,
     'max_depth': 1}
     Best validation score: 1.0
     Test accuracy: 0.9625
     Random state: 467
     The best model parameters are: {'n_estimators': 100, 'max_features': 3,
     'max_depth': 1}
     Best validation score: 1.0
     Test accuracy: 0.9625
     Random state: 4427
     The best model parameters are: {'n_estimators': 100, 'max_features': 3,
     'max_depth': 8}
     Best validation score: 1.0
     Test accuracy: 1.0
     Random state: 5579
     The best model parameters are: {'n_estimators': 100, 'max_features': 5,
     'max_depth': 3}
     Best validation score: 1.0
     Test accuracy: 1.0
     Random state: 8323
     The best model parameters are: {'n_estimators': 100, 'max_features': 3,
     'max depth': 8}
     Best validation score: 1.0
     Test accuracy: 1.0
[28]: print('The average test accuracy for rf is', test score rf.mean())
      print('The test accuracy standard deviation for rf is', test_score_rf.std())
      y_pred_flatten = [item for sublist in y_pred_rf for item in sublist]
      overall_accuracy_rf = accuracy_score(y_test_flatten, y_pred_flatten)
      print('The OVERALL test accuracy for rf is', overall_accuracy_rf)
```

The average test accuracy for rf is 0.97875
The test accuracy standard deviation for rf is 0.022395591083961153

0.4.3 (iii) SVM

Random state: 7271

Best validation score: 1.0

```
[29]: # Initialze empty variables
      best model svm = []
      best_valscore_svm = np.zeros(len(random_states))
      test_score_svm = np.zeros(len(random_states))
      y_pred_svm = []
      for i in range(len(random_states)):
          # Define hyperparameter grid for tuning
          param_grid = {
              'C': [1e-1, 1e0, 1e1, 1e2, 1e3, 1e4],
              'gamma': [1e-4, 1e-3, 1e-2, 1e-1, 1e0, 1e1]
          }
          val_score = np.zeros(len(ParameterGrid(param_grid)))
          models = \Pi
          # train the model SVM
          for p in range(len(ParameterGrid(param_grid))):
              params = ParameterGrid(param_grid)[p]
              clf = SVC(**params, probability = True, random_state =
       →random_states[i]) # use SVC for classification problem
              clf.fit(X_train[i], y_train[i])
              models.append(clf)
              y_val_pred = clf.predict(X_val[i])
              val_score(p] = accuracy_score(y_val[i], y_val_pred)
          # get the best validation score
          best_valscore_svm[i] = np.max(val_score)
          best_model_svm.append(models[np.argmax(val_score)])
          # evaluate the model using test set (test score)
          y_test_pred = best_model_svm[i].predict(X_test[i])
          test_score_svm[i] = accuracy_score(y_test[i], y_test_pred)
          y_pred_svm.append(y_test_pred)
          print('Random state: ', random_states[i])
          print('The best model parameters are: ', ParameterGrid(param_grid)[np.
       →argmax(val_score)])
          print('Best validation score: ', best_valscore_svm[i])
          print('Test accuracy: ', test_score_svm[i])
          print(' ')
```

22

The best model parameters are: {'gamma': 0.01, 'C': 1.0}

Test accuracy: 0.9625

Random state: 861

The best model parameters are: {'gamma': 0.1, 'C': 1.0}

Best validation score: 0.9875

Test accuracy: 0.9625

Random state: 5391

The best model parameters are: {'gamma': 0.01, 'C': 1.0}

Best validation score: 1.0

Test accuracy: 0.9875

Random state: 5192

The best model parameters are: {'gamma': 0.01, 'C': 1.0}

Best validation score: 1.0

Test accuracy: 1.0

Random state: 5735

The best model parameters are: {'gamma': 0.01, 'C': 1.0}

Best validation score: 1.0

Test accuracy: 1.0

Random state: 6266

The best model parameters are: {'gamma': 0.01, 'C': 1.0}

Best validation score: 1.0

Test accuracy: 0.975

Random state: 467

The best model parameters are: {'gamma': 0.1, 'C': 1.0}

Best validation score: 1.0

Test accuracy: 1.0

Random state: 4427

The best model parameters are: {'gamma': 0.1, 'C': 0.1}

Best validation score: 1.0

Test accuracy: 0.925

Random state: 5579

The best model parameters are: {'gamma': 0.1, 'C': 1.0}

Best validation score: 1.0

Test accuracy: 0.9875

Random state: 8323

The best model parameters are: {'gamma': 0.01, 'C': 100.0}

Best validation score: 1.0

Test accuracy: 1.0

```
[30]: print('The average test accuracy for svm is', test_score_svm.mean())
print('The test accuracy standard deviation for svm is', test_score_svm.std())
y_pred_flatten = [item for sublist in y_pred_svm for item in sublist]
overall_accuracy_svm = accuracy_score(y_test_flatten, y_pred_flatten)
print('The OVERALL test accuracy for rf is', overall_accuracy_svm)
```

The average test accuracy for svm is 0.98000000000000000The test accuracy standard deviation for svm is 0.023184046238739247The OVERALL test accuracy for rf is 0.98

0.4.4 (iv) Logistic Regression with Elastic Net Regularization

```
[31]: # Initialze empty variables
      best_model_logR = []
      best_valscore_logR = np.zeros(len(random_states))
      test_score_logR = np.zeros(len(random_states))
      y_pred_logR = []
      for i in range(len(random_states)):
          # Define hyperparameter grid for tuning
          param grid = {
              'penalty': ['elasticnet'],
              'solver': ['saga'],
              'l1_ratio': np.linspace(0,1,8),
              'C': [0.01, 0.05, 0.1, 0.5, 1, 1.5, 2]
          }
          val_score = np.zeros(len(ParameterGrid(param_grid)))
          models = \Pi
          # train the model
          for p in range(len(ParameterGrid(param_grid))):
              params = ParameterGrid(param_grid)[p]
              clf = LogisticRegression(**params, random_state = random_states[i])
              clf.fit(X train[i], y train[i])
              models.append(clf)
              y val pred = clf.predict(X val[i])
              val_score[p] = accuracy_score(y_val[i], y_val_pred)
          # get the best validation score
          best_valscore_logR[i] = np.max(val_score)
          best_model_logR.append(models[np.argmax(val_score)])
          # evaluate the model using test set (test score)
          y_test_pred = best_model_logR[i].predict(X_test[i])
          test_score_logR[i] = accuracy_score(y_test[i], y_test_pred)
          y_pred_logR.append(y_test_pred)
```

```
print('Random state: ', random_states[i])
    print('The best model parameters are: ', ParameterGrid(param grid)[np.
 →argmax(val_score)])
    print('Best validation score: ', best_valscore_logR[i])
    print('Test accuracy: ', test score logR[i])
    print(' ')
Random state: 7271
The best model parameters are: {'solver': 'saga', 'penalty': 'elasticnet',
'l1_ratio': 0.0, 'C': 0.01}
Best validation score: 1.0
Test accuracy: 0.95
Random state: 861
The best model parameters are: {'solver': 'saga', 'penalty': 'elasticnet',
'l1_ratio': 0.0, 'C': 0.05}
Best validation score: 0.9875
Test accuracy: 0.9875
Random state: 5391
The best model parameters are: {'solver': 'saga', 'penalty': 'elasticnet',
'l1_ratio': 0.0, 'C': 0.05}
Best validation score: 0.9875
Test accuracy: 0.975
Random state: 5192
The best model parameters are: {'solver': 'saga', 'penalty': 'elasticnet',
'l1 ratio': 0.0, 'C': 0.01}
Best validation score: 1.0
Test accuracy: 1.0
Random state: 5735
The best model parameters are: {'solver': 'saga', 'penalty': 'elasticnet',
'l1_ratio': 0.0, 'C': 0.05}
Best validation score: 1.0
Test accuracy: 1.0
Random state: 6266
The best model parameters are: {'solver': 'saga', 'penalty': 'elasticnet',
'l1_ratio': 0.0, 'C': 0.01}
Best validation score: 1.0
Test accuracy: 0.95
Random state: 467
The best model parameters are: {'solver': 'saga', 'penalty': 'elasticnet',
'l1 ratio': 0.0, 'C': 0.05}
Best validation score: 1.0
```

```
Test accuracy: 0.95
     Random state: 4427
     The best model parameters are: {'solver': 'saga', 'penalty': 'elasticnet',
     'l1 ratio': 0.14285714285714285, 'C': 0.01}
     Best validation score: 1.0
     Test accuracy: 0.975
     Random state: 5579
     The best model parameters are: {'solver': 'saga', 'penalty': 'elasticnet',
     'l1_ratio': 0.0, 'C': 0.1}
     Best validation score: 1.0
     Test accuracy: 1.0
     Random state: 8323
     The best model parameters are: {'solver': 'saga', 'penalty': 'elasticnet',
     'l1_ratio': 0.0, 'C': 0.01}
     Best validation score: 0.9875
     Test accuracy: 0.9875
[32]: print('The average test accuracy for Logistic Regression with Elastic Net
      →Regularization is', test_score_logR.mean())
      print('The test accuracy standard deviation for Logistic Regression with,
       ⇒Elastic Net Regularization is', test_score_logR.std())
      y_pred_flatten = [item for sublist in y_pred_logR for item in sublist]
      overall_accuracy_logR = accuracy_score(y_test_flatten, y_pred_flatten)
      print('The OVERALL test accuracy for Logistic Regression with Elastic Net⊔

¬Regularization is', overall_accuracy_logR)
     The average test accuracy for Logistic Regression with Elastic Net
     Regularization is 0.9775
     The test accuracy standard deviation for Logistic Regression with Elastic Net
     Regularization is 0.020000000000000025
     The OVERALL test accuracy for Logistic Regression with Elastic Net
     Regularization is 0.9775
     0.4.5 (v) KNN
[33]: # Initialze empty variables
      best model knn = []
      best_valscore_knn = np.zeros(len(random_states))
      test_score_knn = np.zeros(len(random_states))
      y_pred_knn = []
      for i in range(len(random_states)):
          # Define hyperparameter grid for tuning
```

param_grid = {

```
'n_neighbors': [1, 3, 5, 10], # Number of neighbors
        'weights': ['uniform', 'distance'], # Weighting scheme
        'metric': ['euclidean', 'manhattan', 'minkowski'], # Distance metrics
    }
    val_score = np.zeros(len(ParameterGrid(param_grid)))
    models = \Pi
    # train the model
    for p in range(len(ParameterGrid(param_grid))):
        params = ParameterGrid(param grid)[p]
        clf = KNeighborsClassifier(**params)
        clf.fit(X_train[i], y_train[i])
        models.append(clf)
        y_val_pred = clf.predict(X_val[i])
        val_score[p] = accuracy_score(y_val[i], y_val_pred)
    # get the best validation score
    best_valscore_knn[i] = np.max(val_score)
    best_model_knn.append(models[np.argmax(val_score)])
    # evaluate the model using test set (test score)
    y_test_pred = best_model_knn[i].predict(X_test[i])
    test_score_knn[i] = accuracy_score(y_test[i], y_test_pred)
    y_pred_knn.append(y_test_pred)
    print('Random state: ', random_states[i])
    print('The best model parameters are: ', ParameterGrid(param_grid)[np.
 →argmax(val_score)])
    print('Best validation score: ', best_valscore_knn[i])
    print('Test accuracy: ', test_score_knn[i])
    print(' ')
Random state: 7271
The best model parameters are: {'weights': 'uniform', 'n neighbors': 3,
'metric': 'euclidean'}
Best validation score: 1.0
Test accuracy: 0.975
Random state: 861
The best model parameters are: {'weights': 'uniform', 'n_neighbors': 1,
'metric': 'euclidean'}
Best validation score: 0.9875
Test accuracy: 0.9875
Random state: 5391
The best model parameters are: {'weights': 'distance', 'n_neighbors': 10,
'metric': 'euclidean'}
```

```
Best validation score: 0.975
Test accuracy: 0.975
Random state: 5192
The best model parameters are: {'weights': 'uniform', 'n_neighbors': 1,
'metric': 'euclidean'}
Best validation score:
                      0.9875
Test accuracy: 0.9875
Random state: 5735
The best model parameters are: {'weights': 'uniform', 'n_neighbors': 1,
'metric': 'euclidean'}
Best validation score: 1.0
Test accuracy: 0.9625
Random state: 6266
The best model parameters are: {'weights': 'uniform', 'n_neighbors': 1,
'metric': 'euclidean'}
Best validation score: 0.975
Test accuracy: 1.0
Random state: 467
The best model parameters are: {'weights': 'uniform', 'n_neighbors': 1,
'metric': 'euclidean'}
Best validation score: 1.0
Test accuracy: 0.9625
Random state: 4427
The best model parameters are: {'weights': 'uniform', 'n_neighbors': 1,
'metric': 'euclidean'}
Best validation score: 1.0
Test accuracy: 0.975
Random state: 5579
The best model parameters are: {'weights': 'uniform', 'n_neighbors': 1,
'metric': 'manhattan'}
Best validation score: 1.0
Test accuracy: 1.0
Random state: 8323
The best model parameters are: {'weights': 'uniform', 'n_neighbors': 1,
'metric': 'euclidean'}
Best validation score: 0.9875
Test accuracy: 0.9875
```

```
[34]: print('The average test accuracy for KNN is', test_score_knn.mean())
print('The test accuracy standard deviation for KNN is', test_score_knn.std())
y_pred_flatten = [item for sublist in y_pred_knn for item in sublist]
overall_accuracy_knn = accuracy_score(y_test_flatten, y_pred_flatten)
print('The OVERALL test accuracy for Logistic KNN is', overall_accuracy_knn)
```

The average test accuracy for KNN is 0.98125 The test accuracy standard deviation for KNN is 0.0128086884574495 The OVERALL test accuracy for Logistic KNN is 0.98125

0.4.6 Summary of result

Since we have balanced dataset (0.6 percentage and 0.4 percentage for binary target variable), we can use the accuracy score as our baseline score. And the strategy of basaline accuracy score is to take the majority of one class divide by the total number of data points. Here I use y_test from all 10 different random states.

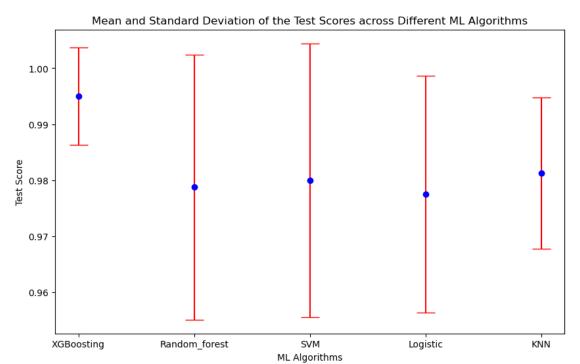
```
[35]: x = sum(y_test_flatten)/len(y_test_flatten)
if x >= 0.5:
    baseline_score = x
else:
    baseline_score = 1 - x
print(baseline_score)
```

0.63125

```
XGBoosting
             Random_forest
                                SVM Logistic
                                                  KNN
0
       1.0000
                     0.9500 0.9625
                                       0.9500 0.9750
                                       0.9875 0.9875
1
       1.0000
                      0.9875
                             0.9625
2
                     0.9875 0.9875
                                       0.9750 0.9750
       0.9875
3
       1.0000
                      0.9375
                             1.0000
                                       1.0000 0.9875
4
       1.0000
                      1.0000
                             1.0000
                                       1.0000 0.9625
5
       0.9875
                     0.9625
                             0.9750
                                       0.9500 1.0000
6
       0.9750
                     0.9625 1.0000
                                       0.9500 0.9625
7
       1.0000
                     1.0000 0.9250
                                       0.9750 0.9750
8
       1.0000
                     1.0000 0.9875
                                       1.0000 1.0000
9
       1.0000
                     1.0000 1.0000
                                       0.9875 0.9875
```

```
[37]: test_mean = test_score_sum.mean()
test_std = test_score_sum.std()
# plot means and standard deviations of test scores
plt.figure(figsize = (10,6))
```

```
plt.errorbar(x = test_mean.index, y = test_mean, yerr = test_std, fmt_\( \) \( \text{\textstar} \) \( \text{\
```

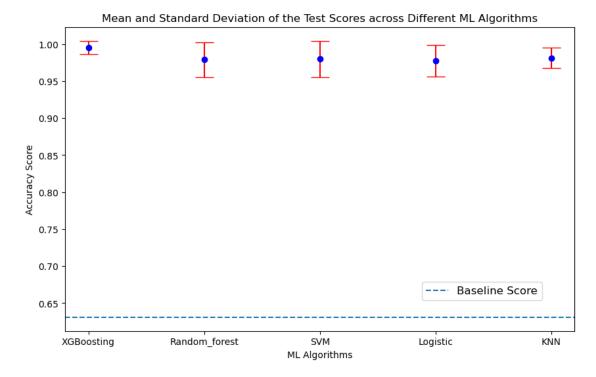


```
XGBoosting 0.99500
Random_forest 0.97875
SVM 0.98000
Logistic 0.97750
KNN 0.98125
```

dtype: float64

```
[38]: # add a baseline on the plot
plt.figure(figsize = (10,6))
plt.errorbar(x = test_mean.index, y = test_mean, yerr = test_std, fmt_u

--'o',color='blue', ecolor='red', capsize=10)
plt.xlabel('ML Algorithms')
plt.ylabel('Accuracy Score')
```



Baseline_score: 0.63125

0.5 Model Inspectation

Here we choose to use XGBoosting model because it has the highest accuracy score, and then we use the y_pred from this model to do the model inspection. ### (i) confusion matrix

```
[39]: # confusion matrix using XGBoosting

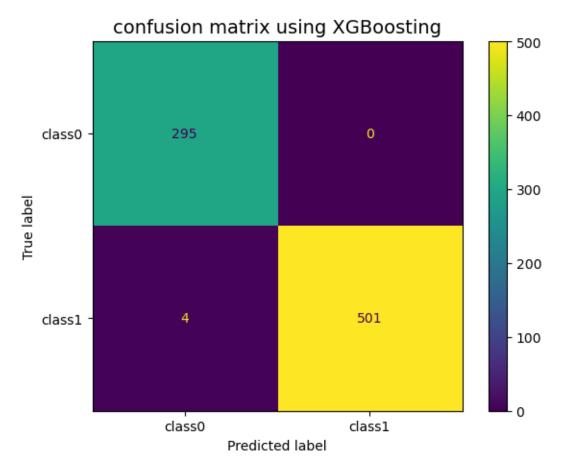
cm = confusion_matrix(y_test_flatten, [item for sublist in y_pred_xgb for item_
in sublist])

disp = ConfusionMatrixDisplay(cm, display_labels = ['class0', 'class1'])

fig, ax = plt.subplots(figsize = (7,5))

disp.plot(ax=ax)

plt.title('confusion matrix using XGBoosting', fontsize = 14)
```



recall: 0.992079207921

precision: 1.0 accuracy: 0.995

f1 score: 0.9960238568588469

Next, we want to investigate some bias:

```
[40]: finaldf = []
      for i in range(len(random_states)):
          combinecol = pd.concat([pd.DataFrame(X_test[i]).reset_index(drop=True),
                                  pd.DataFrame(y_test[i]).reset_index(drop=True),
                                  pd.DataFrame(y_pred_xgb[i]).
       oreset_index(drop=True)], axis = 1)
          colnames = list(X_test[i].columns) + ['y_true', 'y_pred']
          combinecol.columns = colnames
          finaldf.append(combinecol)
[41]: df1 = []
      df2 = []
      df3 = []
      df4 = []
      df5 = []
```

```
df6 = []
df7 = []
df8 = \Pi
df9 = []
df10 = []
for i in range(len(random_states)):
    # Data and metrics for hypertension people
    temp_data1 = finaldf[i][X_test_withoutstd[i]['htn_yes'] == 1][['y_true',__
 df1.append(temp_data1)
    # Data and metrics for diabetes mellitus people
    temp_data2 = finaldf[i][X_test_withoutstd[i]['dm_yes'] == 1][['y_true',_

        'y_pred']]

    df2.append(temp_data2)
    # Data and metrics for coronary artery disease people
    temp_data3 = finaldf[i][X_test_withoutstd[i]['cad_yes'] == 1][['y_true',__

    'y_pred']]

    df3.append(temp_data3)
    # Data and metrics for appetite people
    temp_data4 = finaldf[i][X_test_withoutstd[i]['appet_good'] == 1][['y_true',_

    'y_pred']]

    df4.append(temp_data4)
    # Data and metrics for pedal edema people
    temp_data5 = finaldf[i][X_test_withoutstd[i]['pe_yes'] == 1][['y_true',_

    'y_pred']]
```

```
df5.append(temp_data5)
    # Data and metrics for anemia people
   temp_data6 = finaldf[i][X_test_withoutstd[i]['ane_yes'] == 1][['y_true',__

    'y_pred']]

   df6.append(temp data6)
    # Data and metrics for red blood cells people
   temp_data7 = finaldf[i][X_test_withoutstd[i]['rbc_normal'] == 1][['y_true',__

    'y_pred']]

   df7.append(temp_data7)
    # Data and metrics for pus cell people
   temp_data8 = finaldf[i][X_test_withoutstd[i]['pc_normal'] == 1][['y_true',_

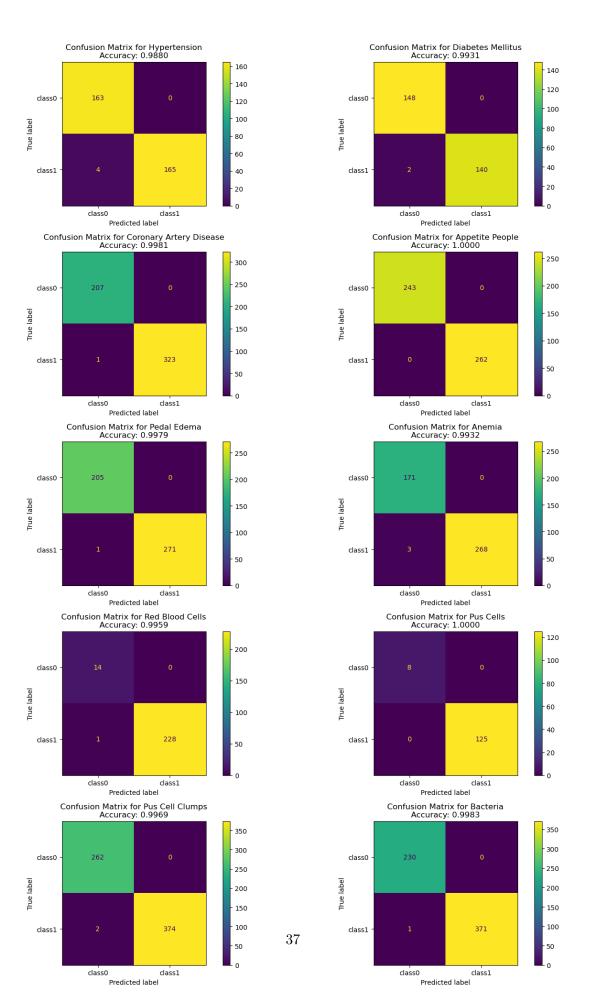
        'y_pred']]

   df8.append(temp_data8)
    # Data and metrics for pus cell clumps people
   temp_data9 = finaldf[i][X_test_withoutstd[i]['pcc_present'] ==__
 df9.append(temp data9)
    # Data and metrics for becteria people
   temp_data10 = finaldf[i][X_test_withoutstd[i]['ba_present'] ==__
 df10.append(temp_data10)
df1 = pd.concat(df1, axis = 0)
df2 = pd.concat(df2, axis = 0)
df3 = pd.concat(df3, axis = 0)
df4 = pd.concat(df4, axis = 0)
df5 = pd.concat(df5, axis = 0)
df6 = pd.concat(df6, axis = 0)
df7 = pd.concat(df7, axis = 0)
df8 = pd.concat(df8, axis = 0)
df9 = pd.concat(df9, axis = 0)
df10 = pd.concat(df10, axis = 0)
ac1 = accuracy_score(df1['y_true'], df1['y_pred'])
cm1 = confusion_matrix(df1['y_true'], df1['y_pred'])
ac2 = accuracy_score(df2['y_true'], df2['y_pred'])
cm2 = confusion_matrix(df2['y_true'], df2['y_pred'])
ac3 = accuracy_score(df3['y_true'], df3['y_pred'])
cm3 = confusion_matrix(df3['y_true'], df3['y_pred'])
ac4 = accuracy_score(df4['y_true'], df4['y_pred'])
cm4 = confusion_matrix(df4['y_true'], df4['y_pred'])
ac5 = accuracy_score(df5['y_true'], df5['y_pred'])
```

```
cm5 = confusion_matrix(df5['y_true'], df5['y_pred'])
ac6 = accuracy_score(df6['y_true'], df6['y_pred'])
cm6 = confusion_matrix(df6['y_true'], df6['y_pred'])
ac7 = accuracy_score(df7['y_true'], df7['y_pred'])
cm7 = confusion_matrix(df7['y_true'], df7['y_pred'])
ac8 = accuracy_score(df8['y_true'], df8['y_pred'])
cm8 = confusion_matrix(df8['y_true'], df8['y_pred'])
ac9 = accuracy_score(df9['y_true'], df9['y_pred'])
cm9 = confusion_matrix(df9['y_true'], df9['y_pred'])
ac10 = accuracy_score(df10['y_true'], df10['y_pred'])
cm10 = confusion_matrix(df10['y_true'], df10['y_pred'])
# Plots
# Create a grid for the plots (5 rows x 2 columns)
fig, axes = plt.subplots(5, 2, figsize=(14, 20)) # Adjust figsize for better_
 \hookrightarrow visualization
# Plot for hypertension people
disp1 = ConfusionMatrixDisplay(cm1, display_labels=['class0', 'class1'])
disp1.plot(ax=axes[0, 0], values_format='d')
axes[0, 0].set title(f'Confusion Matrix for Hypertension\nAccuracy: {ac1:.4f}',,,
 ofontsize=12)
# Plot for diabetes mellitus people
disp2 = ConfusionMatrixDisplay(cm2, display_labels=['class0', 'class1'])
disp2.plot(ax=axes[0, 1], values_format='d')
axes[0, 1].set title(f'Confusion Matrix for Diabetes Mellitus\nAccuracy: {ac2:.
\rightarrow4f}', fontsize=12)
# Plot for coronary artery disease people
disp3 = ConfusionMatrixDisplay(cm3, display_labels=['class0', 'class1'])
disp3.plot(ax=axes[1, 0], values_format='d')
axes[1, 0].set_title(f'Confusion Matrix for Coronary Artery Disease\nAccuracy:
 \rightarrow{ac3:.4f}', fontsize=12)
# Plot for appetite people
disp4 = ConfusionMatrixDisplay(cm4, display_labels=['class0', 'class1'])
disp4.plot(ax=axes[1, 1], values_format='d')
axes[1, 1].set_title(f'Confusion Matrix for Appetite People\nAccuracy: {ac4:.
 \rightarrow4f}', fontsize=12)
# Plot for pedal edema people
disp5 = ConfusionMatrixDisplay(cm5, display_labels=['class0', 'class1'])
disp5.plot(ax=axes[2, 0], values_format='d')
axes[2, 0].set_title(f'Confusion Matrix for Pedal Edema\nAccuracy: {ac5:.4f}', __

fontsize=12)
```

```
# Plot for anemia people
disp6 = ConfusionMatrixDisplay(cm6, display_labels=['class0', 'class1'])
disp6.plot(ax=axes[2, 1], values_format='d')
axes[2, 1].set_title(f'Confusion Matrix for Anemia\nAccuracy: {ac6:.4f}', ___
 ⇔fontsize=12)
# Plot for red blood cells people
disp7 = ConfusionMatrixDisplay(cm7, display_labels=['class0', 'class1'])
disp7.plot(ax=axes[3, 0], values_format='d')
axes[3, 0].set_title(f'Confusion Matrix for Red Blood Cells\nAccuracy: {ac7:.
 \hookrightarrow4f}', fontsize=12)
# Plot for pus cell people
disp8 = ConfusionMatrixDisplay(cm8, display_labels=['class0', 'class1'])
disp8.plot(ax=axes[3, 1], values_format='d')
axes[3, 1].set_title(f'Confusion Matrix for Pus Cells\nAccuracy: {ac8:.4f}', __
 ⇔fontsize=12)
# Plot for pus cell clumps people
disp9 = ConfusionMatrixDisplay(cm9, display_labels=['class0', 'class1'])
disp9.plot(ax=axes[4, 0], values_format='d')
axes[4, 0].set title(f'Confusion Matrix for Pus Cell Clumps\nAccuracy: {ac9:.
 \hookrightarrow4f}', fontsize=12)
# Plot for bacteria people
disp10 = ConfusionMatrixDisplay(cm10, display_labels=['class0', 'class1'])
disp10.plot(ax=axes[4, 1], values_format='d')
axes[4, 1].set_title(f'Confusion Matrix for Bacteria\nAccuracy: {ac10:.4f}', __
 ⇔fontsize=12)
# Adjust layout to avoid overlapping
plt.tight layout()
plt.savefig('/Users/ericdavis/Desktop/data1030/FinalProject1030/
 ⇔confusion_matrix_forbias.png', dpi = 350)
plt.show()
```

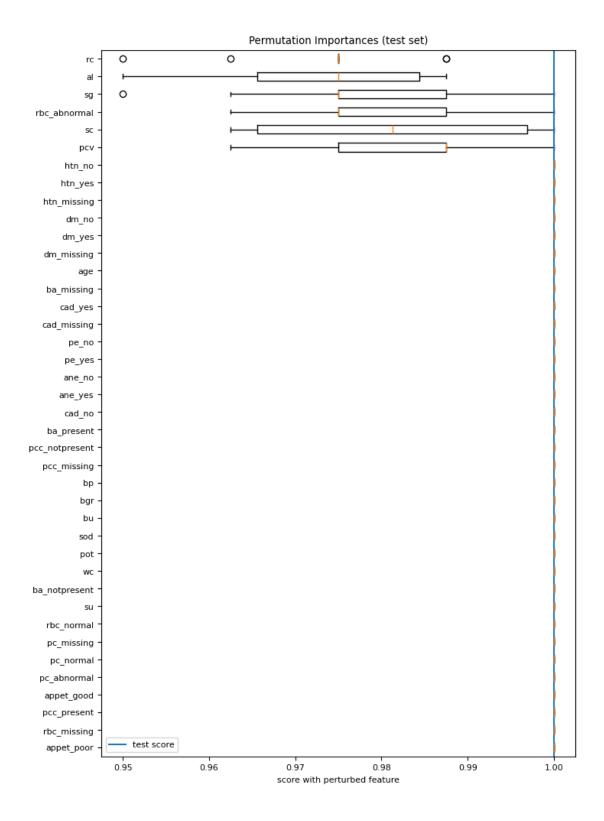


0.5.1 (ii) feature importance

Global importance

method1:

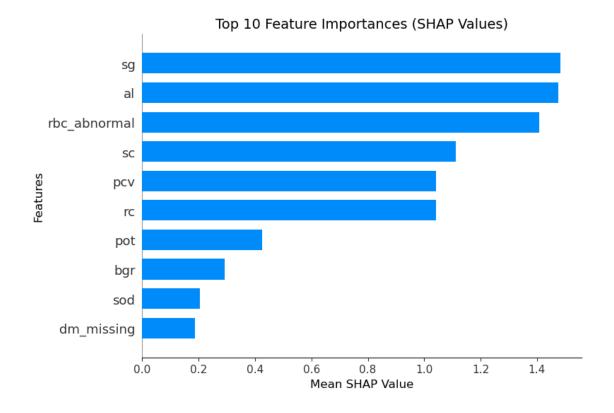
```
[28]: np.random.seed(42) # for reproducibility
      nr runs = 10
      best_index = np.argmax(test_score_xgb)
      feature names = X test[best index].columns
      scores = np.zeros([len(feature_names), nr_runs])
      for i in range(len(feature_names)):
          acc_scores = []
          for j in range(nr_runs):
              X_test_shuffled = X_test[best_index].copy()
              X_test_shuffled[feature_names[i]] = np.random.
       opermutation(X_test[best_index][feature_names[i]].values)
              acc_scores.append(best_model_xgb[best_index].score(X_test_shuffled,_
       y test[best index]))
          scores[i] = acc_scores
      sorted_indcs = np.argsort(np.mean(scores, axis = 1))
      plt.rcParams.update({'font.size': 8})
      plt.figure(figsize=(8,12))
      plt.boxplot(scores[sorted_indcs].T,tick_labels = feature_names[sorted_indcs],_
       →vert=False)
      plt.title("Permutation Importances (test set)")
      plt.xlabel('score with perturbed feature')
      plt.axvline(test_score_xgb[best_index], label='test_score')
      plt.gca().invert_yaxis()
      plt.legend()
      plt.savefig('/Users/ericdavis/Desktop/data1030/FinalProject1030/permutation_imp.
       \hookrightarrowpng', dpi = 350)
      plt.show()
```



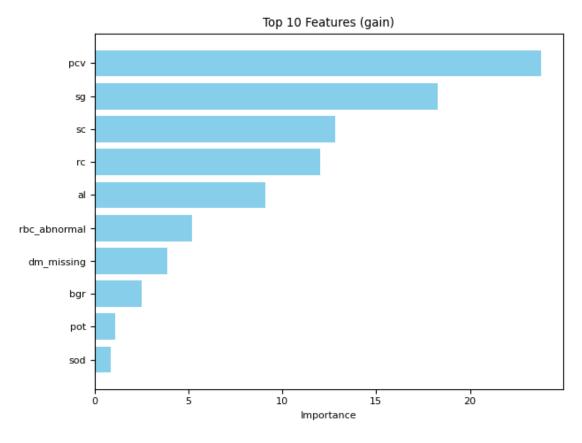
method 2:

<IPython.core.display.HTML object>

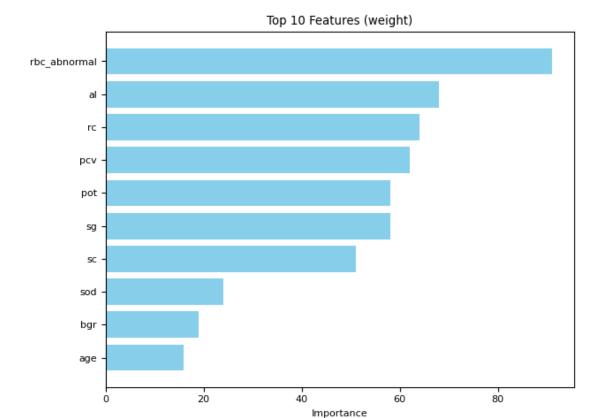
The shape of the shap values is (80, 40).

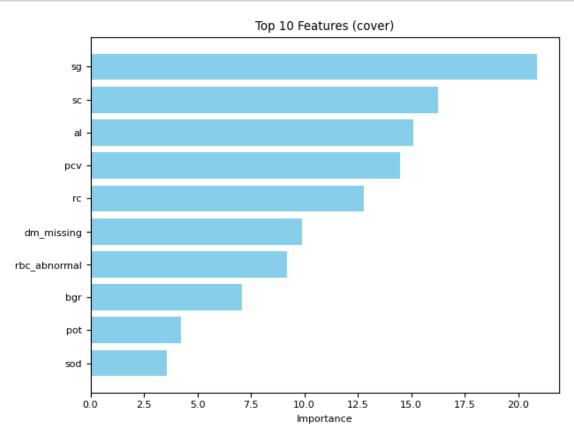


method3:



```
method 4:
[53]: booster = best_model_xgb[best_index].get_booster()
    feature_importance = booster.get_score(importance_type='weight')
```



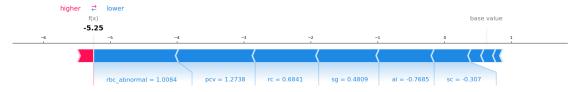


local importance

```
[55]: index = 68 # predict class 0

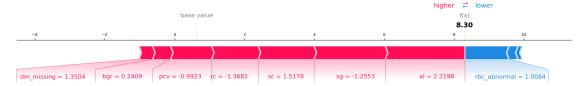
# Extract the SHAP values and feature values for the specific index
shap_vals = np.round(shap_values[index, :],4)
features = np.round(X_test[best_index].iloc[index, :],4) # Feature values for_
the specific data point

# Sort SHAP values and corresponding features by absolute SHAP value
```



<Figure size 640x480 with 0 Axes>

```
features=top_features,
  feature_names=top_feature_names,
  matplotlib = True
)
plt.tight_layout()
plt.show()
```



<Figure size 640x480 with 0 Axes>

```
[]:
```