Proj_01: Finding Donors for CharityML

Source: UdaCity, Machine Learning Intro, Supervided Learning

Introduction

In this project, I will employ several supervised algorithms to accurately model indivisuals' income using data collected from the 1944 U.S. Census.

I will then chose the best candidate algorithm from preliminary results and further optimize this algorithm to best model the data.

Goal

The goal of this project is to construct a model that accurately predicts whether an individual makes more than \$50,000.

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1. Exploring the Dataset

```
In [1]: # Import libraries
   import numpy as np
   import pandas as pd
   from time import time
   from IPython.display import display # this allows the use of display() for DataFrames
   # import visuals as vs #
   import matplotlib.pyplot as plt
%matplotlib inline
```

1-1. Overview the data

- The dataset for this project originates from the UCI Machine Learning Repository.
- From the table below, I found that there are numerical and string features.

```
In [2]: # Load the Census dataset
data = pd.read_csv("./Data/Proj_01/census.csv")

# Display the first 5 records
display(data.head(5))

# Display the unique income values
print("The unique values of INCOME")
display(data["income"].unique())
```

	age	workclass	education_level	education- num	marital- status	occupation	relationship	race	sex	capital- gain	capita lo
0	39	State-gov	Bachelors	13.0	Never- married	Adm- clerical	Not-in- family	White	Male	2174.0	С
1	50	Self-emp- not-inc	Bachelors	13.0	Married- civ- spouse	Exec- managerial	Husband	White	Male	0.0	С
2	38	Private	HS-grad	9.0	Divorced	Handlers- cleaners	Not-in- family	White	Male	0.0	С
3	53	Private	11th	7.0	Married- civ- spouse	Handlers- cleaners	Husband	Black	Male	0.0	С
4	28	Private	Bachelors	13.0	Married- civ- spouse	Prof- specialty	Wife	Black	Female	0.0	С

1-2. Check invalid and missing data

array(['<=50K', '>50K'], dtype=object)

• There are no invalid data.

The unique values of INCOME

```
In [3]: data.isnull().sum(axis=0)
```

Out[3]: age

```
workclass 0
education_level 0
education-num 0
marital-status 0
occupation 0
relationship 0
race 0
sex 0
capital-gain 0
capital-loss 0
hours-per-week 0
native-country 1
income 0
dtype: int64
```

1-3. Data exploration: income

A cursory investigation of the dataset will determine how many individuals fit into either group, and will tell us about the percentage of these individuals making more than \$50,000.

In the code cell below, I will compute the following:

- The total number of records : n_records
- The number of individuals making more than \$50,000 annually n_greater_50k
- The number of individuals making at most \$50,000 annually: n_at_most_50k
- The percentage of individuals making more than \$50,000 annually: greater_percent

```
In [4]: n_records
                      = data.shape[0]
        n greater 50k = (data["income"] == ">50K").sum()
       n at most 50k = (data["income"] == "<=50K").sum()
        greater percent = n greater 50k / n records * 100
        # Print results
       print("Total number of records
                                                                             : {}".format(n rec
                                                                             : {}".format(n gre
       print ("The number of individuals making more than $50,000 annually
       print("The number of individuals making at most $50,000 annually : {}".format(n at
       print("The percentage of individuals making more than $50,000 annually: {0:5.1f} %".form
       Total number of records
                                                                     : 45222
       The number of individuals making more than $50,000 annually
                                                                     : 11208
       The number of individuals making at most $50,000 annually
       The percentage of individuals making more than $50,000 annually: 24.8 %
```

1-4. Data exploration: numerical data

- From the table and the graph below, capital-gain and capital-loss are highly skewed.
- So, some transformation are needed for those features.

```
In [5]: col_names_numerical = ["age", "education-num", "capital-gain", "capital-loss", "hours-pe
    data_numerical = data.loc[:, col_names_numerical]
    data_numerical.describe()
```

Out[5]: age education-num capital-gain capital-loss hours-per-week **count** 45222.000000 45222.000000 45222.000000 45222.000000 45222.000000 38.547941 10.118460 1101.430344 88.595418 40.938017 mean 13.217870 2.552881 7506.430084 404.956092 12.007508 17.000000 1.000000 0.000000 0.000000 1.000000 min

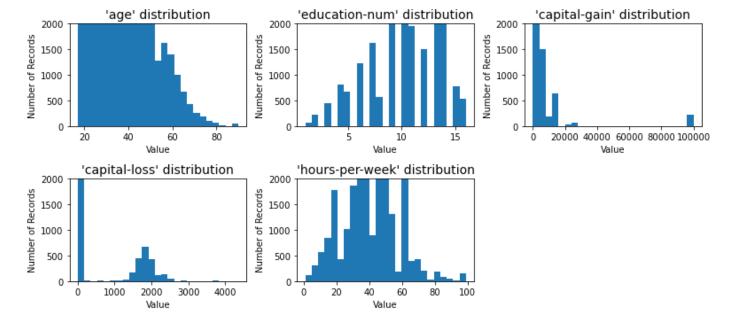
25%	28.000000	9.000000	0.000000	0.000000	40.000000
50%	37.000000	10.000000	0.000000	0.000000	40.000000
75%	47.000000	13.000000	0.000000	0.000000	45.000000
max	90.000000	16.000000	99999.000000	4356.000000	99.000000

```
In [6]: def plot_hist(df, col_names, plot_row_num=2, plot_col_num=3, y_limit=(0,2000)):
    # create figure
    fig = plt.figure(figsize = (11,5));

# plot in bin
for i, feature in enumerate(col_names):
        ax = fig.add_subplot(plot_row_num, plot_col_num, i+1)
        ax.hist(df[feature], bins=25)
        ax.set_title("'%s' distribution"%(feature), fontsize=14)
        ax.set_xlabel("Value")
        ax.set_ylabel("Number of Records")
        ax.set_ylim(y_limit)

# fig.tight_layout()
```

```
In [7]: #plot_hist(data, ["capital-gain", "capital-loss"])
plot_hist(data, col_names_numerical)
```



1-5. Data exploration: string data(other than income)

- I show the unique values of string data.
- A conversion from string to numerical data is needed.

```
for j, unique_name in enumerate(unique_names):
    if j == 0:
        output_name = unique_name
    else:
        output_name = output_name + ", " + unique_name

    print(output_name)

0. [sex]: unique value num = 2
Male, Female

1. [education_level]: unique value num = 16
Bachelors, HS-grad, 11th, Masters, 9th, Some-college, Assoc-acdm, 7th-8th, Doct orate, Assoc-voc, Prof-school, 5th-6th, 10th, Preschool, 12th, 1st-4th

2. [workclass]: unique value num = 7
```

State-qov, Self-emp-not-inc, Private, Federal-qov, Local-qov, Self-emp-inc, Witho

ut-pay

and-Netherlands

3. [relationship] : unique value num = 6
Not-in-family, Husband, Wife, Own-child, Unmarried, Other-relative

4. [occupation] : unique value num = 14
Adm-clerical, Exec-managerial, Handlers-cleaners, Prof-specialty, Other-service, S
ales, Transport-moving, Farming-fishing, Machine-op-inspct, Tech-support, Craft-rep
air, Protective-serv, Armed-Forces, Priv-house-serv

5. [marital-status] : unique value num = 7
Never-married, Married-civ-spouse, Divorced, Married-spouse-absent, Separated, Married-AF-spouse, Widowed

6. [race] : unique value num = 5
White, Black, Asian-Pac-Islander, Amer-Indian-Eskimo, Other

7. [native-country] : unique value num = 41
United-States, Cuba, Jamaica, India, Mexico, Puerto-Rico, Honduras, England, Ca
nada, Germany, Iran, Philippines, Poland, Columbia, Cambodia, Thailand, Ecuador,
Laos, Taiwan, Haiti, Portugal, Dominican-Republic, El-Salvador, France, Guatema
la, Italy, China, South, Japan, Yugoslavia, Peru, Outlying-US(Guam-USVI-etc), Sc
otland, Trinadad&Tobago, Greece, Nicaragua, Vietnam, Hong, Ireland, Hungary, Hol

2. Preprocessing the Dataset

- Fortunately, for this dataset, there are no invalid or missing entries we must deal with.
- However, there are some qualities about certain features that must be adjusted.
- This preprocessing can help tremendously with the outcome and predictive power of nearly all learning algorithms.

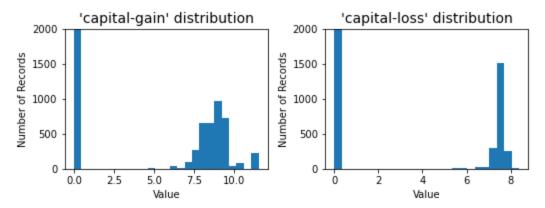
```
In [9]: # Split the dataset into "target label" and "features"
  income_raw = data["income"]
  features_raw = data.drop("income", axis=1)
```

2-1. Transformation: skewed continuous data

• Algorithms can be sensitive hightly skewed feature distribution such as capital-gain and capital-loss, and underperform if the range is not properly normalized.

- I will apply logarithmic transformation) on the data, so that the very large and very small values do not genatively affect the performance of a learning algorithm.
- Using a logarithmic transformation significantly reduces the range of values caused by outliers.
 - The logarighm of 0 is undefined, so we must translate the values by a small amount above 0 to apply the logarithm successfully.

	capital-gain	capital-loss
count	45222.000000	45222.000000
mean	0.740759	0.355489
std	2.466527	1.595914
min	0.000000	0.000000
25%	0.000000	0.000000
50%	0.000000	0.000000
75%	0.000000	0.000000
max	11.512925	8.379539



- From the table above, the data range becomes smaller.
 - capital-gain : 99999.0 --> 11.5
 - capital-loss:4356.0 --> 8.3

2-2. Normalization: numerical data

- Normalization enusures that each feature is treated equally when applying supervised learners.
- I will use sklearn.preprocessing.MinMaxScaler for this.

```
features_log_minmax_transform = pd.DataFrame(data=features_log_transformed)
features_log_minmax_transform[col_names_numerical] = scaler.fit_transform(features_log_t
# Show results
display(features_log_minmax_transform[col_names_numerical].describe())
```

	age	education-num	capital-gain	capital-loss	hours-per-week
count	45222.000000	45222.000000	45222.000000	45222.000000	45222.000000
mean	0.295177	0.607897	0.064342	0.042423	0.407531
std	0.181067	0.170192	0.214240	0.190454	0.122526
min	0.000000	0.000000	0.000000	0.000000	0.000000
25%	0.150685	0.533333	0.000000	0.000000	0.397959
50%	0.273973	0.600000	0.000000	0.000000	0.397959
75%	0.410959	0.800000	0.000000	0.000000	0.448980
max	1.000000	1.000000	1.000000	1.000000	1.000000

2-3. Normalization: non-numerical data

- There are several features for each record that are non-numeric.
- Typically, learning algorithms expect input to be numeric.
- I will use one-hot encoding to convert categorical features.
 - One-hot encoding creates a dummy cariable for each possible category of each non-numerical feature.
 - For exapmle, assume someFeature has three possible entiries: A , B or C .
 - We then endode this feature into someFeatureA, someFeatureB and someFeatureC.
- I will use pandas.get_dummies() to perform one-hot encoding on the feature_log_minmax_transform data.

	someFeature		someFeature_A	someFeature_B	someFeature_C
0	В		0	1	0
1	С	> one-hot encode>	0	0	1
2	А		1	0	0

- Additionally, I will convert the non-numerical target label income to numerical values.
- Since there are only two possible categories for this label(<=50K and >50K), we can avoid using one-hot encoding and simply encode these two categories as 0 and 1.
 - I set records with <=50K to 0 and record with >50K to 1.

```
In [13]: # One-hot encoding : features
  features_final = pd.get_dummies(features_log_minmax_transform)
  display(features_final.describe())
```

count	45222.000000	45222.000000	45222.000000	45222.000000	45222.000000	45222.000000	45222.000000	45222.0
mean	0.295177	0.607897	0.064342	0.042423	0.407531	0.031091	0.068551	0.7
std	0.181067	0.170192	0.214240	0.190454	0.122526	0.173566	0.252691	0.4
min	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.000000	0.0
25%	0.150685	0.533333	0.000000	0.000000	0.397959	0.000000	0.000000	0.0
50%	0.273973	0.600000	0.000000	0.000000	0.397959	0.000000	0.000000	1.0
75%	0.410959	0.800000	0.000000	0.000000	0.448980	0.000000	0.000000	1.0
max	1.000000	1.000000	1.000000	1.000000	1.000000	1.000000	1.000000	1.0

8 rows × 103 columns

```
In [14]: # One-hot encoding : target label
        income = income raw.map(\{"<=50K":0, ">50K":1\})
        display(income.describe())
        count 45222.000000
       mean
                  0.247844
                   0.431766
        std
                  0.000000
       min
        25%
                  0.000000
                  0.000000
        50%
        75%
                   0.000000
       max
                  1.000000
        Name: income, dtype: float64
```

2-4. Shuffle and Split Data

Testing set has 9045 samples.

Now all *categorical features* have been converted into numerical features, and all numerical features have been normalized. As always, I will now split the data into training and test sets.

80% of the data will be used for training and 20% for testing.

```
In [15]: # Import train_test_split
    from sklearn.model_selection import train_test_split

# Spline the features and the label into training and testing sets
    x_train, x_test, y_train, y_test = train_test_split(features_final, income, test_size =

# Show resutls
    print("Training set has {} samples.".format(x_train.shape[0]))
    print("Testing set has {} samples.".format(x_test.shape[0]))

Training set has 36177 samples.
```

3. Modeling

In this section, I will investigate four different algorithms and determine which is best at modeling the data. Three of these algorithms will be supervised learners, and the fourth algorithm is known as a predictor.

3-1. Metrics

CharityML, equipped with theire research, knows individuals that make more than \$50,000 are most likely to donate to their charity. Because of this, *CharityML* is particularly interested in predicting who makes more than \$50,000 accurately.

It would seem that using **accuracy** as a metric for evaluating a particular model's performance would be appropriate.

Additionally, identifying someone that does not make more than \$50,000 as someone who does would be detrimental to *CharityML*, since they are looking to find individuals willing to donate.

Therefore, a model's ability to precisely predict those that make more than \$50,000 is *more important* than the model's ability to recall those individuals.

I can use **F-beta score** as a metric that considers both precision and recall:

$$F_{eta} = (1 + eta^2) \cdot rac{precision \cdot recall}{(eta^2 \cdot precision) + recall}$$

In particular, when $\beta=0.5$, more emphasis is placed on precision. This is called the $F_{0.5}score$ or (F-score for simplicity).

$$egin{aligned} F_1 = (1+1^2) \cdot rac{precision \cdot recall}{\left(1^2 \cdot precision
ight) + recall} &= rac{2 \cdot precision \cdot recall}{precision + recall} \ F_0 = (1+0^2) \cdot rac{precision \cdot recall}{\left(0^2 \cdot precision
ight) + recall} &= precision \end{aligned}$$

3-2. Recap of Metrics

		Prediction Value			
		Yes	No		
American Volus	Yes	TP: True Positive	FN: False Negative		
Answer Value	No	FP: False Positive	TN: True Negative		

Accuracy

- Accurach = TP / (TP + FP + TN + FN)
- Accracy mesures how often the classifier make the correct prediction.
- It's the ratio of the number of correct predictions to the total number of predictions (the number of test data points).

Precision

- Precision = TP / (TP + FP)
- Precision tells us what proportion of messages we classified as spam, actually were spam.
- It is a ratio of true positives to all positives.
- If we want to reduce the ration of FP, we ues precision.
- e.g.
 - true positives : words classified as spam, and which are accually spam

all positives : all words classified as spam, irrespective of whether that was the correct classification

Recall (sensitivity)

- Recall = TP / (TP + FN)
- Recall tells us what proportion of messages that actually were spam were classified by us as spam.
- It is a ration of true positives to all the words that were actually spam.
- If we want to reduce the ration of FN, we use recall. (e.g. cancer detection)

3-3. Naive Predictor Performace

- I create a naive predictor which always predicts 1 (i.e. the individuals makes more than 50k).
- I will evaluate this model's accuracy and F-score.
- The purpose of generating a native predictor is simply to show what a base model without any intelligence would look like.
- This model will have no TN(true negative) or FN(false negative).

```
In [16]: TP = np.sum(income)
        FP = income.shape[0] - TP
        TN = 0
        FN = 0
        accuracy = TP / (TP+FP+TN+FN)
        precision = TP / (TP+FP)
        recall = TP / (TP+FN)
        beta
                  = 0.5
        fscore = (1+beta**2)*(precision*recall)/(beta**2*precision+recall)
        # Show Results
        print("Naive Predictor Performance")
        print(" Accuracy = {:.4f}".format(accuracy))
        print(" F-score = {:.4f}".format(fscore))
        Naive Predictor Performance
           Accuracy = 0.2478
           F-score = 0.2917
```

3-4. Select Supervised Learning Models

I choose three models which are appropriate for this model. I will explain the reasons about the four points below.

- 1. Describe one real-world application in industry where the model can be applied.
- 2. What are the strengths of the model; when does it perform well?
- 3. What are the weaknesses of the model; when does it perform poorly?
- 4. What makes this model a good candidate for the problem, given what you know about the data?

1:Gaussian Naive Bayes

- 1. Real time forecast of stock prices
- 2. Training time and prediction time are small. We use it when features are independent.

- 3. When features are not independent, accuracy will become bad.
- 4. For 1st trial, I use this. I want to find quickly whether features are independent.

2:Decision Tree

- 1. Optimization of the arrangement of workers according to the weather forecast
- 2. By using this, we can easily visualize the result.
- 3. This tends to overfit the data.
- 4. To get good understanding of the data visually.

3:AdaBoost

- 1. classification of image data
- 2. To get better accuracy, this method can select the features which mostly contribute to the true prediction.
- 3. This tends to overfit the data.
- 4. To use this, we can find the necessary features, so we will be able to do better preprocessings.

3-5. Create a Training and Predicting Pipeline

To properly evaluate the performance of each model you've chosen, it's important that I create a training and predicting pipeline that allows me to quickly and effectively train models using various sizes of training data and perform predictions on the testing data.

I will implement the following:

- Import fbeta_score and accuracy_score from sklearn.metrics
- Fit the learner to the sampled training data and record the training time.
- Perform predictions on the test data X_test, and also on the first 300 training points
 X_train[:300]
 - Record the total prediction time.
- Calculate the accuracy score for both the training subset and testing subset
- Calculate the F-score for both the training subset and testing subset

```
results['train_time'] = t_end - t_start

# Prediction
t_start = time()
predictions_train = learner.predict(x_train.iloc[:300])
predictions_test = learner.predict(x_test)
t_end = time()
results['pred_time'] = t_end - t_start
beta = 0.5

# Evaluation : train
results['train_acc'] = accuracy_score(y_train[:300], predictions_train[:300])
results['train_f'] = fbeta_score( y_train[:300], predictions_train[:300], beta=b

# Evaluation : test
results['test_acc'] = accuracy_score(y_test, predictions_test)
results['test_f'] = fbeta_score( y_test, predictions_test, beta=beta)

return results
```

I implement a plot function to evaluate.

```
In [18]: import matplotlib.patches as mpatches
                     def plot evaluate(results, bar width=0.3, figsize=(11,7)):
                              Visualization code to display results of learners.
                              inputs:
                                 - results: a map. results[model name][sample num id][metric]
                              # Create a figure
                              fig, ax = plt.subplots(2,3,figsize=figsize)
                              # Parameters
                              model colors = ["tab:blue", "tab:orange", "tab:green", "tab:red", "tab:purple", \
                                                                  "tab:brown", "tab:pink", "tab:gray", "tab:olive", "tab:cyan"]
                              metric_names = ["train_time", "train_acc", "train_f", "pred_time", "test_acc", "test_acc", "test_acc", "train_f", "pred_time", "test_acc", "test_acc", "train_f", "pred_time", "test_acc", "test_acc", "train_f", "pred_time", "test_acc", "test_
                              ylabels = ["Time[s]", "Accuracy", "F-score", "Time[s]", "Accuracy", "F-sc
                              titles = ["Training:Time", "Training:Accuracy", "Training:F-score", "Pred:Time
                              model names = results.keys()
                              num sampling type = 3;
                              # Loops for bar plot
                              for model id, model name in enumerate(model names):
                                       model color = model colors[model id]
                                       for metric id, metric name in enumerate(metric names):
                                                 plot row = metric id // 3 #//len(model names); # shou
                                                 plot col = metric id % 3 #len(model names); # amari
                                                 tmp ax = ax[plot row, plot col]
                                                 for sample id in range (num sampling type):
                                                          x = sample id + model id * bar width;
                                                          y = results[model name][sample id][metric name]
                                                          tmp ax.bar(x,y, width=bar width, color=model color)
                                                          tmp ax.set xticks([0.45, 1.45, 2.45])
                                                          tmp ax.set xticklabels(["1%", "10%", "100%"])
                                                          tmp ax.set xlabel("Training Set Size")
                                                           tmp ax.set xlim((-0.1, 3.0))
                                                           tmp ax.set title(titles[metric id])
                                                           tmp ax.set ylabel(ylabels[metric id])
```

```
# # end of [for sample_id]
# end of [for metric_id]
# end of [for model_id]

# Create patches for the legend
patches = []
for model_id, model_name in enumerate(model_names):
    patches.append(mpatches.Patch(color=model_colors[model_id], label=model_name))
#
plt.legend(handles=patches, bbox_to_anchor=(1.05, 2.0), loc='upper left', \
    borderaxespad = 0.0, ncol = 1, fontsize = 'x-large')

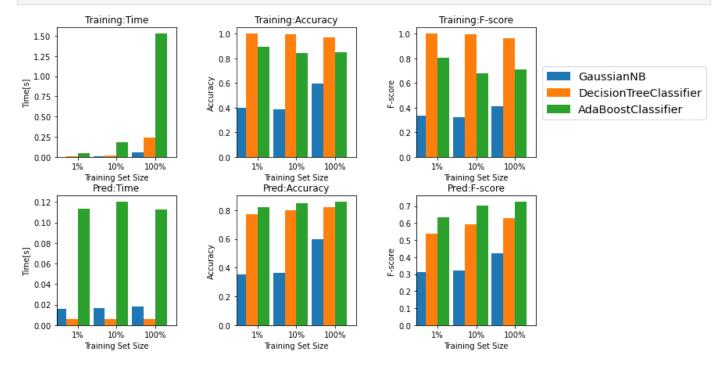
# To avoid overlapping xlabel and title
fig.subplots_adjust(wspace=0.5, hspace=0.3)
```

3-6. Initial Model Evaluation

I will implement the following:

- Import the three supervised learning models which I have discussed in the previous section.
- Initialize the three models
 - I will use the default settings for each model. I will tune one specific model in a later section.
- Calculate the number of records equal to 1%, 10% and 100% of the training data.

```
In [19]: # Import models
        from sklearn.naive bayes import GaussianNB
        from sklearn.ensemble import AdaBoostClassifier
        # Initialize the three models
        models = [GaussianNB(), DecisionTreeClassifier(random state=0), AdaBoostClassifier(random)
        # list of number of sample
        sample list = [int(len(y train)/100), int(len(y train)/10), int(len(y train))]
        # Train models and Predict results
        results = {}
        for model in models:
            model name = model.__class__.__name__
            results[model name] = {}
            for i, sample size in enumerate(sample list):
                t start = time()
                results[model name][i] = train and predict(model, sample size, x train, y train,
                t end = time()
                print("{}[{}]] : {:..3f}[s]".format(model name, i, (t end-t start)))
        print("Finish !")
        GaussianNB[0] : 0.021[s]
        GaussianNB[1] : 0.026[s]
        GaussianNB[2] : 0.072[s]
        DecisionTreeClassifier[0] : 0.012[s]
        DecisionTreeClassifier[1] : 0.027[s]
        DecisionTreeClassifier[2] : 0.246[s]
        AdaBoostClassifier[0] : 0.165[s]
        AdaBoostClassifier[1] : 0.306[s]
        AdaBoostClassifier[2] : 1.642[s]
        Finish!
```



In [22]: show_results_table(results);

Gauss 1 Gauss 10 Gauss 100 train time 0.001966 0.005980 0.051188 pred time 0.015855 0.016969 0.017952 0.400000 0.383333 train acc 0.593333 0.333333 0.325092 0.412500 train f test acc 0.351797 0.366059 0.597678 test f 0.310134 0.320258 0.420899 Decis 1 Decis 10 Decis 100 train_time 0.002992 0.018694 0.237493 pred time 0.005765 0.005985 0.005847 train acc 1.000000 0.996667 0.970000 train f 1.000000 0.997191 0.963855 test acc 0.771918 0.801658 0.818242 test f 0.535978 0.593875 0.627250 AdaBo 1 AdaBo 10 AdaBo 100 1.526668 train time 0.047863 0.182667 pred time 0.113699 0.120073 0.112408 train acc 0.893333 0.840000 0.850000 0.711538 train f 0.801282 0.680147 test acc 0.820674 0.849862 0.857601 test f 0.632757 0.701882 0.724551

3-7. Chose the Best Model

From the table above, I select the Adaboost mode.

- F-score using 100% of traning data(colum= AdaBo_100 , row= test_f) is the best.
- Both of the training and prediction time are small enough.
- From the value of F-score(train_f and test_f), the result of AdaBoost did not overfit.

3-8. Model Tuning

I will fine tune the chosen model using gird search(GridSearchCV). I will implement the following:

- Import sklearn.grid_search.GridSearchCV and sklearn.metrics.make_scorer.
- Initialize the classifier which I've chosen and store it in base model.
- Create a dictionary of parameters I wich to tune for the chosen model.
- Use make_scorer to create an fbeta_score scoring object(with β =0.5)
- Perform grid search on the classifier base_model using the scorer and store it in grid_obj.
- Fit the grid search object to the training data(x_train , y_train), and store it in grid_fit

```
In [23]: import os
         import pickle
         from sklearn.model selection import GridSearchCV
                              import make_scorer, f1_score, fbeta score
        from sklearn.metrics
         # Initialize my model
        base model = AdaBoostClassifier(random state=0)
         file name best model = './Data/Proj 01/best model.sav'
        if os.path.isfile(file name best model):
            print("Load the best model from a file")
             # If there is already a file which contains the best model, load it.
            best model = pickle.load(open(file name best model, 'rb'))
        else:
            print("Train the model with grid search to get the best model")
            # Paremter settings
            params = {}
            params['base estimator'] = [DecisionTreeClassifier(random state=0, max depth=x) for
            params['n estimators'] = [50, 100, 200, 400] # default = 50
            params['learning rate'] = [0.5, 0.6, 0.7, 0.8, 0.9, 1.0, 1.1] # default = 1.0
            # Grid search and Cross Validation
            scorer = make scorer(fbeta score, beta=0.5)
            grid obj = GridSearchCV(base model, params, scoring=scorer)
            grid fit = grid obj.fit(x train, y train)
            best model = grid fit.best estimator
             # save the model to a file
            pickle.dump(best model, open(file name best model, 'wb'))
```

Load the best model from a file

I show the score

```
In [24]: pred base = (base model.fit(x train, y train)).predict(x test)
        pred best = best model.predict(x test)
        # Report the before-and-afterscores
        print("----- Unoptimized model ----")
        print("Accuracy score on testing data: {:.4f}".format(accuracy score(y test, pred base))
        print("F-score on testing data: {:.4f}".format(fbeta score(y test, pred base, bet
        print("----- Optimized model ----")
        print("Final accuracy score on the testing data: {:.4f}".format(accuracy score(y test, p
        print("Final F-score on the testing data:
                                                  {:.4f}".format(fbeta score(y test, pred
        ----- Unoptimized model -----
        Accuracy score on testing data: 0.8576
        F-score on testing data:
                                      0.7246
        ----- Optimized model -----
        Final accuracy score on the testing data: 0.8690
        Final F-score on the testing data: 0.7489
```

• These are parameter of the best_model .

```
In [25]: best_model.get params()
         { 'algorithm': 'SAMME.R',
Out[25]:
         'base estimator ccp alpha': 0.0,
         'base estimator class weight': None,
          'base estimator criterion': 'gini',
         'base estimator max depth': 2,
         'base estimator max features': None,
          'base estimator max leaf nodes': None,
          'base estimator min impurity decrease': 0.0,
         'base_estimator__min_samples leaf': 1,
          'base estimator min samples split': 2,
          'base estimator min weight fraction leaf': 0.0,
          'base_estimator__random_state': 0,
         'base estimator splitter': 'best',
          'base estimator': DecisionTreeClassifier(max depth=2, random state=0),
          'learning rate': 0.5,
          'n estimators': 100,
          'random state': 0}
```

4. Evaluation

4-1. Final Model Evaluation

I show the results.

Metric	Naive Predictor	Unoptimized AdaBoost	Optimized AdaBoost
Accuracy Score	0.2478	0.8576	0.8690
F-score	0.2917	0.7246	0.7489

Compaired to the Navive Predictor, the performance of the Optimized Model is improved by 62.1%. Compaired to the Unoptimized Model, the performance of the Optimized Model is improved by 2.4%.

4-2. Feature Importance

An important task when performing supervised learning on a dataset like the census data is determining which features provide the most predictive power.

By focusing on the relationship between only a few crucial features and the target label, I can simplify our understanding of the phenomenon, which is most always a useful thing to do.

In the case of this project, that means I wish to identify a small number of features that most strongly predict whether an indivisual makes at most or more than \$50,000.

Some scikit-learn classifiers(e.g., adaboost, random forests) have a feature_importance_ attributee, which is a function that ranks the importance of features. In the code cell below, I will use this attribute to determin the top X most important features for the census dataset.

This is a plot function which display the top X most important features

In [26]:

```
def plot features(importances, x train, y train, top x, figsize=(15,5)):
             # Get Top X most important features
            indices = np.argsort(importances)[::-1]
            columns = x train.columns.values[indices[:top x]]
            weights = importances[indices][:top x]
            # Create Figure
            bar x1 = np.arange(top x) - 0.3
            bar x2 = np.arange(top x)
            bar y1 = weights
            bar y2 = np.cumsum(weights)
            colors = ["tab:blue", "tab:orange", "tab:green", "tab:red", "tab:purple"]
            labels = ["Feature Weight", "Cumulative Feature Weight"]
            fig = plt.figure(figsize=figsize)
            plt.title("Normalized Wieghts for Frist Five Most Predictive Features", fontsize=16)
             # Plot Bar
            plt.bar(bar x1, bar y1, width=0.6, align="center", color=colors[0], label=labels[0])
            plt.bar(bar x2, bar y2, width=0.3, align="center", color=colors[1], label=labels[1])
            plt.xticks(bar x1, columns, fontsize=12)
            plt.xlim((-0.5, top x-0.5))
            plt.xlabel("Feature", fontsize=12)
            plt.xlabel("Weight", fontsize=12)
            plt.legend(loc="upper center")
            plt.tight layout()
            return columns
         # Extranct feature importances
In [27]:
         importances = best model.feature importances
         # plot
         top x = 6
```

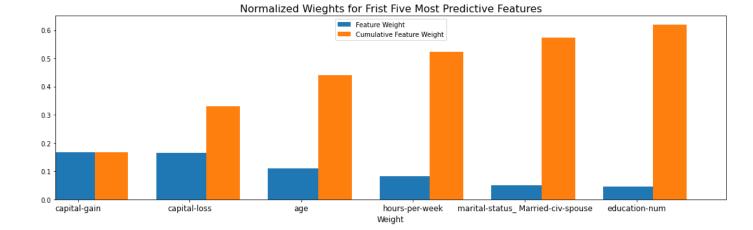
important columns = plot features(importances, x train, y train, top x)

'marital-status Married-civ-spouse', 'education-num'],

array(['capital-gain', 'capital-loss', 'age', 'hours-per-week',

display(important columns)

dtype=object)



Observation about the top 5 features extraction

- 1. capital-gain
- 2. capital-loss
 - I guess that those who can afford to invest can donate a lot.
- 3. age
 - I guess that experience of how to earn relates with investing.
- 4. hours-per-week
 - I guess that who can use more time to invest makes more money.
- 5. marital-status_ Married-civ-spouse
 - I geuss that a lot of citizens with spouse can afford to pay attention not only to themselves but also to their surrounding.

Features Selection

I will train the model with less features. The expectation is that training and prediction time is much lower. From the visualization above, I see that the top 6 most important features contributes more than half of the importance of all features present in the data.

This hins that I can attempt to reduce the feature space and simplify the information required to the model to learn.

```
from sklearn.base import clone
In [28]:
         # create reduced features ant target
         x train reduced = x train[important columns]
         x test reduced = x test[important columns]
         # calculate the training time of best model
         t start = time()
         best model = best model.fit(x train, y train)
         training time best model = time() - t start
         # clone the best model and train
         light model = clone(best model)
         t start = time()
         light model = light model.fit(x train reduced, y train)
         training time light model = time() - t start
         # predict
         pred reduced = light model.predict(x test reduced)
         # Report the before-and-afterscores
```

```
print("----- Optimized model -----")
print("Accuracy score on testing data: {:.4f}".format(accuracy score(y test, pred best))
print("F-score on testing data: {:.4f}".format(fbeta_score(y_test, pred_best, bet
print("Training Time[s]: {:.4f}".format(training_time_best_model))
print("----- Optimized model on reduced data -----")
print("Accuracy score on the testing data: {:.4f}".format(accuracy score(y test, pred re
print("F-score on the testing data: {:.4f}".format(fbeta_score(y_test, pred_reduc
print("Training Time[s]:
                                           {:.4f}".format(training time light model))
----- Optimized model -----
Accuracy score on testing data: 0.8690
F-score on testing data: 0.7489
                               4.3591
Training Time[s]:
----- Optimized model on reduced data -----
Accuracy score on the testing data: 0.8601
F-score on the testing data: 0.7318
Training Time[s]:
                                     0.9891
```

Optimized model used 103 features, and optimized model on reduced data used only 6 features. With reduced features, the F-score reduced only 1.7% and training time is about 4 time faster.

For more information about Features Selection

I will research the detail effect of the number of features which are used on training. I will show the training time, accuracy and f-score chaging the number of features from 1 to 10.

```
In [29]:
        indices = np.argsort(importances)[::-1]
        all columns sorted = x train.columns.values[indices]
        train time list = []
        accuracy list = []
        fscore list = []
        for cnt in range(10):
            top x = cnt + 1
           important columns = all_columns_sorted[:top_x]
            #----#
            # create reduced features ant target
            x train reduced = x train[important columns]
            x_test_reduced = x_test[important columns]
            # clone the best model and train
            light model = clone(best model)
            t start = time()
            light model = light model.fit(x train reduced, y train)
            t end = time()
            pred reduced = light model.predict(x test reduced)
            # output
            train time list.append(t end-t start)
            accuracy list.append(accuracy score(y test, pred reduced))
            fscore list.append(fbeta score(y test, pred reduced, beta = 0.5))
```

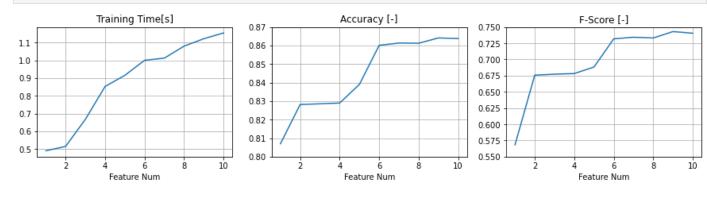
```
In [30]: fig = plt.figure(figsize=(15,3))
x_data = np.arange(10) + 1

ax = fig.add_subplot(1,3,1)
plt.plot(x_data, train_time_list)
plt.grid(1)
plt.title('Training Time[s]')
```

```
plt.xlabel('Feature Num')

ax = fig.add_subplot(1,3,2)
plt.plot(x_data, accuracy_list)
plt.grid(1)
plt.title('Accuracy [-]')
plt.xlabel('Feature Num')
plt.ylim((0.8, 0.87));

ax = fig.add_subplot(1,3,3)
plt.plot(x_data, fscore_list)
plt.grid(1)
plt.title('F-Score [-]')
plt.xlabel('Feature Num')
plt.ylim((0.55, 0.75));
```



In []: