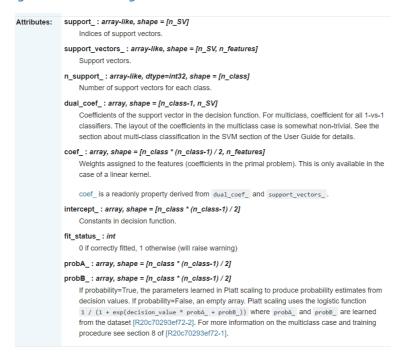
8E and 8F: Finding the Probability P(Y==1|X)

8E: Implementing Decision Function of SVM RBF Kernel

After we train a kernel SVM model, we will be getting support vectors and their corresponsing coefficients a_i Check the documentation for better understanding of these attributes:

https://scikit-learn.org/stable/modules/generated/sklearn.svm.SVC.html



As a part of this assignment you will be implementing the $[decision_function()]$ of kernel SVM, here decision_function() means based on the value return by $[decision_function()]$ model will classify the data point either as positive or negative

Ex 1: In logistic regression After traning the models with the optimal weights w we get, we will find the value $11+e\times p(-(w\times +b))$, if this value comes out to be < 0.5 we will mark it as negative class, else its positive class

Ex 2: In Linear SVM After training the models with the optimal weights w we get, we will find the value of sign(wx + b), if this value comes out to be -ve we will mark it as negative class, else its positive class.

Similarly in Kernel SVM After traning the models with the coefficients a_i we get, we will find the value of $sign(\Sigma_{ni=1}(y_ia_iK(x_i,x_q)) + intercept)$, here $K(x_i,x_q)$ is the RBF kernel. If this value comes out to be -ve we will mark x_q as negative class, else its positive class.

RBF kernel is defined as: $K(x_i, x_q) = \exp(-\gamma ||x_i - x_q||^2)$

For better understanding check this link: https://scikit-learn.org/stable/modules/svm.html#svm-mathematical-formulation font-better understanding check this link: https://scikit-learn.org/stable/modules/svm.html#svm-mathematical-formulation font-better https://scikit-learn.org/stable/modules/svm.html#svm-mathematical-formulation font-better https://scikit-learn.org/stable/modules/svm.html#svm-mathematical-formulation font-better font-better https://scikit-learn.org/stable/modules/svm.html#svm-mathematical-formulation font-better font-better

Task E

- 1. Split the data into $X_{train}(60)$, $X_{cv}(20)$, $X_{test}(20)$
- 2. Train SVC(gamma = 0.001, C = 100.) on the (X_{train} , Y_{train})
- 3. Get the decision boundry values f_{cv} on the X_{cv} data i.e. $f_{cv} = \text{decision_function}(X_{cv})$

```
In [82]:
import numpy as np
import pandas as pd
from sklearn.datasets import make classification
from sklearn.svm import SVC
X, y = make classification(n samples=5000, n features=5, n redundant=2,
                               n classes=2, weights=[0.7], class sep=0.7, random state=1
5)
X.shape
Out[83]:
(5000, 5)
Pseudo code
clf = SVC(gamma=0.001, C=100.)
clf.fit(Xtrain, ytrain)
def decision_function(Xcv, ...): #use appropriate parameters
   for a data point x_q in Xcv:
      #write code to implement (\Sigma_{\text{all the support vectors}} i=1(y_i \alpha_i K(x_i, x_q)) + intercept), here the values y_i, \alpha_i,
and intercept can be obtained from the trained model
return # the decision_function output for all the data points in the Xcv
fcv = decision_function(Xcv, ...) # based on your requirement you can pass any other parameters
Note: Make sure the values you get as fcv, should be equal to outputs of clf.decision_function(Xcv)
In [84]:
# you can write your code here
from sklearn.model selection import train test split
X train, X test, Y train, Y test = train test split(X, y, test size=0.2)
X tr,X cv,Y tr,Y cv = train test split(X train,Y train,test size=0.25)
In [85]:
def decision function(Xcv, sv, y alp, intercept, gam):
  dec fn = []
  #print(sv)
  #print(gam)
  #print(len(Xcv))
  for i in range(len(Xcv)):
    sum = 0
    for j in range(len(sv)):
      sum += (y_alp[j]*np.exp(-1*np.sum((Xcv[i,:]-sv[j,:])**2)*gam))
    dec fn.append((sum+intercept).tolist())
  dec fn = [i[0] for i in dec fn]
  return dec fn
```

```
In [86]:

gam = 0.001
clf = SVC(gamma= gam, C=100)
clf.fit(X tr, Y tr)
```

```
print(len(clf.support_))
y_alpha = clf.dual_coef_
intercept = clf.intercept_
sv = clf.support_vectors_
#print((y_alpha))
dec_fn_implem = decision_function(X_cv,sv,y_alpha[0],intercept,gam)
dec_fn_inbuilt = clf.decision_function(X_cv)
```

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```
In [87]:
```

```
print("Function based decision function values are", dec_fn_implem[:5])
print("sklearn based decision function values are ", dec_fn_inbuilt.tolist()[:5])
```

```
Function based decision function values are [-4.161095825122163, 1.7233399647584635, -2.6499005585185555, 1.3091961267094492, -0.3664902922009814] sklearn based decision function values are [-4.161095825122163, 1.7233399647584635, -2.6499005585185555, 1.3091961267094492, -0.3664902922009814]
```

we can observe that both implementations result in the same values of decision function.

From above we can observe that the function implementation result and inbuilt scikitlearn function results are same.

8F: Implementing Platt Scaling to find P(Y==1|X)

Check this PDF

Let the output of a learning method be f(x). To get calibrated probabilities, pass the output through a sigmoid:

$$P(y = 1|f) = \frac{1}{1 + exp(Af + B)}$$
 (1)

where the parameters A and B are fitted using maximum likelihood estimation from a fitting training set (f_i, y_i) . Gradient descent is used to find A and B such that they are the solution to:

$$\underset{A,B}{argmin} \{ -\sum_{i} y_{i} log(p_{i}) + (1 - y_{i}) log(1 - p_{i}) \}, \quad (2)$$

where

$$p_i = \frac{1}{1 + exp(Af_i + B)} \tag{3}$$

Two questions arise: where does the sigmoid train set come from? and how to avoid overfitting to this training set?

If we use the same data set that was used to train the model we want to calibrate, we introduce unwanted bias. For example, if the model learns to discriminate the train set perfectly and orders all the negative examples before the positive examples, then the sigmoid transformation will output just a 0,1 function. So we need to use an independent calibration set in order to get good posterior probabilities. This, however, is not a draw back, since the same set can be used for model and parameter selection.

To avoid overfitting to the sigmoid train set, an out-ofsample model is used. If there are N_+ positive examples and N_- negative examples in the train set, for each training example Platt Calibration uses target values u_+ and u_- (instead of 1 and 0, respectively), where

$$y_{+} = \frac{N_{+} + 1}{N_{+} + 2}; \ y_{-} = \frac{1}{N_{-} + 2}$$
 (4)

For a more detailed treatment, and a justification of these particular target values see (Platt, 1999).

TASK F

1. Apply SGD algorithm with (f_{CV}, y_{CV}) and find the weight W intercept b Note: here our data is of one dimensional so we will have a one dimensional weight vector i.e W.shape (1,)

Note1: Don't forget to change the values of y_{CV} as mentioned in the above image. you will calculate y+, y- based on data points in train data

Note2: the Sklearn's SGD algorithm doesn't support the real valued outputs, you need to use the code that was done in the 'Logistic Regression with SGD and L2' Assignment after modifying loss function, and use same parameters that used in that assignment.

```
def log_loss(w, b, X, Y):
    N = len(X)
    sum_log = 0
    for i in range(N):
        sum_log += {Y[i]*np.log10(sig(w, X[i], b)) + (1-Y[i])*np.log10(1-sig(w, X[i], b))
    return -1*sum_log/N
```

if Y[i] is 1, it will be replaced with y+ value else it will replaced with y- value

1. For a given data point from X_{test} , $P(Y = 1 | X) = \frac{11 + e^{-(W * f_{test} + b)}}{e^{-(W * f_{test} + b)}}$ where $f_{test} = \frac{11 + e^{-(W * f_{test} + b)}}{e^{-(W * f_{test} + b)}}$ where $f_{test} = \frac{11 + e^{-(W * f_{test} + b)}}{e^{-(W * f_{test} + b)}}$ where $f_{test} = \frac{11 + e^{-(W * f_{test} + b)}}{e^{-(W * f_{test} + b)}}$

```
In [88]:
```

```
def logloss(y_true,y_pred):
    '''In this function, we will compute log loss '''

    n = len(y_true)
    x = np.log10(y_pred)
    x1 = np.log10(np.ones_like(y_pred) - y_pred)
    loss = 0
    for j in range(n):
        loss = loss +((y_true[j]*x[j])+((1-y_true[j])*x1[j]))
    loss*=(-1/n)
    return loss
```

In [89]:

```
def initialize_weights(dim):
    ''' In this function, we will initialize our weights and bias'''
    #initialize the weights to zeros array of (1,dim) dimensions
    #you use zeros_like function to initialize zero, check this link https://docs.sci
py.org/doc/numpy/reference/generated/numpy.zeros_like.html
    #initialize bias to zero

w = np.zeros_like(dim)
b = 0

return w,b
```

```
def sigmoid(z):
    ''' In this function, we will return sigmoid of z'''
    # compute sigmoid(z) and return
    sigma = (1/(1+np.exp(-z)))
    return sigma
In [91]:
```

```
def gradient_dw(x,y,w,b,alpha,N):
    '''In this function, we will compute the gardient w.r.to w '''
    dw = (x*(y - sigmoid(np.matmul(w,x) + b))) - ((alpha/N)*w)
    return dw
```

In [92]:

```
def gradient_db(x,y,w,b):
    '''In this function, we will compute gradient w.r.to b '''
    db = (y - sigmoid(np.matmul(w,x) + b))
    return db
```

In [93]:

```
def train(X train, y train, epochs, alpha, eta0):
   ''' In this function, we will implement logistic regression'''
    #Here eta0 is learning rate
   #implement the code as follows
    # initalize the weights (call the initialize weights(X train[0]) function)
    # for every epoch
        # for every data point(X_train,y_train)
           #compute gradient w.r.to w (call the gradient dw() function)
           #compute gradient w.r.to b (call the gradient db() function)
           #update w, b
        # predict the output of x train[for all data points in X train] using w,b
        #compute the loss between predicted and actual values (call the loss function
        # store all the train loss values in a list
        # predict the output of x test[for all data points in X test] using w,b
        #compute the loss between predicted and actual values (call the loss function
        # store all the test loss values in a list
        # you can also compare previous loss and current loss, if loss is not updatin
g then stop the process and return w,b
   w,b = initialize weights(X train[0])
   N = len(X train)
   loss train = []
   for i in range(epochs):
      w prev,b prev = w,b
     y pred train = []
      for j in range(len(X train)):
       dw = gradient dw(X train[j,:], y train[j], w, b, alpha, N)
        db = gradient_db(X_train[j,:], y_train[j], w, b)
       w = w + (eta0 * dw)
       b = b + (eta0 * db)
      for j in range(len(X train)):
        y pred train.append(sigmoid(np.matmul(w, X train[j,:]) + b))
      present=logloss(y train, y pred train)
      loss train.append(present)
```

```
#print(present)

if(i!=0 and loss_train[i-1]<present):
    print("Minimum loss achieved is",loss_train[i-1])
    return w_prev,b_prev,loss_train[:i]
print("Minimum loss achieved is",present)
return w,b,loss_train</pre>
```

In [94]:

```
N_pos = np.count_nonzero(Y_tr)
N_neg = len(Y_tr)-N_pos
y_pos = (N_pos+1)/(N_pos+2)
y_neg = (1/(N_neg+2))
#print(N_pos,N_neg,y_pos,y_neg)
print(Y_cv.tolist()[:4])
Y_cv_mod = np.zeros_like(Y_cv,dtype =float)
for i in range(len(Y_cv)):
    if Y_cv[i] == 1:
        Y_cv_mod[i] = y_pos
    else:
        Y_cv_mod[i] = y_neg
print(Y_cv_mod.tolist()[:4])
```

```
[0, 1, 0, 1]
[0.00048192771084337347, 0.9989235737351991, 0.00048192771084337347, 0.998923573735199
1]
```

In the above results, list-1 indicates true or actual labels of Y_cv . list-2 indicates the modified labels of Y_cv according to the given formulae.

We can observe that 0 label is replaced with 0.0004766 and label 1 is replaced with 0.9988962

In [95]:

```
alpha=0.0001
eta0=0.001
N=len(X_tr)
epochs=1000
#print(np.array(dec_fn_implem).reshape(-1,1),np.array(Y_cv_mod).reshape(-1,1))
f_cv = np.array(dec_fn_implem).reshape(-1,1)
Y_cv_mod = np.array(Y_cv_mod).reshape(-1,1)
w,b,train_loss = train(f_cv,Y_cv_mod,epochs,alpha,eta0)
```

Minimum loss achieved is [0.07559336]

In [96]:

```
import matplotlib.pyplot as plt

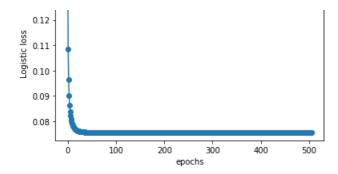
epochs=[i for i in range(len(train_loss))]
plt.plot(epochs,train_loss,label= 'Train Loss')

plt.scatter(epochs,train_loss)

plt.legend()
plt.title('Epochs vs Logistic Loss graph')
plt.xlabel('epochs')
plt.ylabel('Logistic loss')
plt.show()
```

```
0.14 Epochs vs Logistic Loss graph

— Train Loss
```



In [97]:

```
f_test = decision_function(X_test, sv, y_alpha[0], intercept, gam)
print(f_test[:4])
proba_1=[]
for i in range(len(X_test)):
    proba_1.append(sigmoid(np.dot(w, f_test[i]) + b))

proba_1 = [i.tolist() for i in proba_1]
proba_1 = [i[0] for i in proba_1]
print(proba_1[:4])
```

[-2.392145229826883, -2.9514350778685126, -1.6627732024593085, -4.619833147669826] [0.011427693208354434, 0.003914680589311406, 0.04507717531280103, 0.000157258920370584 34]

we can observe that if decision function value is negative it's probability of being class 1 P(y = 1/x) < 0.5 and is low and if it is positive it's P(y = 1/x) > 0.5 and high.

Note: in the above algorithm, the steps 2, 4 might need hyper parameter tuning, To reduce the complexity of the assignment we are excluding the hyerparameter tuning part, but intrested students can try that

If any one wants to try other calibration algorithm istonic regression also please check these tutorials

- 1. http://fa.bianp.net/blog/tag/scikit-learn.html#fn:1
- 2. https://drive.google.com/open?id=1MzmA7QaP58RDzocB0RBmRiWfl7Co_VJ7
- 3. https://drive.google.com/open?id=133odBinMOIVb_rh_GQxxsyMRyW-Zts7a
- 4. https://stat.fandom.com/wiki/Isotonic_regression#Pool_Adjacent_Violators_Algorithm