Final Project Submission

Please fill out:

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Scheduled project review date/time:

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· Blog post URL:

Business Problem

Tanzania has water pumps installed throughout the country in order to make pottable water accessible to it's residents. Maintaining these water pumps can prove to be a challenge, since many are in remote areas that are difficult to monitor.

The goal of this project is to build a model that can predict which water pumps are in need of repair in order to efficiently dispatch technicians to the sites most in need and provide as many people as possible with clean water.

The Data

A dataset of Tanzania water pumps is maintained by Taarifa in a waterpoints dashboard by aggregating data from the Tanzania Ministry of Water. This data set keeps track of information regarding the location of the water pump, information about the pump (contruction year, extraction type, etc.), the people involved in the pump (installers, management, population), as well as the information regarding the source of water feeding the pump.

Additionally, the training test set provided has another test set containing the id nubmer of every pump and wheter it is functional, functional but in need of repair, or non-functional. This will allow us to train and validate our model.

Prior to cleaning, there are 59,400 water pumps in the training data set.

Setup

Import relevant packages

```
In [1]: import sqlite3
import pandas as pd
import numpy as np
import scipy.stats as stats
import statsmodels.api as sm
import json

from sklearn.preprocessing import OneHotEncoder
from sklearn.tree import DecisionTreeRegressor, DecisionTreeClassifier
from sklearn.model_selection import train_test_split
from sklearn.metrics import f1_score, mean_squared_error, r2_score, roc_curv
from sklearn.ensemble import RandomForestClassifier

import seaborn as sns
import matplotlib.pyplot as plt
# plt.style.use('seaborn-v0_8-whitegrid')
%matplotlib inline
```

Load and Clean Data

Training values set

Preview Data

```
In [2]: training_values = pd.read_csv('data/Training_set_values.csv')
    print(f'number of rows: {len(training_values)}')
    training_values.head()
```

number of rows: 59400

Out[2]:

	id	amount_tsh	date_recorded	funder	gps_height	installer	longitude	latitude	wpt_name
0	69572	6000.0	2011-03-14	Roman	1390	Roman	34.938093	-9.856322	none
1	8776	0.0	2013-03-06	Grumeti	1399	GRUMETI	34.698766	-2.147466	Zahanati
2	34310	25.0	2013-02-25	Lottery Club	686	World vision	37.460664	-3.821329	Kwa Mahundi
3	67743	0.0	2013-01-28	Unicef	263	UNICEF	38.486161	-11.155298	Zahanati Ya Nanyumbu
4	19728	0.0	2011-07-13	Action In A	0	Artisan	31.130847	-1.825359	Shuleni

5 rows × 40 columns

Investigate Missing Data

In [3]: training_values.info()

<class 'pandas.core.frame.DataFrame'>
RangeIndex: 59400 entries, 0 to 59399
Data columns (total 40 columns):

Column		ull Count	Dtype				
id	59400	non-null	 int64				
			float64				
			object				
funder	55765		object				
gps height	59400		int64				
installer	55745	non-null	object				
longitude	59400	non-null	float64				
latitude	59400	non-null	float64				
wpt_name	59400	non-null	object				
num_private	59400	non-null	int64				
basin	59400	non-null	object				
subvillage	59029	non-null	object				
region	59400	non-null	object				
region_code	59400	non-null	int64				
district_code	59400	non-null	int64				
lga	59400	non-null	object				
ward	59400	non-null	object				
population	59400	non-null	int64				
<pre>public_meeting</pre>	56066	non-null	object				
recorded_by	59400	non-null	object				
scheme_management	55523	non-null	object				
scheme_name	31234	non-null	object				
permit	56344	non-null	object				
construction_year			int64				
		non-null	object				
		non-null	object				
			object				
_			object				
			object				
•			object				
			object				
			object				
			object				
			object				
			object				
			object				
_ • •			object				
_			object				
			object				
waterpoint_type_group			object				
memory usage: 18.1+ MB							
	Column id amount_tsh date_recorded funder gps_height installer longitude latitude wpt_name num_private basin subvillage region region_code district_code lga ward population public_meeting recorded_by scheme_management scheme_name permit construction_year extraction_type extraction_type_class management management payment payme	id 59400 amount_tsh 59400 funder 55765 gps_height 59400 installer 55745 longitude 59400 wpt_name 59400 num_private 59400 basin 59400 region_code 59400 region_code 59400 district_code 59400 ga 59400 ward 59400 population 59400 population 59400 poblic_meeting 56066 recorded_by 59400 scheme_management 55523 scheme_name 31234 permit 56344 construction_year 59400 extraction_type 59400 extraction_type_group 59400 extraction_type_group 59400 management 59400 management 59400 management 59400 payment 59400 pa	Column Non-Null Count id 59400 non-null amount_tsh 59400 non-null funder 55765 non-null gps_height 59400 non-null statler 55745 non-null latitude 59400 non-null wpt_name 59400 non-null wpt_name 59400 non-null wpt_name 59400 non-null subvillage 59400 non-null subvillage 59029 non-null region 59400 non-null region_code 59400 non-null ga 59400 non-null population 59400 non-null population 59400 non-null recorded_by 59400 non-null recorded_by 59400 non-null scheme_management 55523 non-null scheme_name 31234 non-null permit 56344 non-null extraction_type 59400 non-null extraction_type_group 59400 non-null management 59400 non-null management 59400 non-null payment 59400 non-null source 59400 non-null sourc				

Clean data

Investigate funder, installer, subvillage, public_meeting, scheme management, scheme_name, permit

- funder:
 - 3635 missing values (6.1%)
 - Not Known, Unknown, categories
 - 0 category (same as none?)
- installer:
 - 3655 missing values (6.1%)
 - uknown, Unknown, Not known categories
- subvillage:
 - 371 missing values (0.6%)
 - relatively few missing values
 - 19287 unique values too many to consolidate inot meaningful categories
 - enough other location based categories
 - drop column
- public_meeting:
 - 3334 missing values (5.6%)
 - binary true/false
 - drop missing rows want to keep it binary
- · scheme management:
 - 3877 missing values (6.5%)
 - will likeley create an "unknown" column
 - one occurance of None remove (unable to have an example in both test and training sets)
- · scheme name:
 - 28166 missing values (47.4%)
 - too much missing data drop this
- permit:
 - 3056 missing values (5.1%)
 - binary true/false
 - drop missing rows want to keep it binary

After cleaning, 53,277 data points (89% of original set)

permit, and public_meeting columns:

remove NaN values

In [4]: training_values = training_values.dropna(axis=0, subset=['permit', 'public_m'

Drop unnecessary columns:

columns with majority unique values or data that is not relevant or redundant

- scheme_name, wpt_name, date_recorded, subvillage, ward
- quantity_group and quantity are the same, drop one
- extraction_type, extraction_type_group, extraction_type_class are varying hierarchies of the same feature, take the middle one.
- source, source_class, source_type are varying hierarchies of the same feature. take the middle one.
- waterpoint_type and waterpoint_type_group: virtually the same, except waterpoint_type splits communal standpipe into communal standpipe and communal statepipe multiple. keep waterpoint_type.
- payment and payment_type are identical. drop payment_type.

```
In [5]: drop_cols = ['scheme_name', 'wpt_name', 'date_recorded', 'subvillage', 'ward
training_values = training_values.drop(columns=drop_cols, axis=1)
```

Manage "Unknown" and "Other" Categories

- Remove NaN values or convert to "unknown".
- Combine any "not known" variations into one "unknown" column
- Combine categories with low value counts into a single "other" column

```
In [6]: ## clean up permit, and public_meeting columns
    training_values = training_values.dropna(axis=0, subset=['permit', 'public_m

In [7]: ## clean up funder column
    # combine Not Known, Unknown, and Nan categories into one 'Unknown'
    training_values['funder'] = training_values['funder'].fillna('Unknown')
    training_values['funder'] = training_values['funder'].map(lambda x: 'Unknown

# create an "other" column for all funders with less than 10 wells
    # Note: 156 was chosen as it combined sufficient small categories withou
    small_funders = []
    for key, value in training_values['funder'].value_counts().to_dict().items()
        if value < 16:
            small_funders.append(key)

training_values['funder'] = training_values['funder'].map(lambda x: 'Other'</pre>
```

```
In [8]: ean up installer column
       bine uknown, Unknown, Not known, and Nan categories into one 'Unknown'
       ing_values['installer'] = training_values['installer'].fillna('Unknown')
       ing_values['installer'] = training_values['installer'].map(lambda x: 'Unknown')
       ing_values['installer'] = training_values['installer'].map(lambda x: 'Unknown')
       ate an "other" column for all installers with less than 15 wells
        Note: 15 was chosen as it combined sufficient small categories without making
        installers = []
       ey, value in training values['installer'].value counts().to dict().items():
       f value < 15:
          small installers.append(key)
       ing values['installer'] = training values['installer'].map(lambda x: 'Other'
In [9]: ## clean up scheme management column
        # convert Nan values into one 'Unknown'
        training_values['scheme_management'] = training_values['scheme_management'].
```

Remove categories with only 1 or 2 occurances

A few columns have categories with very few examples, not enough to be reliably split between the training and test data set, so those categories will be removed

```
In [10]: # Scheme management: remove one row with "None"
         training_values = training_values[training_values['scheme_management'] != 'N
In [11]: # lga: remove one row with "Nyamagana"
         training values = training values[training values['lga'] != 'Nyamagana']
```

Investigate value counts of final columns

Visually inspect the number of categories in each column to confirm that no column will create a large number of columns

- any categorical columns should have a "managable" number of columns
- any numberical columns do not need to have low value counts, as they will not be encoded

```
In [12]: for col in list(training values.columns):
             print(f'{col}: {len(training_values[col].value_counts())}')
         # not dummy variables
         # id, gps_height, population, construction_year
         id: 53279
         amount_tsh: 91
         funder: 302
         gps_height: 2426
         installer: 304
         longitude: 51779
         latitude: 51781
         num_private: 61
         basin: 9
         region: 21
         region code: 27
         district_code: 20
         lga: 121
         population: 1006
         public_meeting: 2
         recorded_by: 1
         scheme_management: 12
         permit: 2
         construction_year: 55
         extraction_type_group: 13
         management: 12
         management_group: 5
         payment: 7
         water_quality: 8
         quality_group: 6
         quantity_group: 5
         source: 10
         source_type: 7
         waterpoint_type: 7
```

Training Set Labels

Load the training set lables and restrict to only contain ids that are included in the cleaned training values dataset

```
In [14]: # create a mask for ids present in the filtered training values set
    mask = training_labels['id'].isin(list(training_values.id))

# apply the mask
    training_labels = training_labels[mask]
    print(len(training_values))
```

53279

Full Data Set Visualizations

Initial visualizations to get a look into the data set

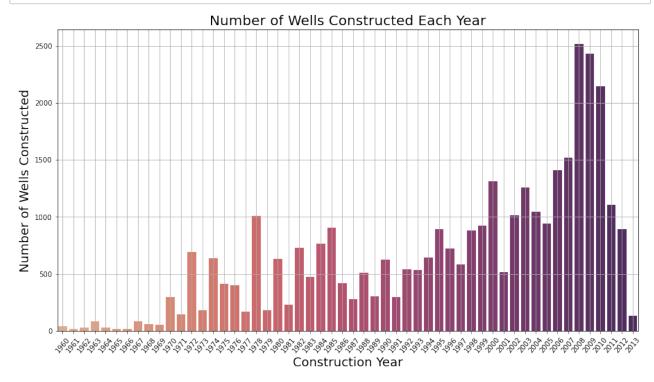
View the number of wells constructed per year

```
In [15]: # group by year constucted
    years = training_values.construction_year.value_counts().index[1:]
    year_counts = training_values.construction_year.value_counts().values[1:]

# count the number of wells in each year
    year_data = {
        'year': training_values.construction_year.value_counts().index[1:],
        'count': training_values.construction_year.value_counts().values[1:]
}

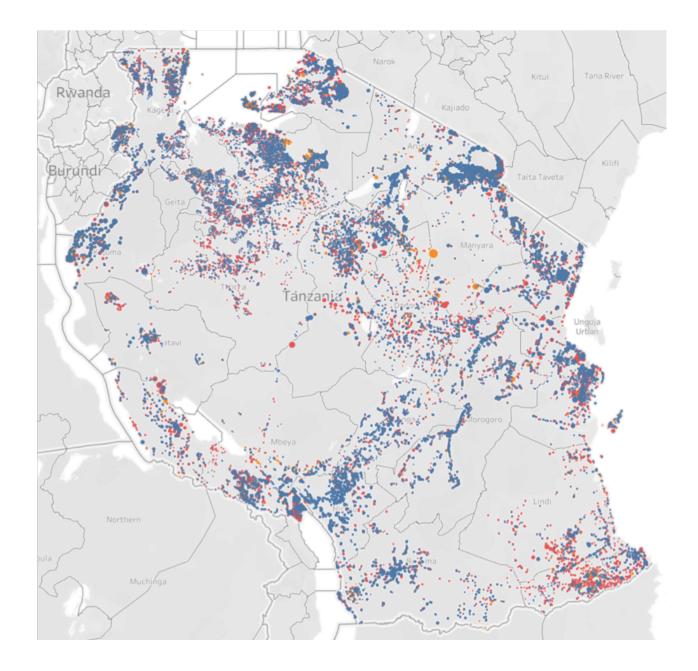
year_df = pd.DataFrame(year_data)

plt.figure(figsize=(15,8))
    sns.barplot(data=year_df, x='year', y='count', palette='flare')
    plt.xticks(rotation=50)
    plt.xlabel('Construction Year', fontsize=18)
    plt.ylabel('Number of Wells Constructed', fontsize=18)
    plt.title('Number of Wells Constructed Each Year', fontsize=20)
    plt.grid()
```



Locations of wells across Tanzania

- map generated using Tableau using well coordinates and inserted into this notbook
 - blue: functional
 - orange: functional, needs repair
 - red: non functional



Data Prep

Prepare the data to be compatible with classification

Split the given training set into a train set and a test set

In [20]: # initialize the One Hot Encoder
ohe = OneHotEncoder()

```
Inputs:
                     - X: dataframe to encode
                     - non_categorical_cols: list of column names to exclude from end
                     - X_ohe_final: final encoded dataframe containing both encoded d
             # create two dataframes: categorical and non categorical
             categorical_df = X.drop(columns=non_categorical_cols, axis=1)
             non categorical df = X[non categorical cols]
             # encode the categorical dataframe and convert back to a dataframe
             X_ohe_categorical = ohe.fit_transform(categorical_df).toarray()
             ohe df = pd.DataFrame(X ohe categorical, columns=ohe.get feature names(d
             # combine the encoded categorical dataframe and the non categorical data
             X_ohe = pd.concat([ohe_df, non_categorical_df], axis=1)
             # reset the index
             ohe df['ind'] = categorical df.index
             ohe_df = ohe_df.set_index('ind')
             ohe_df.head()
             # combine the encoded categorical dataframe and the non categorical data
             X ohe final = pd.concat([ohe df, non categorical df], axis=1)
             # check the number of columns in train and test sets to ensure they are
             print(len(X_ohe_final.columns))
             return X_ohe_final
In [22]: # encode the train and test data sets
         X_train_ohe = encode_X(X_train, non_categorical_cols)
         X_test_ohe = encode_X(X_test, non_categorical_cols)
```

In [21]: # create a function to perform the one hot encoding of the X data

def encode_X(X, non_categorical_cols):

function to encode a given X data set

Functions for Evaluating Model Performance

882 882 Model evaluation will occur several times throughout this notebook. These functions allow the model perfromance evaluation to be standareized easily applied for each model iteration

4 functions are defined:

- create_y_ohe: encodes the data labels to allow metrics to be calculated for each label (there are 3 labels)
- performance_metrics: calcualtes performance metrics between a data set of test data labels and their predicted values
- evaluate_sbuset_performance: evaluates the performance of each subset of labels
- evaluate_model_performance: evaluate the performance of a model

```
In [23]: tion to perform one hot encoding on the y results
        eate_y_ohe(y, columns):
        nction to encode a given y data set
          Inputs:
               - y: dataframe to encode
              - columns: list of column names included in the encoding
          Outputs:
              - y ohe df: final encoded dataframe
        initialize the encoder
        e = OneHotEncoder()
        create a y dataframe
        df = pd.DataFrame(y, columns=columns)
        encode the dataframe
        ohe = ohe.fit_transform(y_df).toarray()
        converte the encoded y back into a dataframe
        ohe df = pd.DataFrame(y ohe, columns=ohe.get feature names(y df.columns))
        turn y_ohe_df
        tional subset accuracy
        rformance_metrics(y_test, y_pred):
        nction to calcualte performance metrics between a data set of test data labe
          Inputs:
               y_test: known labels for the test data set
              - y_pred: predicted labels for the test data set
          Outputs:
              - results: dictionary containing performance metircs including:
                   accuracy, precision, recall, mean-squared_error, f1, and auc
        calculate metrics
        curacy = accuracy_score(y_test, y_pred)
```

```
ecision = precision_score(y_test, y_pred)
call = recall_score(y_test, y_pred)
e = mean_squared_error(y_test, y_pred)
= f1_score(y_test, y_pred)
lse_positive_rate, true_positive_rate, thresholds = roc_curve(y_test, y_pred)
c_auc = auc(false_positive_rate, true_positive_rate)
organize metrics into a dictionary
sults = {
  'accuracy': round(accuracy, 4),
  'precision': round(precision, 4),
  'recall': round(recall ,4),
  'mse': round(mse, 4),
  'f1': round(f1, 4),
  'auc': round(roc_auc, 4)
turn results
aluate_sbuset_performance(y_test_ohe_df, y_pred_ohe_df):
nction to evaluate the performance of each subset of labels
 Inputs:
      - y_test_ohe_df: encoded dataframe of known labels for the test data se
      - y_pred_ohe_dt: encoded dataframe of predicted labels for the test data
 Outputs:
      - subset_results: dictionary containing performance metircs for each s
          - each set of results within a key named for the subset
              'subset': {
                  accuracy: val,
                  precision: val,
                  recall: val,
                  mse: val,
                  auc: val
              }
define the subsets withing the labels
bsets = ['status_group_functional', 'status_group_functional needs repair',
bset_results = {}
loop through each subset label and create
r subset in subsets:
 # if one subset is not present in the predicted labels given, create a coll
     y_pred = y_pred_ohe_df[subset]
 except KeyError:
      y_pred_ohe_df[subset] = np.zeros(len(y_pred_ohe_df))
 y_test = y_test_ohe_df[subset]
```

```
# calculate the subset results
  subset results[subset] = performance_metrics(y_test, y_pred)
turn subset_results
aluate_model_performance(y_test, y_pred):
nction to evaluate the performance of a model
  Inputs:
      - y test: array of known labels for the test data set
      - y_pred: array of predicted labels for the test data set
 Outputs:
      - results: dictionary containing performance metircs for each subset a
          - each set of results within a key named for the subset
              'subset': {
                  accuracy: val,
                  precision: val,
                  recall: val,
                  mse: val,
                  auc: val
              }
convert the label arrays into encoded dataframes in order to calculate the me
test_ohe_df = create_y_ohe(y_test, columns=['status_group'])
pred_ohe_df = create_y_ohe(y_pred, columns=['status_group'])
calculate the performance for each label subset
bset_results = evaluate_sbuset_performance(y_test_ohe_df, y_pred_ohe_df)
average the subset results together to get an evaluation for the whole model
curacy = []
ecision = []
call = []
e = []
= []
c = []
r subset in subset_results.keys():
 accuracy.append(subset_results[subset]['accuracy'])
 precision.append(subset_results[subset]['precision'])
  recall.append(subset results[subset]['recall'])
 mse.append(subset results[subset]['mse'])
 f1.append(subset results[subset]['f1'])
 auc.append(subset_results[subset]['auc'])
g_accuracy = round(np.array(accuracy).mean(), 4)
g_precision = round(np.array(precision).mean(), 4)
g_recall = round(np.array(recall).mean(), 4)
g mse = round(np.array(mse).mean(), 4)
g_f1 = round(np.array(f1).mean(), 4)
g_auc = round(np.array(auc).mean(), 4)
sults = {
```

10/10/23, 7:00 PM data_analysis - Jupyter Notebook

```
accuracy : avg_accuracy,
  'precision': avg_precision,
  'recall': avg_recall,
  'auc': avg_auc,
  'mse': avg_mse,
  'f1': avg f1,
  'subsets': subset_results
turn results
```

Base Model

Train the tree

- the criterion "gini" was chosen, as it is typically faster than "entropy"
- the same random state is used in the decision tree as the train/test split

```
In [24]: criterion='gini'
         # intialize and fit the tree
         clf_base = DecisionTreeClassifier(criterion=criterion, random_state=SEED)
         clf_base.fit(X_train_ohe, y_train)
```

Out[24]: DecisionTreeClassifier(random_state=20)

Use the tree to predict labels for the test data set

```
In [25]: y_pred_base = clf_base.predict(X_test_ohe)
In [26]: X_test_ohe.head()
```

Out [26]:

	Tunder_U	tunder_Aar	Tunder_Abasia	Tunder_Acra	Tunder_Adb	Tunaer_Aap	Tunder_Adra	tunder
47562	0.0	0.0	0.0	0.0	0.0	0.0	0.0	
29566	0.0	0.0	0.0	0.0	0.0	0.0	0.0	
10873	0.0	0.0	0.0	0.0	0.0	0.0	0.0	
31529	0.0	0.0	0.0	0.0	0.0	0.0	0.0	
5826	0.0	0.0	0.0	0.0	0.0	0.0	0.0	

funder O funder Aar funder Abasia funder Aara funder Adh funder Ada funder Adra funder

5 rows × 882 columns

Evaluate the base model performance

Which value will be our primary metric?

- the goal of this model is to identify which water wells are not functioning as intended, and therefore need to be repaired or fixed.
- We want to avoid identifying a well as functioning when it is not functioning (false positives)
 - We would rather visit a well that is working rather than not visit a well that isn't, assuming we have the man power available
- Therefore, the metric we will be prioritizing is precision, as it prioritizes minimizing false positives rather than false negatives

```
In [27]: results = evaluate_model_performance(y_test.values, y_pred_base)
         print(f'Model 1 Performance - Average')
         print(f' accuracy = {results["accuracy"]}')
         print(f' ** precision = {results["precision"]} **')
         print(f' ** f1 = {results["f1"]} **')
         print(f' recall = {results["recall"]}')
         print(f' auc = {results["auc"]}')
         print(f'Model 1 Performance - Functional')
         print(f' accuracy = {results["subsets"]["status group functional"]["accuracy
         print(f' ** precision = {results["subsets"]["status group functional"]["prec
         print(f' ** f1 = {results["subsets"]["status_group_functional"]["f1"]} **')
         print(f' recall = {results["subsets"]["status_group_functional"]["recall"]}
                   auc = {results["subsets"]["status_group_functional"]["auc"]}')
         print(f'
         Model 1 Performance - Average
           accuracy = 0.8359
          ** precision = 0.6419 **
          ** f1 = 0.6404 **
           recall = 0.639
           auc = 0.7464
         Model 1 Performance - Functional
           accuracy = 0.7809
          ** precision = 0.7947 **
          ** f1 = 0.7998 **
           recall = 0.805
           auc = 0.7786
```

We have an precision of 80% with no model tailoring

Hyperparemeter Tuning

Tune the model by determining the optimal value for 4 model parameters:

- · max features
- min_samples_split
- min_samples_leaf
- max_features

Tune the model based on functional precision

- · this is the precision in identifying whether or not a well is functional
- combines non-functional and functional-needs-repair into one category
 - both pumps will need to be visited
 - not many examples of functional-needs-repair, difficult to train to that metric

Function to perform the tuning

Grid search major - use precision as the metric to define what the best model 4 nested for loops - maybe try another algorighms - use random forest, XG Boost

Include a classification report (call to scikit learn) - compare to my function

• get a confustion matrix in there (communicate false positives)

Feature Importance:

- at the end of the analysis, calculate feature importance of the best performing model (decision tree vs forest, etc)
- areas for the NGO to focus on to repair and help with longevity
- kinda like analyzing coefficients in linear regression

3 recs:

model, then 2 most important features

Function to perform the tuning

Define a function to investigate a range of values for a given parameter on the effect they have on the model.

 This function will be used to perform tuning on the base model and determine the best values for each parameter

```
III [20]. | WEI INVESTIGATE PALAMETER (PALAME, PALAME, PALAMETER)
                                    max_depth=None, min_samples_leaf=1, min_samples_sp
             function to investigate a range of values for a given parameter on the e
                 Inputs:
                     - param name: name of the parameter associated with the range of
                     - param values: range of values to test
                     - max depth: option to set a specific value for the max depth (d
                     - max_samples_leaf: option to set a specific value for the max s
                     - min_samples_split: option to set a specific value for the max
                     - max_features: option to set a specific value for the max featu
             .....
             # initialize metric storage
             train_auc = []
             test_auc = []
             train accuracy = []
             test accuracy = []
             train precision = []
             test precision = []
             train_recall = []
             test_recall = []
             train_f1 = []
             test_f1 = []
             # loop through each value and calculate the performance metrics for the
             for value in param_values:
         #
                   print(value)
                 # select the decision tree that uses the range of values specified
                 if param_name == 'max_depth':
                     dt = DecisionTreeClassifier(criterion=criterion,
                                                  max_depth=value,
                                                  min samples leaf=min samples leaf,
                                                  min_samples_split=min_samples_split,
                                                  max_features=max_features,
                                                  random state=SEED)
                 elif param name == 'min samples leaf':
                     dt = DecisionTreeClassifier(criterion=criterion,
                                                  max depth=max depth,
                                                  min_samples_leaf=value,
                                                  min_samples_split=min_samples_split,
                                                  max_features=max_features,
                                                  random state=SEED)
                 elif param_name == 'min_samples_split':
                     dt = DecisionTreeClassifier(criterion=criterion,
                                                  max depth=max depth,
                                                  min_samples_leaf=min_samples_leaf,
                                                  min samples split=value,
                                                  may features-may features
```

```
man_reacures-man_reacures,
                                    random_state=SEED)
   elif param name == 'max features':
        dt = DecisionTreeClassifier(criterion=criterion,
                                    max_depth=max_depth,
                                    min samples leaf=min samples leaf,
                                    min samples split=min samples split,
                                    max_features=value,
                                    random_state=SEED)
   # fit the tree
   dt.fit(X_train_ohe, y_train)
   # predict the train and test labels
   train pred = dt.predict(X train ohe)
   test_pred = dt.predict(X_test_ohe)
   # evaluate model performance for both the train and test data sets
   test_performance = evaluate_model_performance(y_test, test_pred)
   train performance = evaluate model performance(y train, train pred)
   # accumulate metrics for each parameter value to plot
   train_auc.append(train_performance["subsets"]["status_group_function
   test_auc.append(test_performance["subsets"]["status_group_functional
   train_precision.append(train_performance["subsets"]["status_group_fu
   test_precision.append(test_performance["subsets"]["status_group_fund
   train_recall.append(train_performance["subsets"]["status_group_funct
   test recall.append(test performance["subsets"]["status group function
   train accuracy.append(train_performance["subsets"]["status_group_fun"]
   test_accuracy.append(test_performance["subsets"]["status_group_funct
   train_f1.append(train_performance["subsets"]["status_group_functiona
   test_f1.append(test_performance["subsets"]["status_group_functional"
# determine the parameter value with the hightest precision on the test
max precision = max(test precision)
max value = param values[test precision.index(max precision)]
# plot the auc value for train and test results
plt.figure(figsize=(15, 8))
plt.plot(param_values, train_auc, label='train auc')
plt.plot(param_values, test_auc, label='test auc')
plt.grid()
plt.xlabel(param_name)
plt.vlabel('auc')
plt.legend()
plt.show()
# plot the precision, recall, and accuracy for the test results
plt.figure(figsize=(15, 8))
plt.plot(param_values, test_precision, label='test precision')
plt.plot(param_values, test_recall, label='test recall')
nlt_nlot(naram values test accuracy label='test accuracy')
```

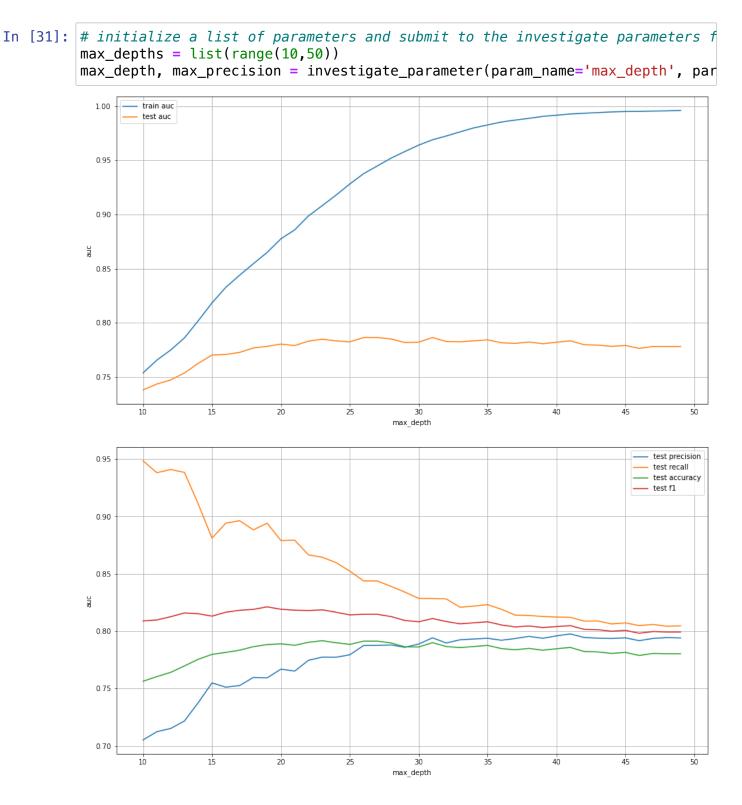
```
plt.plot(param_values, test_f1, label='test f1')
plt.grid()
plt.xlabel(param_name)
plt.ylabel('auc')
plt.legend()
plt.show()
return max_value, max_precision
```

Investigate parameters

Investigate ranges for each parameter.

- select the value that returns the highest precision
- select a range of values including that value to be used in a grid search

Max Depth



Results

```
In [30]: # select the max depth value with the highest precision
    selected_max_depth = max_depth # 19
# define a range of max depth values to in put in the grid search via inspec
    selected_max_depth_range = list(range(30, 46))
    selected_max_depth_range.extend([None]) # include the default value in the
```

```
In [32]: # evaluate the new model and print key metrics
         results2 = evaluate_model_performance(y_test.values, y_pred_2)
         print(f'Model 2 Performance - Average')
         print(f'
                   accuracy = {results2["accuracy"]}')
         print(f'
                   precision = {results2["precision"]}')
         print(f'
                   f1 = {results2["f1"]}')
         print(f' recall = {results2["recall"]}')
                   auc = {results2["auc"]}')
         print(f'
         print(f'Model 2 Performance - Functional')
                   accuracy = {results2["subsets"]["status_group_functional"]["accura
         print(f'
                   precision = {results2["subsets"]["status_group_functional"]["preci
         print(f'
         print(f'
                   f1 = {results2["subsets"]["status_group_functional"]["f1"]}')
                   recall = {results2["subsets"]["status group functional"]["recall"]
         print(f'
                   auc = {results2["subsets"]["status group functional"]["auc"]}')
         print(f'
```

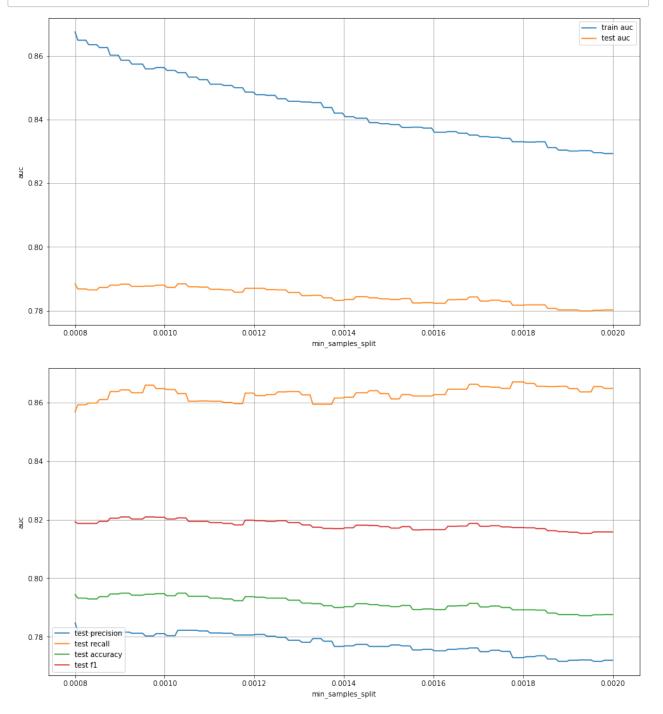
```
Model 2 Performance - Average
accuracy = 0.8421
precision = 0.6535
f1 = 0.6456
recall = 0.6392
auc = 0.7481
Model 2 Performance - Functional
accuracy = 0.79
precision = 0.7942
f1 = 0.811
recall = 0.8285
auc = 0.7863
```

Interpret Results

- precision increased as the max depth increased, until it plateaued around 30.
- the average and function precision values both increased
- the highest precision found with a max depth at 35
- recall started high then fell as max depth increased until it plateaued around 30

Min Samples Split

In [35]: # initialize a list of parameters and submit to the investigate parameters f
min_samples_splits = np.linspace(0.0008, 0.002, 200, endpoint=True)
min_samples_split, max_precision_split = investigate_parameter(param_name='m



Results

```
In [33]: # select the min samples split value with the highest precision
    selected_min_samples_split = min_samples_split # 0.0008
    print(selected_min_samples_split)
    selected_min_samples_split_range = list(np.linspace(0.0008, 0.0014, 7, endposelected_min_samples_split_range.append(2)
    0.0008
In [34]: # create, fit, and predict the model iteration for this parameter tuning
```

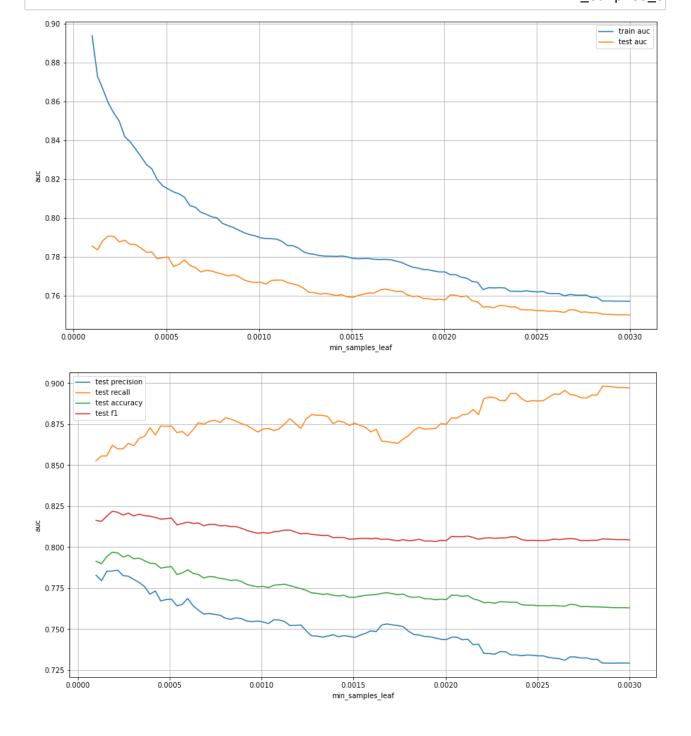
```
In [35]: # evaluate the new model and print key metrics
         results3 = evaluate_model_performance(y_test.values, y_pred_3)
         print(f'Model 3 Performance - Average')
         print(f'
                   accuracy = {results3["accuracy"]}')
                   precision = {results3["precision"]}')
         print(f'
         print(f'
                   precision = {results3["precision"]}')
         print(f'
                   recall = {results3["recall"]}')
                   auc = {results3["auc"]}')
         print(f'
         print(f'Model 3 Performance - Functional')
                   accuracv = {results3["subsets"]["status_group_functional"]["accura
         print(f'
                   precision = {results3["subsets"]["status_group_functional"]["preci
         print(f'
         print(f'
                   f1 = {results3["subsets"]["status_group_functional"]["f1"]}')
         print(f'
                   recall = {results3["subsets"]["status group functional"]["recall"]
                   auc = {results3["subsets"]["status_group_functional"]["auc"]}')
         print(f'
```

```
Model 3 Performance - Average
accuracy = 0.8496
precision = 0.6937
precision = 0.6937
recall = 0.6385
auc = 0.7481
Model 3 Performance - Functional
accuracy = 0.7955
precision = 0.7835
f1 = 0.8209
recall = 0.8621
auc = 0.7891
```

Interpret Results

- there is not a reliable increase throughout the range of min samples split given
 - the optimal value results in an increase in average precision, but not functional precision
- seeing as there is no disntinct increase in precision, it is likely that the default value will be the
 most accurate, however, a range of values will still be submitted to the grid search to investigate
 how they are affected by the other parameters

Min Samples Leaf



Results

```
In [66]: # select the min samples leaf value with the highest precision
         selected min samples leaf = min samples leaf
         selected min samples leaf range = list(np.linspace(13e-5, 26e-5, 14, endpoin
         selected min samples leaf range.append(1)
In [67]: # create, fit, and predict the model iteration for this parameter tuning
         clf4 = DecisionTreeClassifier(criterion=criterion,
                                       max depth=selected max depth,
                                       min_samples_split=selected_min samples split.
                                       min samples leaf=selected min samples leaf,
                                       random_state=SEED)
         clf4.fit(X_train_ohe, y_train)
         y_pred_4 = clf4.predict(X_test_ohe)
In [68]: # evaluate the new model and print key metrics
         results4 = evaluate model performance(y test.values, y pred 4)
         print(f'Model 4 Performance - Average')
         print(f' accuracy = {results4["accuracy"]}')
                   precision = {results4["precision"]}')
         print(f'
         print(f' f1 = {results4["f1"]}')
         print(f' recall = {results4["recall"]}')
                   auc = {results4["auc"]}')
         print(f'
         print(f'Model 4 Performance - Functional')
                   accuracy = {results4["subsets"]["status_group_functional"]["accura
         print(f'
         print(f'
                   precision = {results4["subsets"]["status_group_functional"]["preci
         print(f' f1 = {results4["subsets"]["status_group_functional"]["f1"]}')
                   recall = {results4["subsets"]["status group functional"]["recall"]
         print(f'
                   auc = {results4["subsets"]["status group functional"]["auc"]}')
         print(f'
         Model 4 Performance - Average
           accuracy = 0.851
           precision = 0.7065
           f1 = 0.6552
           recall = 0.6323
           auc = 0.7448
         Model 4 Performance - Functional
           accuracy = 0.7967
           precision = 0.7818
           f1 = 0.8228
           recall = 0.8684
           auc = 0.7898
```

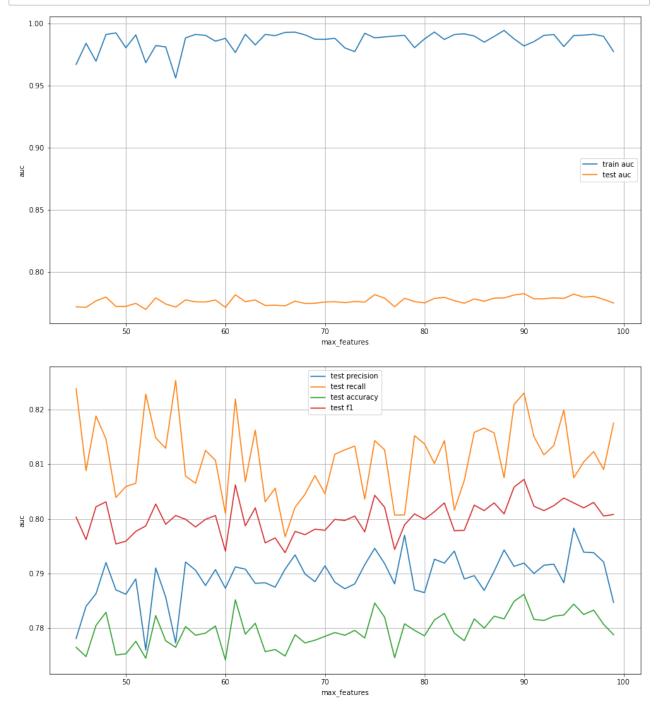
Interpretation

• there is a decrease in precision as min_samples_leaf increases, indicating that smaller values will return higher precision

• with the optimal value selected, average precision increased, while functional precision decreased

Max Features

In [53]: # initialize a list of parameters and submit to the investigate parameters f
max_features = list(range(45, 100))
max_feature, max_precision_features = investigate_parameter(param_name='max_
max_depth=selec



Results

```
In [39]: # select the max features value with the highest precision
         selected_max_features = max_feature
         selected_max_features_range = list(range(73, 100))
         selected max features range.extend([None])
In [40]: # create, fit, and predict the model iteration for this parameter tuning
         clf5 = DecisionTreeClassifier(criterion=criterion,
                                       max depth=selected max depth,
                                       max_features=selected_max_features,
                                       random state=SEED)
         clf5.fit(X_train_ohe, y_train)
         y_pred_5 = clf5.predict(X_test_ohe)
In [41]: | # evaluate the new model and print key metrics
         results5 = evaluate_model_performance(y_test.values, y_pred_5)
         print(f'Model 5 Performance - Average')
                   accuracy = {results5["accuracy"]}')
         print(f'
         print(f'
                   precision = {results5["precision"]}')
         print(f' f1 = {results5["f1"]}')
         print(f' recall = {results5["recall"]}')
         print(f' auc = {results5["auc"]}')
         print(f'Model 5 Performance - Functional')
         print(f'
                   accuracy = {results5["subsets"]["status_group_functional"]["accura
                   precision = {results5["subsets"]["status_group_functional"]["preci
         print(f'
         print(f'
                   f1 = {results5["subsets"]["status_group_functional"]["f1"]}')
         print(f'
                   recall = {results5["subsets"]["status_group_functional"]["recall"]
                   auc = {results5["subsets"]["status group functional"]["auc"]}')
         print(f'
         Model 5 Performance - Average
           accuracy = 0.8388
           precision = 0.6541
           f1 = 0.6481
           recall = 0.6432
           auc = 0.7485
         Model 5 Performance - Functional
           accuracy = 0.783
           precision = 0.7867
           f1 = 0.8051
           recall = 0.8244
           auc = 0.779
```

Interpret Results

- there is a large difference between precision as max features increase without a clear trend.
- the optimal value selection seems more random rather than a distinct peak.

Grid Search

Using the ranges of values selected in the inital parameter tuning iterations, iterate through all possible combinations in a search for the combination of paramters that results in the highest functional precision.

```
In [47]:
        hat iterates between all possible parameter combinations
        les_leafs, max_features):
        keep track of progress
        urrent min samples split to keep track of progress
        irrent combination of paramters
        n=criterion,
        pth,
         =min_samples_leaf,
         t=min_samples_split,
        _feature,
        ED)
        ie)
        rformance(y_test.values, test_pred)
         _performance(y_train, train_pred)
        lance)
        berformance['subsets']['status_group_functional']['precision']
        end(functional_binary_precision)
        st_performance['subsets']['status_group_non functional']['precision']
        l.append(non_functional_binary_precision)
```

```
['precision']
je_precision)
lit, min_samples_leaf, max_feature]
rerage precision
lsion grid.index(max avg precision)]
nctional precision
nary precision grid)
[[functional_binary_precision_grid.index(max_functional_binary_precision)]
n-functional precision
lonal_binary_precision_grid)
grid[non_functional_binary_precision_grid.index(max_non_functional_binary_pr
 2)}%: params = {max avg precision params}')
binary_precision*100, 2)}%: params = {max_functional_binary_precision_params
lctional_binary_precision*100, 2)}%: params = {max_non_functional_binary_prec
rams,
l_binary_precision,
inctional_binary_precision_params,
functional_binary_precision,
x_non_functional_binary_precision_params,
lected_min_samples_split_range, selected_min_samples_leaf_range, selected_max
30
    0.0008
    0.0009
    0.001
    0.0011
    0.00120000000000000001
    0.0013
    0.0014
    2
31
    0.0008
    0.0009
    a a a a 1
```

```
0 . UU T
    0.0011
    0.0012000000000000001
    0.0013
    0.0014
    2
32
    0.0008
    0.0009
    0.001
    0.0011
    0.0012000000000000001
    0.0013
    0.0014
    2
33
    0.0008
    0.0009
    0.001
    0.0011
    0.0012000000000000001
    0.0013
    0.0014
    2
34
    0.0008
    0.0009
    0.001
    0.0011
    0.0012000000000000001
    0.0013
    0.0014
    2
35
    0.0008
    0.0009
    0.001
    0.0011
    0.00120000000000000001
    0.0013
    0.0014
    2
36
    0.0008
    0.0009
    0.001
    0.0011
    0.00120000000000000001
    0.0013
    0.0014
    2
37
    0.0008
    0.0009
    a aa1
```

```
0 I UU I
    0.0011
    0.00120000000000000001
    0.0013
    0.0014
    2
38
    0.0008
    0.0009
    0.001
    0.0011
    0.00120000000000000001
    0.0013
    0.0014
    2
39
    0.0008
    0.0009
    0.001
    0.0011
    0.0012000000000000001
    0.0013
    0.0014
    2
40
    0.0008
    0.0009
    0.001
    0.0011
    0.0012000000000000001
    0.0013
    0.0014
    2
41
    0.0008
    0.0009
    0.001
    0.0011
    0.00120000000000000001
    0.0013
    0.0014
    2
42
    0.0008
    0.0009
    0.001
    0.0011
    0.00120000000000000001
    0.0013
    0.0014
    2
43
    0.0008
    0.0009
    a aa1
```

```
0 . UUT
   0.0011
   0.00120000000000000001
   0.0013
   0.0014
   2
44
   0.0008
   0.0009
   0.001
   0.0011
   0.0012000000000000001
   0.0013
   0.0014
   2
45
   0.0008
   0.0009
   0.001
   0.0011
   0.00120000000000000001
   0.0013
   0.0014
   2
None
   0.0008
   0.0009
   0.001
   0.0011
   0.00120000000000000001
   0.0013
   0.0014
   2
avg precision: 72.27%: params = [38, 0.0014, 0.000209999999999999, 95]
functional precision: 80.07%: params = [42, 2, 1, 89]
91]
```

Interpret Results

- prior to hyperparameter tuning, the model functional precision was 80.25%.
- after hyperparameter tuning, the model functional precision is 80.69%
- there is an increase in functional precision of 0.44%

Final Model - Decision Tree

Run the final model chosen to maximize functional precision

```
In [48]: # import results
         with open("data/grid_results_precision.json", "r") as f:
             grid results = ison.load(f)
In [49]: # select the final model parameters
         chosen_final_model_params = grid_results['max_functional_binary_precision_pa'
         final_max_depth = chosen_final_model_params[0]
         final_min_samples_split = chosen_final_model_params[1]
         final min samples leaf = chosen final model params[2]
         final max features = chosen final model params[3]
In [50]: # define and train a the decision tree, and predict the test labels
         clf final = DecisionTreeClassifier(criterion=criterion,
                                            max depth=final max depth,
                                            min_samples_split=final_min_samples_split
                                            min samples leaf=final min samples leaf,
                                            max_features=final_max_features,
                                            random_state=SEED)
         clf_final.fit(X_train_ohe, y_train)
         y_pred_final = clf_final.predict(X_test_ohe)
In [51]: final results = evaluate model performance(y test.values, y pred final)
         print(f'Final Model Performance - Average')
         print(f'
                   accuracy = {final results["accuracy"]}')
         print(f'
                   precision = {final_results["precision"]}')
         print(f'
                   f1 = {final_results["f1"]}')
         print(f' recall = {final results["recall"]}')
                   auc = {final results["auc"]}')
         print(f'
         print(f'Final Model Performance - Functional')
                   accuracy = {final_results["subsets"]["status_group_functional"]["a
         print(f'
         print(f'
                   precision = {final results["subsets"]["status group functional"]["
                   f1 = {final_results["subsets"]["status_group_functional"]["f1"]}')
         print(f'
                   recall = {final results["subsets"]["status group functional"]["rec
         print(f'
                   auc = {final results["subsets"]["status group functional"]["auc"]}
         print(f'
         Final Model Performance - Average
           accuracy = 0.8395
           precision = 0.6531
           f1 = 0.6514
           recall = 0.6499
           auc = 0.7534
         Final Model Performance - Functional
           accuracy = 0.7878
           precision = 0.8007
           f1 = 0.8061
           recall = 0.8117
           auc = 0.7855
```

Random Forest Using Final Decision Tree Parameters

Input the final decision tree parameters into a random forest.

```
In [52]: # define random forest parameters
         n estimators=100
         n_jobs=None
         # define the random forest and train it
         forest = RandomForestClassifier(n_estimators=n_estimators, criterion=criteri
                                        max depth=final max depth, max features=final
                                        min samples split=final min samples split, mi
                                        n jobs=n jobs, random state=SEED)
         forest.fit(X_train_ohe, y_train)
         # use the forest to predict the test labels
         y_forest = forest.predict(X_test_ohe)
In [53]: forest_results = evaluate_model_performance(y_test.values, y_forest)
         print(f'Final Model Performance - Average')
                   accuracy = {forest results["accuracy"]}')
         print(f'
         print(f'
                   precision = {forest_results["precision"]}')
         print(f'
                   f1 = {forest results["f1"]}')
         print(f' recall = {forest_results["recall"]}')
         print(f'
                   auc = {forest results["auc"]}')
         print(f'Final Model Performance - Functional')
         print(f'
                   accuracy = {forest_results["subsets"]["status_group_functional"]["
         print(f'
                   precision = {forest results["subsets"]["status group functional"][
                   f1 = {forest_results["subsets"]["status_group_functional"]["f1"]}'
         print(f'
                   recall = {forest results["subsets"]["status group functional"]["re
         print(f'
         print(f'
                   auc = {forest_results["subsets"]["status_group_functional"]["auc"]
         Final Model Performance - Average
           accuracy = 0.8686
           precision = 0.7221
           f1 = 0.6893
           recall = 0.6698
           auc = 0.7727
         Final Model Performance - Functional
           accuracy = 0.8204
           precision = 0.805
           f1 = 0.8426
           recall = 0.8839
           auc = 0.8143
```

Interpretaion

- after hyperparameter tuning, the model functional precision is 80.69%
- using the same optimized hyperparameters, the model functional precision is increased to 80.86%
- there is an increase in functional precision of 0.61% from the original base model

Model Improvements

```
In [69]: # functional

precision_increase = forest_results["subsets"]["status_group_functional"]["pr
f1_increase = forest_results["subsets"]["status_group_functional"]["f1"] - re
recall_increase = forest_results["subsets"]["status_group_functional"]["recal
accuracy_increase = forest_results["subsets"]["status_group_functional"]["acc

print(f'increase in precision from base model to final model: {round(precision)
print(f'increase in f1 score from base model to final model: {round(f1_increan)
print(f'increase in recall from base model to final model: {round(accuracy)

increase in precision from base model to final model: 1.03%
increase in f1 score from base model to final model: 4.28%
increase in recall from base model to final model: 7.89%
increase in accuracy from base model to final model: 3.95%
```

The optimized model generated an increase in all major metrics, with the largest increae in recall.

```
In [55]: # final model classification report
final_report = classification_report(y_test, y_forest)
print(final_report)
```

	precision	recall	f1–score	support
functional functional needs repair non functional	0.80 0.52 0.84	0.88 0.35 0.78	0.84 0.42 0.81	7242 984 5094
accuracy macro avg weighted avg	0.72 0.80	0.67 0.80	0.80 0.69 0.80	13320 13320 13320

Confusion Matrix

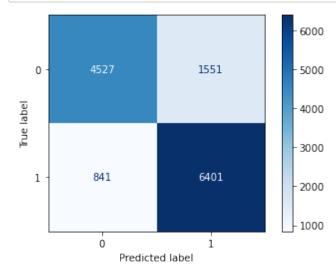
Generate a confusion matrix to illustrate the numbers of true postives, true negaties, false positives, and fasle negatives

Final Model

```
In [56]: y_test_ohe_df = create_y_ohe(y_test, columns=['status_group'])
    y_pred_ohe_df = create_y_ohe(y_forest, columns=['status_group'])

y_test_functional = y_test_ohe_df['status_group_functional'].values
    y_pred_functional = y_pred_ohe_df['status_group_functional'].values

cfm = confusion_matrix(y_test_functional, y_pred_functional)
    disp = ConfusionMatrixDisplay(confusion_matrix=cfm)
    disp.plot(cmap=plt.cm.Blues)
    plt.show()
```

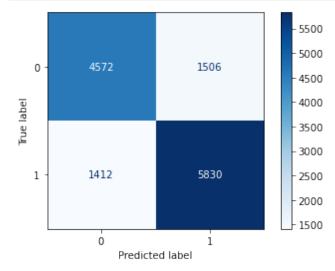


Base Model

```
In [57]: y_test_ohe_df = create_y_ohe(y_test, columns=['status_group'])
y_pred_ohe_df = create_y_ohe(y_pred_base, columns=['status_group'])

y_test_functional = y_test_ohe_df['status_group_functional'].values
y_pred_functional = y_pred_ohe_df['status_group_functional'].values

cfm = confusion_matrix(y_test_functional, y_pred_functional)
disp = ConfusionMatrixDisplay(confusion_matrix=cfm)
disp.plot(cmap=plt.cm.Blues)
plt.show()
```



Interpretation

Our goal was to minimize the number of false positives by using precision as out primary metric, while the number of false positives increased, the number of true positives increased at a greater rate, which in turn raised the precision by 1%.

Random Forest Feature Importance

Investigate the features to determine the which characteristics of the pumps have the largest effect on their functionality.

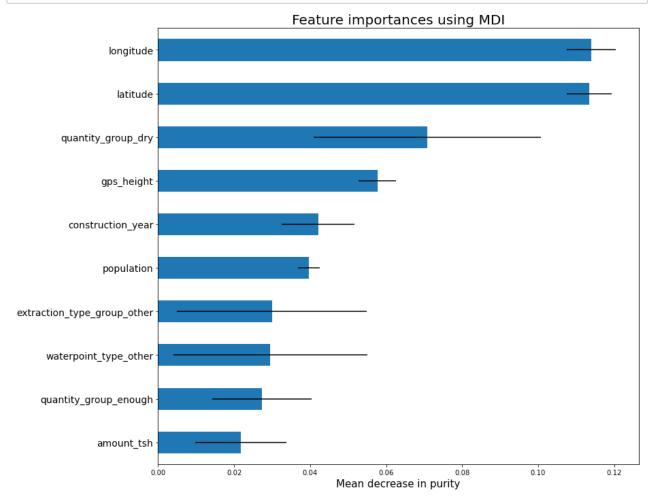
```
In [58]: # define the importances and their standard deivations
importances = forest.feature_importances_
std = np.std([tree.feature_importances_ for tree in forest.estimators_], axi
```

```
In [59]: # define feature names, importances, and their standard deviations
    feature_names = np.array(X_train_ohe.columns)
    importances = forest.feature_importances_
    std = np.std([tree.feature_importances_ for tree in forest.estimators_], axi

# create a data frame for all the importances
data={
      'feature': feature_names,
      'importance': importances,
      'std': std
}
forest_importances = pd.DataFrame(data)

# filter the importances to only include the top features, then sort
forest_importances_filtered = forest_importances[forest_importances['importancestation]
forest_importances_filtered = forest_importances_filtered.set_index('feature forest_importances_filtered = forest_importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='importances_filtered.sort_values(by='import
```

```
In [60]: # plot feature importances
    fig, ax = plt.subplots(figsize=(13,10))
    forest_importances_filtered['importance'].plot.barh(xerr=forest_importances_f
    ax.set_title("Feature importances using MDI", fontsize=20)
    ax.set_xlabel("Mean decrease in purity", fontsize=15)
    ax.set_ylabel('', )
    plt.yticks(fontsize=14)
    fig.tight_layout()
```



Interpretaion

The top features that are most important in determining if a water pump is functional:

- 1. location (latitude/longitude/elevation)
- 2. water quantity
- 3. construction year

As we develop a plan to make maintenance visits to water pumps, we should prioritize the pumps by location first, identifying regions with higher rates of non-functioning pumps. After that, focus on older pumps with "dry" water quantities.