AdaBoost - Lab

September 16, 2021

Modify the AdaBoost scratch code in our lecture such that: - Notice that if $\operatorname{err}=0$, then α will be undefined, thus attempt to fix this by adding some very small value to the lower term - Notice that sklearn version of AdaBoost has a parameter learning_rate. This is in fact the $\frac{1}{2}$ in front of the α calculation. Attempt to change this $\frac{1}{2}$ into a parameter called eta, and try different values of it and see whether accuracy is improved. Note that sklearn default this value to 1. - Observe that we are actually using sklearn DecisionTreeClassifier. If we take a look at it closely, it is actually using weighted gini index, instead of weighted errors that we learn above. Attempt to write your own class of class Stump that actually uses weighted errors, instead of weighted gini index. To check whether your stump really works, it should give you still relatively the same accuracy. In addition, if you do not change y to -1, it will result in very bad accuracy. Unlike sklearn version of DecisionTree, it will STILL work even y is not change to -1 since it uses gini index - Put everything into a class

```
[1]: from sklearn.model_selection import train_test_split
from sklearn.datasets import make_moons
import numpy as np
import matplotlib.pyplot as plt
```

```
[2]: from sklearn.datasets import make_classification

X, y = make_classification(n_samples=500, random_state=1)
y = np.where(y==0,-1,1) #change our y to be -1 if it is 0, otherwise 1

X_train, X_test, y_train, y_test = train_test_split(
    X, y, test_size=0.3, random_state=42)
```

```
[3]: class AdaBoost():
    def __init__(self, S=5, eta=0.5):
        self.S = S
        self.eta = eta
        self.stump_params = {'max_depth': 1, 'max_leaf_nodes': 2}
        self.models = models = [DecisionTreeClassifier(**self.stump_params) for___
    in range(S)]

def fit(self, X, y):
    m = X_train.shape[0]

#initially, we set our weight to 1/m
```

```
self.W = np.full(m, 1/m)
             #keep collection of a_j
             self.a_js = np.zeros(self.S)
             for j, model in enumerate(self.models):
                 #train weak learner
                 model.fit(X_train, y_train, sample_weight = self.W)
                 #compute the errors
                 yhat = model.predict(X_train)
                 err = self.W[(yhat != y_train)].sum()
                 #compute the predictor weight a_j
                 #if predictor is doing well, a_j will be big
                 peb = 1e-5 # prevent divide by zero
                 a_j = np.log ((1 - err) / (err + peb))
                 self.a_js[j] = a_j
                 #update sample weight; divide sum of W to normalize
                 self.W = (self.W * np.exp(-a_j * y_train * yhat))
                 self.W = self.W / sum (self.W)
         def predict(self, X_test):
             #make weighted predictions
             Hx = 0
             for i, model in enumerate(self.models):
                 yhat = model.predict(X_test)
                 Hx += self.a_js[i] * yhat
             yhat = np.sign(Hx)
             return yhat
[4]: class Stump():
         def __init__(self):
             # Determines whether threshold should be evaluated as < or >
             self.polarity = 1
             self.feature_index = None
             self.threshold = None
             # Voting power of the stump
             self.alpha = None
[5]: class AdaBoostStump():
         def init (self, S=5, eta=0.5):
             self.S = S
             self.eta = eta
```

```
self.clfs = []
   def fit(self, X, y):
       m, n = X.shape
       #initially, we set our weight to 1/m
       self.W = np.full(m, 1/m)
       #keep collection of a j
       self.a_js = np.zeros(self.S)
       for _ in range(self.S):
           clf = Stump()
           #set initially minimum error to infinity
           #so at least the first stump is identified
           min_err = np.inf
           #previously we don't need to do this
           #since sklearn learn does it
           #but now we have to loop all features, all threshold
           #and all polarity to find the minimum weighted errors
           for feature in range(n):
               feature_vals = np.sort(np.unique(X[:, feature]))
               thresholds = (feature_vals[:-1] + feature_vals[1:])/2
               for threshold in thresholds:
                   for polarity in [1, -1]:
                       yhat = np.ones(len(y)) #set all to 1
                       yhat[polarity * X[:, feature] < polarity * threshold] =__</pre>
\hookrightarrow-1 #polarity=1 rule
                       err = self.W[(yhat != y)].sum()
                       #save the best stump
                       if err < min_err:</pre>
                            clf.polarity = polarity
                            clf.threshold = threshold
                            clf.feature_index = feature
                            min_err = err
           #compute the predictor weight a_j
           #if predictor is doing well, a_j will be big
           peb = 1e-5 # prevent divide by zero
           a_j = np.log ((1 - err) / (err + peb))
           clf.alpha = a_j
           #update sample weight; divide sum of W to normalize
           self.W = (self.W * np.exp(-clf.alpha * y_train * yhat))
```

```
[6]: from sklearn.metrics import accuracy_score
    from sklearn.tree import DecisionTreeClassifier
    from sklearn.metrics import classification_report

model = AdaBoost(S=20, eta=0.5)
    model.fit(X_train, y_train)
    yhat = model.predict(X_test)
    print(classification_report(y_test, yhat))
```

support	f1-score	recall	precision	
79	0.95	0.99	0.92	-1
71	0.94	0.90	0.98	1
150	0.05			
150 150	0.95 0.95	0.94	0.95	accuracy
150	0.95	0.94	0.95	macro avg weighted avg

```
[7]: model = AdaBoostStump(S=20, eta=0.5)
model.fit(X_train, y_train)
yhat = model.predict(X_test)
print(classification_report(y_test, yhat))
```

support	f1-score	recall	precision	
79	0.94	0.95	0.94	-1
71	0.94	0.93	0.94	1
150	0.94			accuracy
150	0.94	0.94	0.94	macro avg
150	0.94	0.94	0.94	weighted avg

```
[8]: from sklearn.ensemble import AdaBoostClassifier

#SAMME.R - a variant of SAMME which relies on class probabilities
#rather than predictions and generally performs better
ada_clf = AdaBoostClassifier(
    DecisionTreeClassifier(max_depth=1), n_estimators=200,
    learning_rate=0.07, random_state=42)
ada_clf.fit(X_train, y_train)
y_pred = ada_clf.predict(X_test)
print("Ada score: ", accuracy_score(y_test, y_pred))
```

Ada score: 0.97333333333333334

[]: