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# CSE4020 MACHINE LEARNING ETH THEORY

## TOPIC: SEMI – SUPERVISED LEARNING (WEB CONTENT)

### **Abstract**

Semi-supervised learning is a machine learning technique that combines both labeled and unlabeled data to improve the accuracy of models. In the context of web content, semisupervised learning can be used to classify and categorize large amounts of data, such as text and images, that may not have explicit labels. By leveraging the relationships and patterns between labeled and unlabeled data, semi-supervised learning can effectively identify and classify web content with higher accuracy than traditional supervised learning methods. This approach is particularly useful for applications such as web search, recommendation systems, and content filtering, where there is a vast amount of unstructured data that needs to be organized and analyzed. However, as with any machine learning technique, semi-supervised learning has its limitations and requires careful consideration of the data and model architecture to achieve optimal results. This can be particularly useful for applications such as sentiment analysis, topic modeling, and content recommendation systems. Overall, semi-supervised learning can help to improve the efficiency and accuracy of web content analysis, making it a valuable tool for businesses and researchers alike.

### Introduction

Web content analysis has become an essential part of businesses and researchers' strategies to understand their target audience, identify trends, and generate insights. However, analyzing vast amounts of text data can be a daunting task, and manually labeling every data point can be extremely time-consuming and resource intensive. Semi-supervised learning offers an alternative approach to analyzing web content, which can significantly reduce the need for manual labeling and improve the accuracy of analysis. In semi-supervised learning, models are trained using a combination of labeled and unlabeled data, allowing them to learn patterns and relationships in the data and classify new, unlabeled data accurately. This approach can be particularly useful in analyzing web content, where a vast amount of data is available, but only a small portion of it may be labeled. In this paper, we will explore the functionalities, benefits, and challenges of using semi-supervised learning for web content analysis.

### <u>Functionalities of Semi – Supervised Learning</u>

Semi-supervised learning is a machine learning technique that falls in between supervised and unsupervised learning. In this case, some of the training data is labeled, while the rest is unlabeled. This allows the model to learn from both labeled and unlabeled data, which can be beneficial when it's labeled data is scarce or expensive to obtain. One potential application of semi-supervised learning is in the field of web content analysis. With the vast amount of information available on the internet, it can be challenging to classify and organize web content. However, with the help of semi-supervised learning, it is possible to classify web content with high accuracy.

In this project we use a combination of supervised and unsupervised learning techniques to classify web pages. Initially, a small set of web pages are physically labeled to provide a starting point for the model. This labeled data is then used to train a supervised learning model. The model is then used to predict the category of other web pages. However, since labeled data is often limited in web content analysis, the model is also trained using unsupervised learning techniques. This involves clustering similar web pages together based on their content. The model then uses these clusters to infer the category of other web pages. The project's functionality is based on a web crawler that scans the internet for web pages and extracts their content. The content is then fed into the model, which predicts the category of the web page. The categories are then displayed on a user interface, which allows users to search for web pages based on their category.

In summary, the Web Content Classification Project is a semi-supervised learning project that aims to classify web pages based on their content. The project uses a combination of supervised and unsupervised learning techniques to classify web pages accurately. The project's functionality is based on a web crawler that extracts web page content and feeds it into the model. The results are then displayed on a user interface, which allows users to search for web pages based on their category.

Semi-supervised learning can be a valuable tool in analyzing and categorizing web content. Some of the key functionalities of semi-supervised learning in the context of web content include:

- 1. <u>Data labeling:</u> In semi-supervised learning, labeled data is used to train a model, which can then be used to categorize larger amounts of unlabeled data. With web content, this can involve manually describing a small set of data, which can then be used to train a model that can categorize much larger amounts of unlabeled data.
- 2. <u>Text classification</u>: One of the main applications of semi-supervised learning in web content analysis is text classification. This involves categorizing text data into different topics or themes. For example, a model might be trained to categorize news articles into different categories such as politics, sports, and entertainment.
- 3. **Sentiment analysis:** Another application of semi-supervised learning in web content analysis is sentiment analysis. This involves categorizing text data based on the sentiment expressed, such as positive, negative, or neutral. This can be particularly useful for analyzing social media data, online reviews, and customer feedback.
- 4. <u>Topic modeling:</u> Semi-supervised learning can also be used for topic modeling, which involves identifying the underlying themes or topics in a set of text data. This can be useful for understanding the content of large amounts of web content, such as news articles or blog posts.
- 5. <u>Content recommendation:</u> Finally, semi-supervised learning can be used for content recommendation systems, which suggest relevant content to users based on their browsing history or other behavior. This can involve training a model to recommend similar content to a user based on the content they have previously viewed or interacted with.

Overall, semi-supervised learning can be a powerful tool in web content analysis, enabling businesses and researchers to more efficiently and accurately categorize and analyze large amounts of text data.

### **Observations**

The Web Content Classification Project is an interesting and useful application of semi-supervised learning in the field of web content analysis. It demonstrates the potential of semi-supervised learning to improve the accuracy of web content classification by combining both labeled and unlabeled data.

One possible additional functionality for the project could be to allow users to provide feedback on the accuracy of the classification. This feedback can be used to improve the model and make it more accurate over time. For example, users could be given the option to suggest a different category for a misclassified web page, or to flag a web page that is incorrectly classified.

Another possible functionality could be to incorporate user preferences into the classification process. For instance, users could specify their preferred categories for certain topics, and the model could prioritize these preferences when classifying web pages. This would make the classification process more personalized and tailored to individual user needs.

Overall, there are many possibilities for additional functionalities that could be added to the Web Content Classification Project to improve its accuracy and usefulness. By incorporating user feedback and preferences, the project could become even more effective at classifying web content and providing relevant search results.

### Code

```
import pandas as pd
import numpy as np
import seaborn as sns
import matplotlib.pyplot as plt
%matplotlib inline
sns.set()
df = pd.read_csv('oasis_longitudinal.csv')
df.head()
df = df.loc[df['Visit']==1]
df = df.reset index(drop=True)
df['M/F'] = df['M/F'].replace(['F','M'], [0,1])
df['Group'] = df['Group'].replace(['Converted'], ['Demented'])
df['Group'] = df['Group'].replace(['Demented', 'Nondemented'], [1,0])
df = df.drop(['MRI ID', 'Visit', 'Hand'], axis=1)
def bar chart(feature):
  Demented = df[df['Group']==1][feature].value counts()
  Nondemented = df[df['Group']==0][feature].value_counts()
  df bar = pd.DataFrame([Demented,Nondemented])
  df bar.index = ['Demented','Nondemented']
  df bar.plot(kind='bar',stacked=True, figsize=(8,5))
```

```
bar_chart('M/F')
plt.xlabel('Group')
plt.ylabel('Number of patients')
plt.legend()
plt.title('Gender and Demented rate')
facet= sns.FacetGrid(df,hue="Group", aspect=3)
facet.map(sns.kdeplot,'MMSE',shade= True)
facet.set(xlim=(0, df['MMSE'].max()))
facet.add_legend()
plt.xlim(15.30)
facet= sns.FacetGrid(df,hue="Group", aspect=3)
facet.map(sns.kdeplot,'ASF',shade= True)
facet.set(xlim=(0, df['ASF'].max()))
facet.add_legend()
plt.xlim(0.5, 2)
facet= sns.FacetGrid(df,hue="Group", aspect=3)
facet.map(sns.kdeplot,'eTIV',shade= True)
facet.set(xlim=(0, df['eTIV'].max()))
facet.add legend()
plt.xlim(900, 2100)
facet= sns.FacetGrid(df,hue="Group", aspect=3)
facet.map(sns.kdeplot,'nWBV',shade= True)
facet.set(xlim=(0, df['nWBV'].max()))
```

```
facet.add_legend()
plt.xlim(0.6,0.9)
facet= sns.FacetGrid(df,hue="Group", aspect=3)
facet.map(sns.kdeplot,'Age',shade= True)
facet.set(xlim=(0, df['Age'].max()))
facet.add_legend()
plt.xlim(50,100)
facet= sns.FacetGrid(df,hue="Group", aspect=3)
facet.map(sns.kdeplot,'EDUC',shade= True)
facet.set(xlim=(df['EDUC'].min(), df['EDUC'].max()))
facet.add_legend()
plt.ylim(0, 0.16)
pd.isnull(df).sum()
df dropna = df.dropna(axis=0, how='any')
pd.isnull(df dropna).sum()
df_dropna['Group'].value_counts()
x = df['EDUC']
y = df['SES']
ses_not_null_index = y[~y.isnull()].index
x = x[ses not null index]
```

```
y = y[ses_not_null_index]
z = np.polyfit(x, y, 1)
p = np.poly1d(z)
plt.plot(x, y, 'go', x, p(x), "r--")
plt.xlabel('Education Level(EDUC)')
plt.ylabel('Social Economic Status(SES)')
plt.show()
df.groupby(['EDUC'])['SES'].median()
df["SES"].fillna(df.groupby("EDUC")["SES"].transform("median"), inplace=True)
pd.isnull(df['SES']).value counts()
import pandas as pd
import numpy as np
from sklearn.model selection import train test split
from sklearn import preprocessing
from sklearn.preprocessing import MinMaxScaler
from sklearn.model selection import cross val score
Y = df['Group'].values
X = df[['M/F', 'Age', 'EDUC', 'SES', 'MMSE', 'eTIV', 'nWBV', 'ASF']]
X trainval, X test, Y trainval, Y test = train test split(X, Y, random state=0)
```

```
scaler = MinMaxScaler().fit(X_trainval)
X_trainval_scaled = scaler.transform(X_trainval)
X test scaled = scaler.transform(X test)
Y = df_dropna['Group'].values
X = df dropna[['M/F', 'Age', 'EDUC', 'SES', 'MMSE', 'eTIV', 'nWBV', 'ASF']]
X trainval dna, X test dna, Y trainval dna, Y test dna = train test split(X, Y, random state=0)
scaler = MinMaxScaler().fit(X trainval dna)
X_trainval_scaled_dna = scaler.transform(X_trainval_dna)
X test scaled dna = scaler.transform(X test dna)
from sklearn.linear model import LogisticRegression
from sklearn.svm import SVC
from sklearn.tree import DecisionTreeClassifier
from sklearn.ensemble import RandomForestClassifier
from sklearn.ensemble import AdaBoostClassifier
from sklearn.metrics import confusion matrix, accuracy score, recall score, roc curve, auc
acc = []
best score=0
kfolds=5
for c in [0.001, 0.1, 1, 10, 100]:
  logRegModel = LogisticRegression(C=c)
```

```
scores = cross_val_score(logRegModel, X_trainval, Y_trainval, cv=kfolds, scoring='accuracy')
  score = np.mean(scores)
  if score > best score:
    best score = score
    best_parameters = c
SelectedLogRegModel = LogisticRegression(C=best_parameters).fit(X_trainval_scaled, Y_trainval)
test score = SelectedLogRegModel.score(X test scaled, Y test)
PredictedOutput = SelectedLogRegModel.predict(X test scaled)
test recall = recall score(Y test, PredictedOutput, pos label=1)
fpr, tpr, thresholds = roc curve(Y test, PredictedOutput, pos label=1)
test auc = auc(fpr, tpr)
print("Best accuracy on validation set is:", best score)
print("Best parameter for regularization (C) is: ", best parameters)
print("Test accuracy with best C parameter is", test score)
print("Test recall with the best C parameter is", test recall)
print("Test AUC with the best C parameter is", test auc)
m = 'Logistic Regression (w/ imputation)'
acc.append([m, test_score, test_recall, test_auc, fpr, tpr, thresholds])
best score=0
kfolds=5
for c in [0.001, 0.1, 1, 10, 100]:
  logRegModel = LogisticRegression(C=c)
```

```
scores = cross val score(logRegModel, X trainval scaled dna, Y trainval dna, cv=kfolds,
scoring='accuracy')
  score = np.mean(scores)
  if score > best_score:
    best score = score
    best parameters = c
SelectedLogRegModel = LogisticRegression(C=best_parameters).fit(X trainval scaled dna,
Y_trainval_dna)
test score = SelectedLogRegModel.score(X test scaled dna, Y test dna)
PredictedOutput = SelectedLogRegModel.predict(X test scaled)
test recall = recall score(Y test, PredictedOutput, pos label=1)
fpr, tpr, thresholds = roc curve(Y test, PredictedOutput, pos label=1)
test auc = auc(fpr, tpr)
print("Best accuracy on validation set is:", best score)
print("Best parameter for regularization (C) is: ", best parameters)
print("Test accuracy with best C parameter is", test_score)
print("Test recall with the best C parameter is", test recall)
print("Test AUC with the best C parameter is", test auc)
m = 'Logistic Regression (w/ dropna)'
acc.append([m, test score, test recall, test recall, fpr, tpr, thresholds])
best score = 0
```

```
for c paramter in [0.001, 0.01, 0.1, 1, 10, 100, 1000]:
  for gamma paramter in [0.001, 0.01, 0.1, 1, 10, 100, 1000]:
    for k parameter in ['rbf', 'linear', 'poly', 'sigmoid']:
      svmModel = SVC(kernel=k parameter, C=c paramter, gamma=gamma paramter)
      scores = cross_val_score(svmModel, X_trainval_scaled, Y_trainval, cv=kfolds,
scoring='accuracy')
      score = np.mean(scores)
      if score > best score:
        best score = score
        best parameter c = c paramter
        best_parameter_gamma = gamma_paramter
        best parameter k = k parameter
SelectedSVMmodel = SVC(C=best parameter c, gamma=best parameter gamma,
kernel=best parameter k).fit(X trainval scaled, Y trainval)
test score = SelectedSVMmodel.score(X test scaled, Y test)
PredictedOutput = SelectedSVMmodel.predict(X test scaled)
test_recall = recall_score(Y_test, PredictedOutput, pos_label=1)
fpr, tpr, thresholds = roc curve(Y test, PredictedOutput, pos label=1)
test auc = auc(fpr, tpr)
print("Best accuracy on cross validation set is:", best score)
print("Best parameter for c is: ", best parameter c)
print("Best parameter for gamma is: ", best parameter gamma)
```

```
print("Best parameter for kernel is: ", best_parameter_k)
print("Test accuracy with the best parameters is", test_score)
print("Test recall with the best parameters is", test recall)
print("Test recall with the best parameter is", test auc)
m = 'SVM'
acc.append([m, test score, test recall, test auc, fpr, tpr, thresholds])
from sklearn.tree import DecisionTreeClassifier
best score = 0
for md in range(1, 9):
  treeModel = DecisionTreeClassifier(random state=0, max depth=md, criterion='gini')
  scores = cross_val_score(treeModel, X_trainval_scaled, Y_trainval, cv=kfolds,
scoring='accuracy')
  score = np.mean(scores)
  if score > best_score:
    best score = score
    best parameter = md
SelectedDTModel = DecisionTreeClassifier(max_depth=best_parameter).fit(X_trainval_scaled,
Y trainval)
```

```
test score = SelectedDTModel.score(X test scaled, Y test)
PredictedOutput = SelectedDTModel.predict(X test scaled)
test recall = recall score(Y test, PredictedOutput, pos label=1)
fpr, tpr, thresholds = roc curve(Y test, PredictedOutput, pos label=1)
test_auc = auc(fpr, tpr)
print("Best accuracy on validation set is:", best score)
print("Best parameter for the maximum depth is: ", best parameter)
print("Test accuracy with best parameter is ", test score)
print("Test recall with best parameters is ", test recall)
print("Test AUC with the best parameter is ", test auc)
m = 'Decision Tree'
acc.append([m, test score, test recall, test auc, fpr, tpr, thresholds])
print("Feature importance: ")
np.array([X.columns.values.tolist(), list(SelectedDTModel.feature importances )]).T
from sklearn.tree import export graphviz
import graphviz
dot data=export graphviz(SelectedDTModel,
feature names=X trainval.columns.values.tolist(),out file=None)
graph = graphviz.Source(dot_data)
graph
best score = 0
for M in range(2, 15, 2):
```

```
for d in range(1, 9):
    for m in range(1, 9):
      forestModel = RandomForestClassifier(n estimators=M, max features=d, n jobs=4,
                       max_depth=m, random_state=0)
      scores = cross val score(forestModel, X trainval scaled, Y trainval, cv=kfolds,
scoring='accuracy')
      score = np.mean(scores)
      if score > best score:
        best_score = score
        best M = M
        best d = d
        best m = m
SelectedRFModel = RandomForestClassifier(n_estimators=M, max_features=d,
                       max depth=m, random state=0).fit(X trainval scaled, Y trainval)
PredictedOutput = SelectedRFModel.predict(X test scaled)
test_score = SelectedRFModel.score(X_test_scaled, Y_test)
test recall = recall score(Y test, PredictedOutput, pos label=1)
fpr, tpr, thresholds = roc curve(Y test, PredictedOutput, pos label=1)
test_auc = auc(fpr, tpr)
print("Best accuracy on validation set is:", best score)
print("Best parameters of M, d, m are: ", best M, best d, best m)
```

```
print("Test accuracy with the best parameters is", test_score)
print("Test recall with the best parameters is:", test_recall)
print("Test AUC with the best parameters is:", test auc)
m = 'Random Forest'
acc.append([m, test score, test recall, test auc, fpr, tpr, thresholds])
print("Feature importance: ")
np.array([X.columns.values.tolist(), list(SelectedRFModel.feature importances )]).T
best score = 0
for M in range(2, 15, 2):
  for Ir in [0.0001, 0.001, 0.01, 0.1, 1]:
    # train the model
    boostModel = AdaBoostClassifier(n_estimators=M, learning_rate=Ir, random_state=0)
    scores = cross val score(boostModel, X trainval scaled, Y trainval, cv=kfolds,
scoring='accuracy')
    score = np.mean(scores)
    if score > best score:
      best score = score
      best_M = M
      best Ir = Ir
```

```
SelectedBoostModel = AdaBoostClassifier(n estimators=M, learning rate=Ir,
random_state=0).fit(X_trainval_scaled, Y_trainval)
PredictedOutput = SelectedBoostModel.predict(X test scaled)
test score = SelectedRFModel.score(X test scaled, Y test)
test_recall = recall_score(Y_test, PredictedOutput, pos_label=1)
fpr, tpr, thresholds = roc curve(Y test, PredictedOutput, pos label=1)
test auc = auc(fpr, tpr)
print("Best accuracy on validation set is:", best_score)
print("Best parameter of M is: ", best M)
print("best parameter of LR is: ", best Ir)
print("Test accuracy with the best parameter is", test score)
print("Test recall with the best parameters is:", test_recall)
print("Test AUC with the best parameters is:", test auc)
m = 'AdaBoost'
acc.append([m, test score, test recall, test auc, fpr, tpr, thresholds])
print("Feature importance: ")
np.array([X.columns.values.tolist(), list(SelectedBoostModel.feature importances )]).T
result = pd.DataFrame(acc, columns=['Model', 'Accuracy', 'Recall', 'AUC', 'FPR', 'TPR', 'TH'])
result[['Model', 'Accuracy', 'Recall', 'AUC']]
```

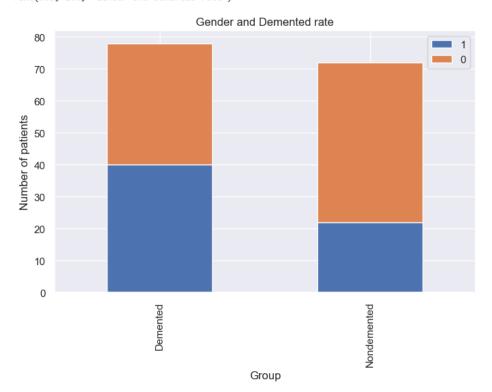
### **Output**

```
In [1]: import pandas as pd
              import numpy as np
import seaborn as sns
import matplotlib.pyplot as plt
              %matplotlib inline
              sns.set()
              df = pd.read_csv('oasis_longitudinal.csv')
              df.head()
Out[1]:
                                                                   Group Visit MR Delay M/F Hand Age EDUC SES MMSE CDR eTIV nWBV ASF
                     Subject ID
                                                MRI ID
               0 OAS2_0001 OAS2_0001_MR1 Nondemented 1 0 M R 87 14 2.0 27.0 0.0 1987 0.696 0.883
               1 OAS2_0001 OAS2_0001_MR2 Nondemented 2
                                                                                                             R 88
               2 OAS2_0002 OAS2_0002_MR1 Demented 1 0 M R 75 12 NaN 23.0 0.5 1678 0.736 1.046
              3 OAS2_0002 OAS2_0002_MR2 Demented 2 560 M R 76 12 NaN 28.0 0.5 1738 0.713 1.010
4 OAS2_0002 OAS2_0002_MR3 Demented 3 1895 M R 80 12 NaN 22.0 0.5 1698 0.701 1.034
In [2]: df = df.loc[df['Visit']=-1]
    df = df.reset_index(drop=True)
    df['M/r'] = df['M/r'].replace(['f','M'], [0,1])
    df['Group'] = df['Group'].replace(['Converted'], ['Demented'])
    df['Group'] = df['Group'].replace(['Demented', 'Nondemented'], [1,0])
    df = df.drop(['MRI ID', 'Visit', 'Hand'], axis=1)
In [3]:

def bar_chart(feature):
    Demented = df[df['Group']==1][feature].value_counts()
    Nondemented = df[df['Group']==0][feature].value_counts()
    df_bar = pd.DataFrame([Demented,Nondemented])
    df_bar.index = ['Demented','Nondemented']
    df_bar.plot(kind='bar',stacked=True, figsize=(8,5))
```

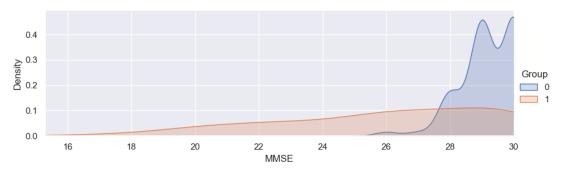
```
In [4]: bar_chart('M/F')
  plt.xlabel('Group')
  plt.ylabel('Number of patients')
  plt.legend()
  plt.title('Gender and Demented rate')
```

Out[4]: Text(0.5, 1.0, 'Gender and Demented rate')



```
In [5]: facet= sns.FacetGrid(df,hue="Group", aspect=3)
facet.map(sns.kdeplot,'MMSE',shade= True)
facet.set(xlim=(0, df['MMSE'].max()))
facet.add_legend()
plt.xlim(15.30)
```

Out[5]: (15.3, 30.0)

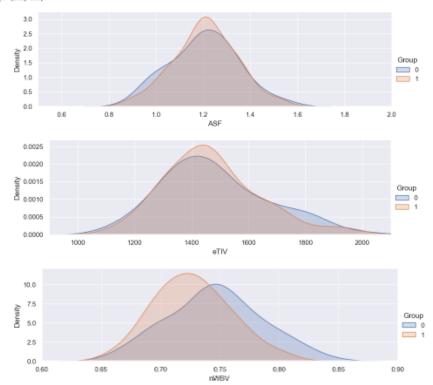


```
In [6]: facet= sns.FacetGrid(df,hue="Group", aspect=3)
facet.map(sns.kdeplot,'ASF',shade= True)
facet.set(xlim=(0, df['ASF'].max()))
facet.add_legend()
plt.xlim(0.5, 2)

facet= sns.FacetGrid(df,hue="Group", aspect=3)
facet.map(sns.kdeplot,'eTIV',shade= True)
facet.set(xlim=(0, df['eTIV'].max()))
facet.add_legend()
plt.xlim(900, 2100)

facet= sns.FacetGrid(df,hue="Group", aspect=3)
facet.map(sns.kdeplot,'nNBV',shade= True)
facet.set(xlim=(0, df['nNBV'].max()))
facet.add_legend()
plt.xlim(0.6,0.9)
```

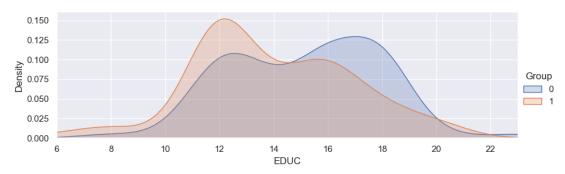
Out[6]: (0.6, 0.9)



```
In [7]: facet= sns.FacetGrid(df,hue="Group", aspect=3)
facet.map(sns.kdeplot, 'Age',shade= True)
facet.set(xlim=(0, df['Age'].max()))
facet.add_legend()
plt.xlim(50,100)
Out[7]: (50.0, 100.0)
                     0.05
                     0.04
                0.03
0.02
                                                                                                                                                                                                                   Group
                                                                                                                                                                                                                  0
                                                                                                                                                                                                                  ____ 1
                     0.01
                     0.00
                             50
                                                                 60
                                                                                                    70
                                                                                                                                        80
                                                                                                                                                                           90
                                                                                                                                                                                                              100
                                                                                                                    Age
```

```
In [8]:
facet= sns.FacetGrid(df,hue="Group", aspect=3)
facet.map(sns.kdeplot,'EDUC',shade= True)
facet.set(xlim=(df['EDUC'].min(), df['EDUC'].max()))
facet.add_legend()
plt.ylim(0, 0.16)
```

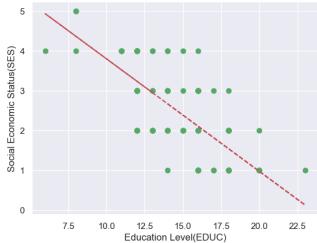




### In [9]: pd.isnull(df).sum()

```
Out[9]: Subject ID 0
Group 0
MR Delay 0
MF 0
Age 0
EDUC 0
SES 8
MMSE 0
CDR 0
eTIV 0
nWBV 0
ASF 0
dtype: int64
```

```
In [10]: df_dropna = df.dropna(axis=0, how='any')
pd.isnull(df_dropna).sum()
   Out[10]: Subject ID
                  Group
                  MR Delay
M/F
                                         0
0
0
                  Age
EDUC
                  SES
MMSE
                                         0
                                         0
                  CDR
                  eTIV
                  nWBV
ASF
                                         0
                                         0
                  dtype: int64
   In [11]: df_dropna['Group'].value_counts()
  Out[11]: 0 72
1 70
                  Name: Group, dtype: int64
In [12]: x = df['EDUC']
y = df['SES']
              ses_not_null_index = y[~y.isnull()].index
x = x[ses_not_null_index]
y = y[ses_not_null_index]
              z = np.polyfit(x, y, 1)
p = np.poly1d(z)
plt.plot(x, y, 'go', x, p(x), "r--")
plt.xlabel('Education Level(EDUC)')
plt.ylabel('Social Economic Status(SES)')
               plt.show()
                      5
```



```
In [13]: df.groupby(['EDUC'])['SES'].median()
Out[13]: EDUC
            8
                    5.0
            11
                  4.0
            12
                   3.0
            13
                   2.0
             15
                   2.0
             16
                    2.0
             17
                    1.0
                    2.0
             20
                   1.0
             23
                   1.0
             Name: SES, dtype: float64
In [14]: df["SES"].fillna(df.groupby("EDUC")["SES"].transform("median"), inplace=True)
In [15]: pd.isnull(df['SES']).value_counts()
Out[15]: False 150
Name: SES, dtype: int64
In [16]: import pandas as pd
             import numpy as np
             from sklearn.model_selection import train_test_split
            from sklearn import preprocessing
from sklearn preprocessing import MinMaxScaler
            from sklearn.model_selection import cross_val_score
In [17]:
Y = df['Group'].values
X = df[['M/F', 'Age', 'EDUC', 'SES', 'MMSE', 'eTIV', 'nWBV', 'ASF']]
X_trainval, X_test, Y_trainval, Y_test = train_test_split(X, Y, random_state=0)
             scaler = MinMaxScaler().fit(X_trainval)
            X_trainval_scaled = scaler.transform(X_trainval)
X_test_scaled = scaler.transform(X_test)
    In [18]: Y = df_dropna['Group'].values
X = df_dropna[['M/F', 'Age', 'EDUC', 'SES', 'MMSE', 'eTIV', 'nWBV', 'ASF']]
X_trainval_dna, X_test_dna, Y_trainval_dna, Y_test_dna = train_test_split(X, Y, random_state=0)
                scaler = MinMaxScaler().fit(X_trainval_dna)
                X_trainval_scaled_dna = scaler.transform(X_trainval_dna)
X_test_scaled_dna = scaler.transform(X_test_dna)
    In [19]: from sklearn.linear model import LogisticRegression
                 from sklearn.svm import SVC
                from sklearn.tree import DecisionTreeClassifier
from sklearn.ensemble import RandomForestClassifier
                 from sklearn.ensemble import AdaBoostClassifier
                from sklearn.metrics import confusion_matrix, accuracy_score, recall_score, roc_curve, auc
    In [20]: acc = []
```

```
In [21]: best score=0
          kfolds=5
          for c in [0.001, 0.1, 1, 10, 100]:
              logRegModel = LogisticRegression(C=c)
               scores = cross_val_score(logRegModel, X_trainval, Y_trainval, cv=kfolds, scoring='accuracy')
               score = np.mean(scores)
               if score > best_score:
                   best score = score
                   best parameters = c
          SelectedLogRegModel = LogisticRegression(C=best\_parameters).fit(X\_trainval\_scaled, Y\_trainval)
          test score = SelectedLogRegModel.score(X test scaled, Y test)
          PredictedOutput = SelectedLogRegModel.predict(X_test_scaled)
          test_recall = recall_score(Y_test, PredictedOutput, pos_label=1)
          fpr, tpr, thresholds = roc_curve(Y_test, PredictedOutput, pos_label=1)
          test_auc = auc(fpr, tpr)
          print("Best accuracy on validation set is:", best_score)
print("Best parameter for regularization (C) is: ", best_parameters)
          print("Test accuracy with best C parameter is", test_score)
          print("Test recall with the best C parameter is", test_recall)
          print("Test AUC with the best C parameter is", test_auc)
          m = 'Logistic Regression (w/ imputation)
          acc.append([m, test_score, test_recall, test_auc, fpr, tpr, thresholds])
         Increase the number of iterations (max_iter) or scale the data as shown in:
               https://scikit-learn.org/stable/modules/preprocessing.html
          Please also refer to the documentation for alternative solver options: 
https://scikit-learn.org/stable/modules/linear_model.html#logistic-regression
            n_iter_i = _check_optimize_result(
          Best accuracy on validation set is: 0.724901185770751
          Best parameter for regularization (C) is: 10
Test accuracy with best C parameter is 0.7631578947368421
Test recall with the best C parameter is 0.7
          Test AUC with the best C parameter is 0.7666666666666667
In [22]: best score=0
           for c in [0.001, 0.1, 1, 10, 100]:
               logRegModel = LogisticRegression(C=c)
               scores = cross_val_score(logRegModel, X_trainval_scaled_dna, Y_trainval_dna, cv=kfolds, scoring='accuracy')
               score = np.mean(scores)
               if score > best_score:
                   best score = score
                   best_parameters = c
           SelectedLogRegModel = LogisticRegression(C=best_parameters).fit(X_trainval_scaled_dna, Y_trainval_dna)
           test_score = SelectedLogRegModel.score(X_test_scaled_dna, Y_test_dna)
           PredictedOutput = SelectedLogRegModel.predict(X_test_scaled)
           test_recall = recall_score(Y_test, PredictedOutput, pos_label=1)
           fpr, tpr, thresholds = roc_curve(Y_test, PredictedOutput, pos_label=1)
           test_auc = auc(fpr, tpr)
           print("Best accuracy on validation set is:", best_score)
          print("Best parameter for regularization (C) is: ", best_parameters)
print("Test accuracy with best C parameter is", test_score)
print("Test recall with the best C parameter is", test_recall)
          print("Test AUC with the best C parameter is", test_auc)
           m = 'Logistic Regression (w/ dropna)'
          acc.append([m, test_score, test_recall, test_recall, fpr, tpr, thresholds])
           Best accuracy on validation set is: 0.725974025974026
           Best parameter for regularization (C) is: 10
           Test accuracy with best C parameter is 0.80555555555556
           Test recall with the best C parameter is 0.75
```

Test AUC with the best C parameter is 0.819444444444443

```
In [23]: best score = 0
                    for c_paramter in [0.001, 0.01, 0.1, 1, 10, 100, 1000]:
                            c_parameter in [0.001, 0.01, 0.1, 1, 100, 1000]:
for gamma_parameter in [0.001, 0.01, 0.1, 1, 100, 1000]:
    for k_parameter in ['rbf', 'linear', 'poly', 'sigmoid']:
        symModel = SVC(kernel=k_parameter, C=c_parameter, gamma=gamma_parameter)
        scores = cross_val_score(symModel, X_trainval_scaled, Y_trainval, cv=kfolds, scoring='accuracy')
                                             score = np.mean(scores)
                                             if score > best score:
                                                     best_score = score
                                                     best_parameter_c = c_paramter
                                                     best_parameter_gamma = gamma_paramter
                                                     best_parameter_k = k_parameter
                   SelectedSVM model = SVC(C-best\_parameter\_c, \ gamma=best\_parameter\_gamma, \ kernel=best\_parameter\_k).fit(X\_trainval\_scaled, Y\_trainval\_scaled, Y
                    test score = SelectedSVMmodel.score(X test scaled, Y test)
                    PredictedOutput = SelectedSVMmodel.predict(X_test_scaled)
                    test_recall = recall_score(Y_test, PredictedOutput, pos_label=1)
                    fpr, tpr, thresholds = roc_curve(Y_test, PredictedOutput, pos_label=1)
                   test_auc = auc(fpr, tpr)
                   print("Best accuracy on cross validation set is:", best_score)
                   print("Best parameter for c is: ", best_parameter_c)
                  print("Best parameter for gamma is: ", best_parameter_gamma)
print("Best parameter for kernel is: ", best_parameter_k)
print("Test accuracy with the best parameters is", test_score)
print("Test recall with the best parameters is", test_recall)
                   print("Test recall with the best parameter is", test_auc)
                   m = 'SVM
                   acc.append([m, test score, test recall, test auc, fpr, tpr, thresholds])
                   4
                    Best accuracy on cross validation set is: 0.7687747035573123
                    Best parameter for c is: 100
                    Best parameter for gamma is: 0.1
                   Best parameter for kernel is: rbf
Test accuracy with the best parameters is 0.8157894736842105
                    Test recall with the best parameters is 0.7
                    Test recall with the best parameter is 0.822222222222222
 In [24]: from sklearn.tree import DecisionTreeClassifier
                    best_score = 0
                    for md in range(1, 9):
                             treeModel = DecisionTreeClassifier(random_state=0, max_depth=md, criterion='gini')
                            scores = cross_val_score(treeModel, X_trainval_scaled, Y_trainval, cv=kfolds, scoring='accuracy')
                            score = np.mean(scores)
                            if score > best_score:
                                     best score = score
                                     best parameter = md
                    Selected DTModel = Decision Tree Classifier (\verb|max_depth| = best_parameter|).fit (X_trainval_scaled, Y_trainval_s) \\
                    test score = SelectedDTModel.score(X_test_scaled, Y_test)
                    PredictedOutput = SelectedDTModel.predict(X_test_scaled)
                     test_recall = recall_score(Y_test, PredictedOutput, pos_label=1)
                     fpr, tpr, thresholds = roc_curve(Y_test, PredictedOutput, pos_label=1)
                    test_auc = auc(fpr, tpr)
                    print("Best accuracy on validation set is:", best_score)
print("Best parameter for the maximum depth is: ", best_parameter)
                    print("Test accuracy with best parameter is ", test_score)
print("Test recall with best parameters is ", test_recall)
print("Test AUC with the best parameter is ", test_auc)
                    acc.append([m, test_score, test_recall, test_auc, fpr, tpr, thresholds])
                    Best accuracy on validation set is: 0.7770750988142293
                    Best parameter for the maximum depth is: 1
                     Test accuracy with best parameter is 0.8157894736842105
                     Test recall with best parameters is 0.65
                    Test AUC with the best parameter is 0.825
```

```
In [25]: print("Feature importance: ")
                      np.array([X.columns.values.tolist(), list(SelectedDTModel.feature_importances_)]).T
                       Feature importance:
 Out[25]: array([['M/F', '0.0'],
                                      ['M/F', '0.0'],
['Age', '0.0'],
['EDUC', '0.0'],
['SES', '0.0'],
['MMSE', '1.0'],
['eTIV', '0.0'],
['nWBV', '0.0'],
['ASF', '0.0']], dtype='<U32')
  In [27]: from sklearn.tree import export_graphviz
                           import graphviz
                          dot_data=export_graphviz(SelectedDTModel, feature_names=X_trainval.columns.values.tolist(),out_file=None)
                          graph = graphviz.Source(dot_data)
                          graph
In [28]: best_score = 0
                     for M in range(2, 15, 2):
                             for d in range(1, 9):
   for m in range(1, 9):
                                                forestModel = RandomForestClassifier (n\_estimators=M, max\_features=d, n\_jobs=4, max\_features=d, n\_jobs=4, max\_features=d, max_features=d, ma
                                                                                                                    max depth=m, random state=0)
                                                scores = cross_val_score(forestModel, X_trainval_scaled, Y_trainval, cv=kfolds, scoring='accuracy')
                                                score = np.mean(scores)
                                                if score > best_score:
                                                         best_score = score
                                                         best_M = M
                                                         best d = d
                                                         best m = m
                     SelectedRFModel = RandomForestClassifier(n_estimators=M, max_features=d,
                                                                                                                     max_depth=m, random_state=0).fit(X_trainval_scaled, Y_trainval )
                     PredictedOutput = SelectedRFModel.predict(X test scaled)
                     test_score = SelectedRFModel.score(X_test_scaled, Y_test)
                     test_recall = recall_score(Y_test, PredictedOutput, pos_label=1)
                     fpr, tpr, thresholds = roc_curve(Y_test, PredictedOutput, pos_label=1)
test_auc = auc(fpr, tpr)
print("Best accuracy on validation set is:", best_score)
                     print("Best parameters of M, d, m are: ", best_M, best_d, best_m)
print("Test accuracy with the best parameters is", test_score)
print("Test recall with the best parameters is:", test_recall)
                     print("Test AUC with the best parameters is:", test auc)
                     m = 'Random Forest'
                     acc.append([m, test_score, test_recall, test_auc, fpr, tpr, thresholds])
                     Best accuracy on validation set is: 0.8035573122529645
                     Best parameters of M, d, m are: 2 5 7
Test accuracy with the best parameters is 0.868421052631579
                     Test recall with the best parameters is: 0.8
                     Test AUC with the best parameters is: 0.872222222222222
  In [29]: print("Feature importance: ")
                       np.array([X.columns.values.tolist(), list(SelectedRFModel.feature_importances_)]).T
                        Feature importance:
```

```
In [30]: best_score = 0
                      for M in range(2, 15, 2):
                               for lr in [0.0001, 0.001, 0.01, 0.1, 1]:
                                        # train the model
                                        boostModel = AdaBoostClassifier(n_estimators=M, learning_rate=lr, random_state=0)
                                        scores = cross_val_score(boostModel, X_trainval_scaled, Y_trainval, cv=kfolds, scoring='accuracy')
                                        score = np.mean(scores)
                                        if score > best_score:
                                                  best_score = score
                                                 best M = M
                                                 best lr = lr
                      Selected Boost Model = AdaBoost Classifier (n\_estimators=M, learning\_rate=lr, random\_state=0).fit (X\_trainval\_scaled, Y\_trainval\_) (In the first of the first o
                      PredictedOutput = SelectedBoostModel.predict(X test scaled)
                     test_score = SelectedRFModel.score(X_test_scaled, Y_test)
test_recall = recall_score(Y_test, PredictedOutput, pos_label=1)
                      fpr,\ tpr,\ thresholds\ =\ roc\_curve(Y\_test,\ PredictedOutput,\ pos\_label=1)
                     rest_auc = auc(fpr, tpr)

print("Best accuracy on validation set is:", best_score)

print("Best parameter of M is: ", best_M)

print("best parameter of LR is: ", best_Ir)
                     print("Test accuracy with the best parameter is", test_score)
print("Test recall with the best parameters is:", test recall)
                      print("Test AUC with the best parameters is:", test_auc)
                      m = 'AdaBoost'
                     acc.append([m, test_score, test_recall, test_auc, fpr, tpr, thresholds])
                      Best accuracy on validation set is: 0.7770750988142293
                      Best parameter of M is: 2
best parameter of LR is: 0.0001
                      Test accuracy with the best parameter is 0.868421052631579
                      Test recall with the best parameters is: 0.65
                      Test AUC with the best parameters is: 0.825
   In [31]: print("Feature importance: ")
                        np.array([X.columns.values.tolist(), list(SelectedBoostModel.feature_importances_)]).T
                        Feature importance:
   Out[31]: array([['M/F', '0.07142857142857142'],
                                        [['M/F', '0.07142857142857142'],
['Age', '0.14285714285714285'],
['EDUC', '0.21428571428571427'],
['SES', '0.071428571428571425'],
['MNSE', '0.14285714285714285'],
['eTIV', '0.21428571428571427'],
['nNBV', '0.14285714285714285'],
['ASF', '0.0']], dtype='<U32')
   In [32]: result = pd.DataFrame(acc, columns=['Model', 'Accuracy', 'Recall', 'AUC', 'FPR', 'TPR', 'TH'])
                        result[['Model', 'Accuracy', 'Recall', 'AUC']]
   Out[32]:
                                                                             Model Accuracy Recall
                                                                                                                                    AUC
                         0 Logistic Regression (w/ imputation) 0.763158 0.70 0.766667
                          1 Logistic Regression (w/ dropna) 0.805556 0.75 0.750000
                                                                          SVM 0.815789 0.70 0.822222
                                                               Decision Tree 0.815789 0.65 0.825000
                          3
                                                              Random Forest 0.868421 0.80 0.872222
                          4
                          5
                                                                       AdaBoost 0.868421 0.65 0.825000
```

### Github Links

## Project:

Vaishnav2535/Detecting-Early-Alzheimer-s-using-Semi---Supervised-Learning (github.com)

### Code:

https://github.com/Vaishnav2535/Detecting-Early-Alzheimer-s-using-Semi---Supervised-Learning/blob/fa2dab865fd5762280428e1bfaccfb74a5ee23dd/Detecting%20Early%20Alzheimers.ipynb

Detecting Early Alzheimers.ipynb