

# **ELL409: MACHINE LEARNING**

## **ASSIGNMENT 3**

Implementation of Support Vector Regression (SVR) using  
CVXOPT and using sklearn

**HARSH KUMAR**  
**2017MT10729**

Dataset is Bosten House Pricing Data, consisting of 506 rows and 13 features and a class containing the Median value of owner-occupied homes to be predicted later.

First will look after SVR implementation using CVXOPT

### Implementation of SVR using CVXOPT:

We need to solve a dual problem in this, as discussed in the class.

$$\begin{aligned}\tilde{L}(\mathbf{a}, \hat{\mathbf{a}}) = & -\frac{1}{2} \sum_{n=1}^N \sum_{m=1}^N (a_n - \hat{a}_n)(a_m - \hat{a}_m)k(\mathbf{x}_n, \mathbf{x}_m) \\ & -\epsilon \sum_{n=1}^N (a_n + \hat{a}_n) + \sum_{n=1}^N (a_n - \hat{a}_n)t_n\end{aligned}$$

We need to maximize this equation with respect to constraints

$$\sum_{n=1}^N (a_n - \hat{a}_n) = 0$$

$$0 \leq a_n \leq C$$

$$0 \leq \hat{a}_n \leq C$$

Where C and epsilon are hyperparameters.

Using the value of “a” and “a^” found we can find the value of y(x), which is as follows:

$$y(\mathbf{x}) = \sum_{n=1}^N (a_n - \hat{a}_n)k(\mathbf{x}, \mathbf{x}_n) + b$$

Where  $b$  is

$$\begin{aligned} b &= t_n - \epsilon - \mathbf{w}^T \phi(\mathbf{x}_n) \\ &= t_n - \epsilon - \sum_{m=1}^N (a_m - \hat{a}_m) k(\mathbf{x}_n, \mathbf{x}_m) \end{aligned}$$

Then we took the average of all the  $b$  for which points are support vectors.

The above dual problem is solved using a QP Solver, whose standard form is given below:

$$\begin{aligned} \min_x \quad & \frac{1}{2} x^T P x + q^T x \\ \text{subject to} \quad & Gx \preceq h \\ & Ax = b \end{aligned}$$

We can get the value of  $x$  using the following equations:

```
sol = solvers.qp(P,q,G,h,A,b)
sol['x']
```

Hence we need to convert our dual problem in the above form. So, took  $x$  as a matrix of  $[a, a^*]$  and then changes each equation accordingly, which is shown below.

- First, data normalization is done.

```
# normalisation of data
data = np.array(data)
data -= np.mean(data, axis=0)
data /= np.std(data, axis=0)
```

- The different kernel function is introduced as:

```
# define the Kernal Function for it is RBF
def kernal_rbf(x, y, gamma):
    k = math.exp((-1)*gamma*(LA.norm(x-y)*(LA.norm(x-y))))
    return k
```

```
# define the polynomial kernel function, d id the degree of freedom pf polynomial and cons is the constant term
def kernal_polynomial(x, y, constant, d):
    k = pow(np.dot(x, y)+constant, d)
    return k
```

```
# define a linear kernal function
def kernal_linear(x, y, cons):
    k = np.dot(x, y)+cons
    return k
```

- Then a kernel matrix is introduced for each kernel function such that  $K = K[X_i, X_j]$  for  $0 \leq i, j \leq n$ , where  $n$  is the number of data vectors.

```
def kernal_matrix_rbf(X, gamma):
    n = len(X);
    k = np.zeros([n, n])
    for i in range(len(X)):
        for j in range(len(X)):
            k[i][j] = kernal_rbf(X[i], X[j], gamma)
    return k
```

```
def kernal_matrix_polynomial(X, constant, d):
    n = len(X);
    k = np.zeros([n, n])
    for i in range(len(X)):
        for j in range(len(X)):
            k[i][j] = kernal_polynomial(X[i], X[j], constant, d)
    return k
```

```
def kernal_matrix_linear(X, cons):
    n = len(X);
    k = np.zeros([n, n])
    for i in range(len(X)):
        for j in range(len(X)):
            k[i][j] = kernal_linear(X[i], X[j], cons)
    return k
```

- The P matrix will be equal to  $\begin{bmatrix} k & -k \\ -k & k \end{bmatrix}$ , which is implemented as below for each of the kernel functions

```
# calculate P matrix for rbf
def P_Matrix_rbf(X, gamma):
    k = kernal_matrix_rbf(X, gamma)
    neg_k = (-1)*k
    temp1 = np.concatenate((k, neg_k))
    temp2 = np.concatenate((neg_k, k))
    P = np.concatenate((temp1, temp2), axis = 1)
    return P
```

```
# calculate P matrix for linear
def P_Matrix_linear(X, cons):
    k = kernal_matrix_linear(X, cons)
    neg_k = (-1)*k
    temp1 = np.concatenate((k, neg_k))
    temp2 = np.concatenate((neg_k, k))
    P = np.concatenate((temp1, temp2), axis = 1)
    return P
```

```
# calculate P matrix for polynomial
def P_Matrix_polynomial(X, constant, d):
    k = kernal_matrix_polynomial(X, constant, d)
    neg_k = (-1)*k
    temp1 = np.concatenate((k, neg_k))
    temp2 = np.concatenate((neg_k, k))
    P = np.concatenate((temp1, temp2), axis = 1)
    return P
```

- Q matrix is implemented as follows:

```
# calculate q matrix
def q_matrix(y, elp):
    temp1 = -y+elp
    temp2 = y+elp
    q = np.concatenate((temp1, temp2))
    return q
```

- G Matrix:

```
# calculate G matrix
def G_matrix(n):
    temp1 = np.identity(2*n)
    temp2 = (-1.0)*temp1
    g = np.concatenate((temp1, temp2))
    return g
```

- H Matrix:

```
# calculate h matrix
def h_matrix(C, n):
    temp1 = np.ones(2*n)
    temp2 = np.zeros(2*n)
    h = np.concatenate((temp1, temp2))
    return h
```

- A matrix:

```
# calculate A matrix
def A_matrix(n):
    temp1 = np.ones(n)
    temp2 = (-1.0)*temp1
    temp = np.concatenate((temp1, temp2))
    temp = temp.reshape(1,-1)
    return temp
```

- B matrix:

```
# calculate b matrix
def b_matrix():
    b = np.zeros(1)
    return b
```



- Now, the prediction function for different kernel function is implemented as follows:

```
# prediction of y for rbf
def Y_predicted_rbf(X, n, lamda_1, lamda_2, X_train, b, gamma):
    total = 0
    for i in range(n):
        total = total+(lamda_1[i][0]-lamda_2[i][0])*kernel_rbf(X, X_train[i], gamma)
    y = total+b
    return y
```

```
# prediction of y for linear
def Y_predicted_linear(X, n, lamda_1, lamda_2, X_train, b, cons):
    total = 0
    for i in range(n):
        total = total+(lamda_1[i][0]-lamda_2[i][0])*kernel_linear(X, X_train[i], cons)
    y = total+b
    return y
```

```
# prediction of y for polynomial
def Y_predicted_polynomial(X, n, lamda_1, lamda_2, X_train, b, constant, d):
    total = 0
    for i in range(n):
        total = total+(lamda_1[i][0]-lamda_2[i][0])*kernel_polynomial(X, X_train[i], constant, d)
    y = total+b
    return y
```

- Math Square Error is calculated as follows:

```
# calculation of Math square Error
def MSE(Y_predict, Y_test, elp):
    total = 0
    for i in range(len(Y_test)):
        total = total + (abs(Y_predict[i]-Y_test[i])-elp)*(abs(Y_predict[i]-Y_test[i])-elp)
    error = total/(len(Y_test))
    return error
```

- Then used 5 fold cross-validation to find the value of y for the test cases and also predict the error.
- The matrices are then initialized as follows:

```
n = len(X_train)
P = P_Matrix_linear(X_train, cons)
q = q_matrix(Y_train, elp)
G = G_matrix(n)
h = h_matrix(C, n)
A = A_matrix(n)
b = b_matrix()
```

- Then converted the matrices in the form of cvxopt matrices:

```
P = cvxopt_matrix(P)
q = cvxopt_matrix(q)
G = cvxopt_matrix(G)
h = cvxopt_matrix(h)
A = cvxopt_matrix(A)
b = cvxopt_matrix(b)
```

- Using the below equations, solved our problem and got our x, which is lamda here, and then split it into lamda\_1(a) and lamda\_2(a^)

```
sol = cvxopt_solvers.qp(P, q, G, h, A, b)
lamda = np.array(sol['x'])
l = int(len(lamda)/2)
lamda_1 = lamda[0:l, :]
lamda_2 = lamda[l:,:]

```

- Calculated the support vector indices as follows:

```
# the supp_vector array will contain all indicies of vectors which are support vectors
supp_vector = []
for j in range(len(lamda_1)):
    temp = lamda_1[j][0]
    if temp>0 and temp<C:
        supp_vector.append(j)
```

- Then calculate the value of b to be used in the prediction of y.

```
# calculation of b
b = 0
for i1 in supp_vector:
    b = b+Y_train[i1]-elp-b_term_linear(X_train[i1], n, lamda_1, lamda_2, X_train, cons)
b = b/(len(supp_vector))
```

- These codes are used for visualization of data points. Since our dataset is not linear so, I took the **first feature** and plotted the points corresponding to that feature.

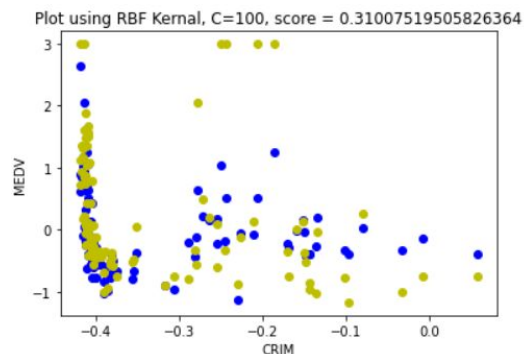
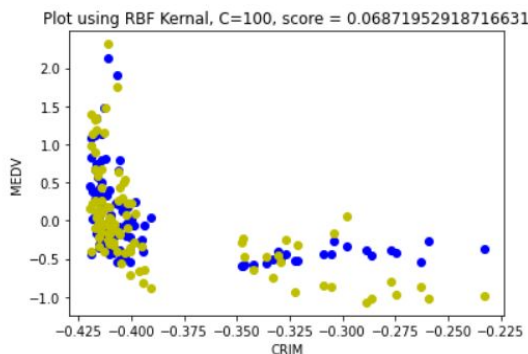


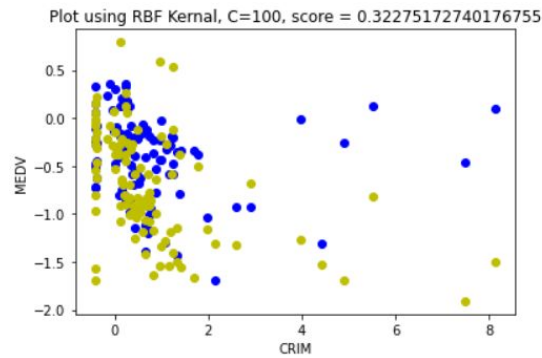
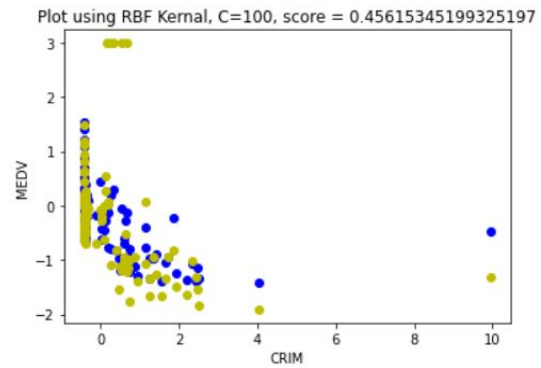
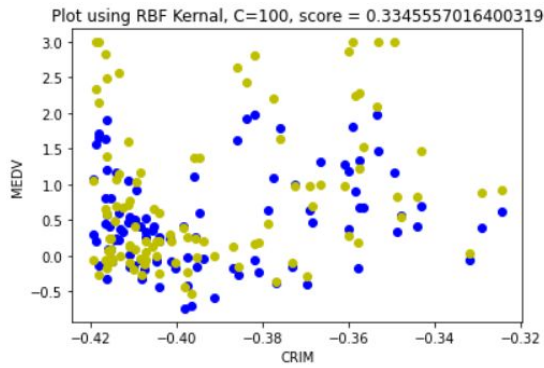
```
# data visualisation
plt.scatter(X_test[:,0], Y_predict_linear, label='Y_predict', color = 'b')
plt.scatter(X_test[:,0], Y_test, label='Y_test', color = 'y')
plt.title("Plot using Linear Kernel, C = 1000, Score = "+str(score_linear[0]))
plt.xlabel("CRIM")
plt.ylabel("MEDV")
plt.show()
```

- The error got in :  
 SVR using RBF Kernel: 0.298 (Using  $c = 100$ ,  $\epsilon = 0.1$ ,  $\gamma = 0.1$ )  
 SVR using Linear Kernel: 0.334 (using  $C = 100$ )  
 SVE using Polynomial Kernel: 17.550 ( $d = 3$ ,  $C = 100$ ,  $\epsilon = 0.1$ )

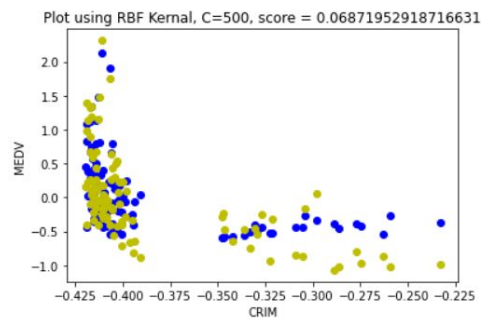
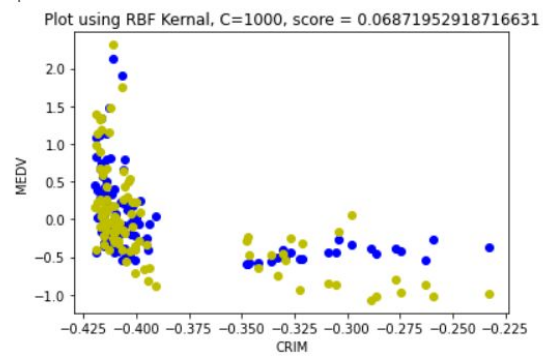
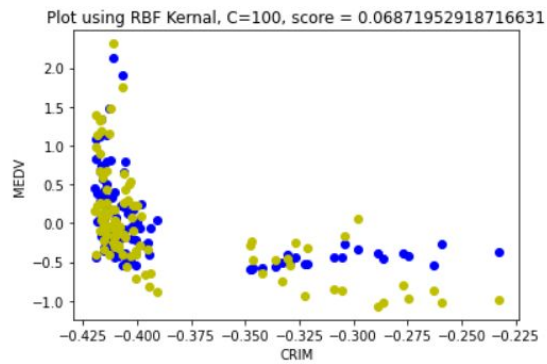
## PLOTTING OF GRAPHS

- Plot for a graph using **RBF Kernel** using  $\gamma = 0.1$ ,  $C = 100$ ,  $\epsilon = 0.1$ , all 5 figures of cross-validation

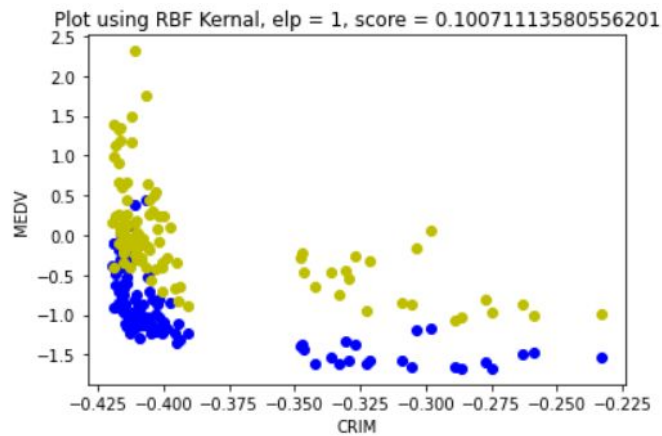
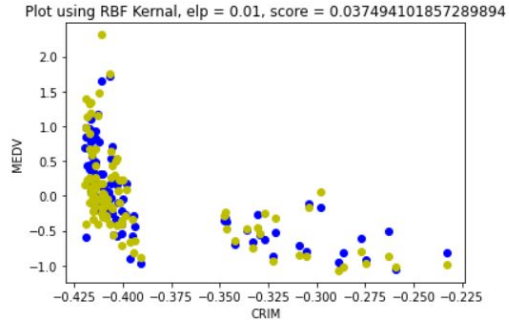
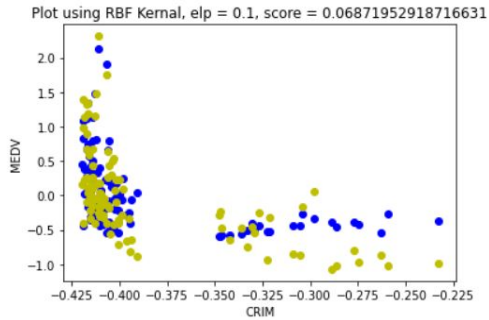




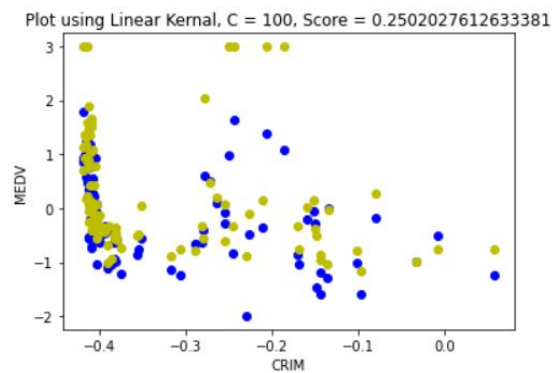
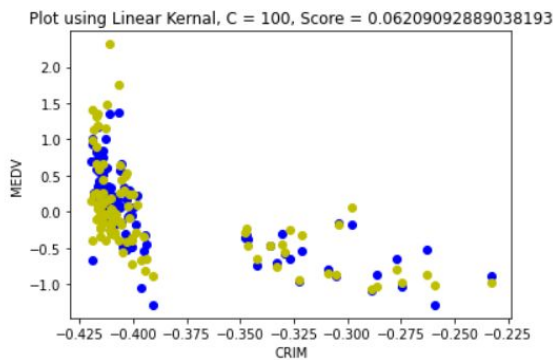
- Plot for different C, keeping epsilon = 0.1, gamma = 0.1

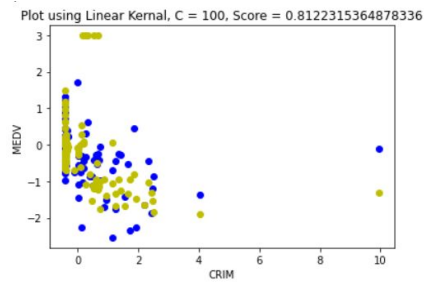
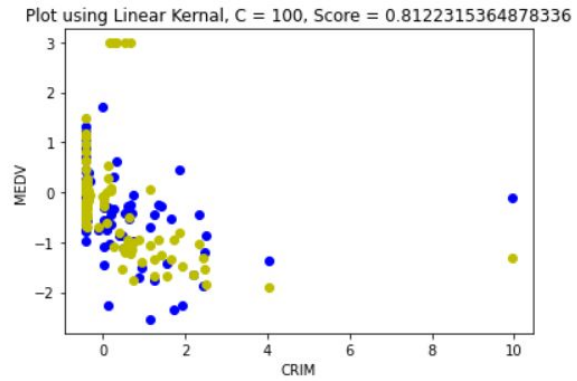
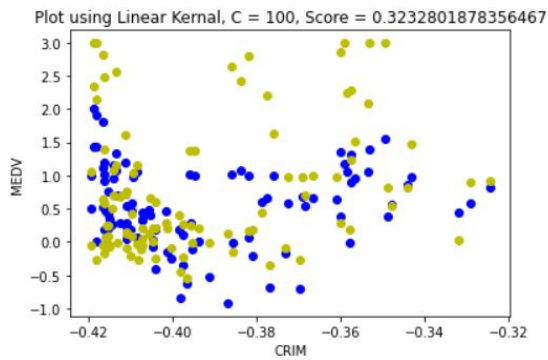


- For different epsilon, keeping gamma = 0.1, C = 100

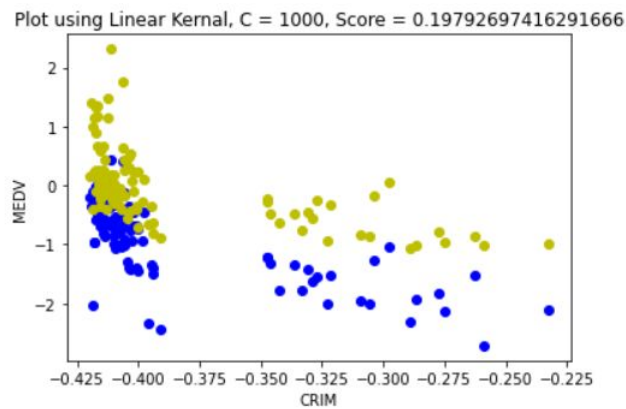
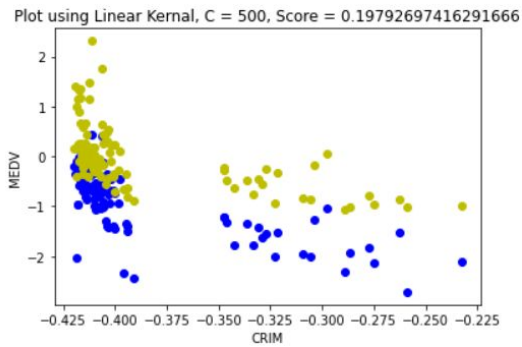
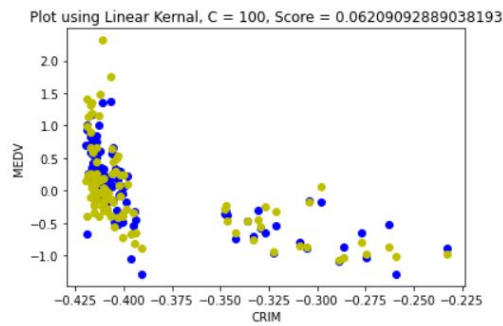


- Plot for a graph using **Linear Kernel** using C = 100

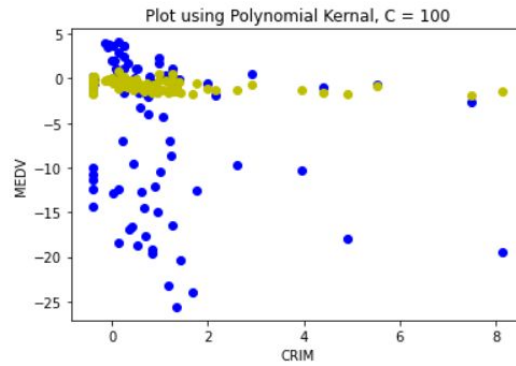
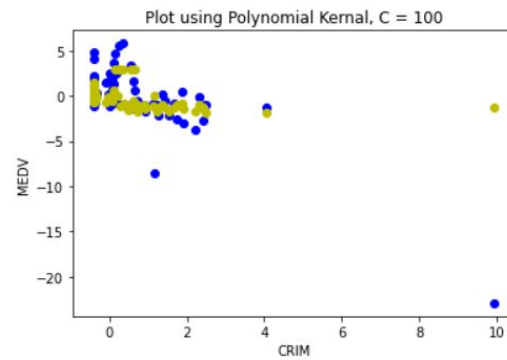
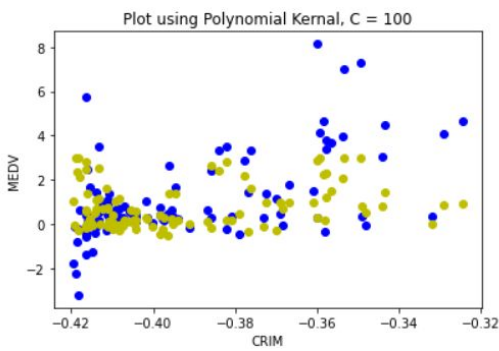
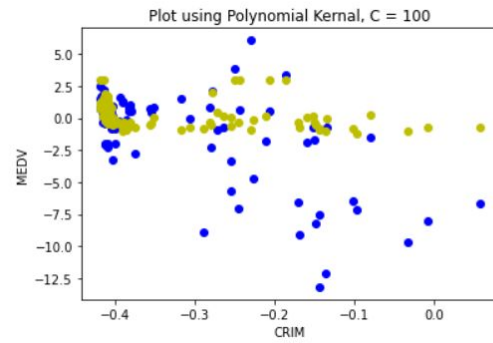
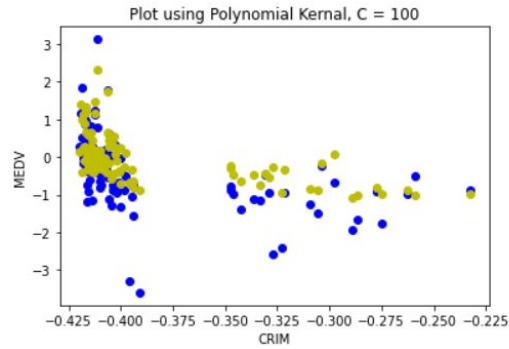




## ● Plot of graph for different C

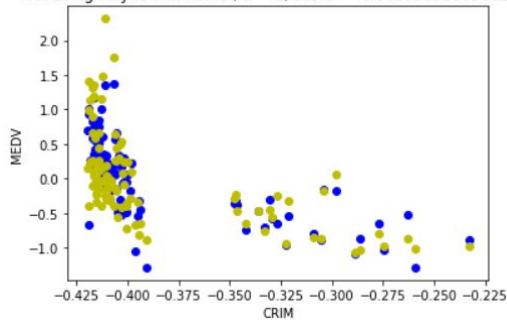


- Plot of Graph for **kernel polynomial** Function

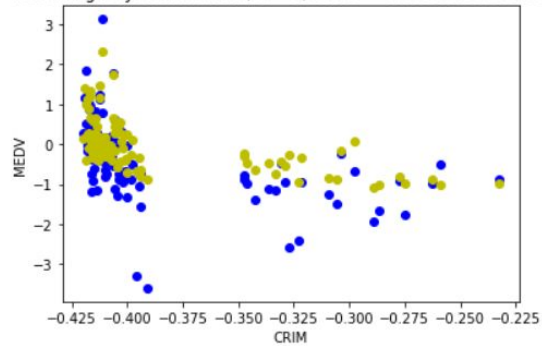


- Polynomial Kernel, for different degree of freedom( $d$ )

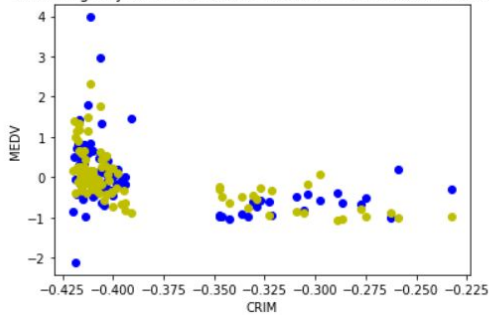
Plot using Polynomial Kernel,  $d = 1$ , score = 0.06209092889062161



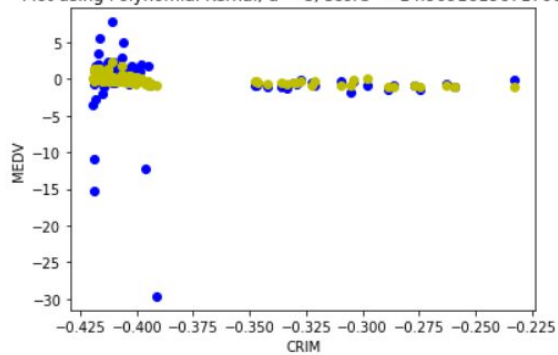
Plot using Polynomial Kernel,  $d = 3$ , score = 0.4632043548588621



Plot using Polynomial Kernel,  $d = 5$ , score = 0.303617285867739

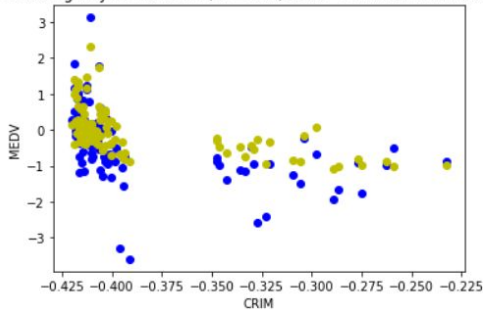


Plot using Polynomial Kernel,  $d = 8$ , score = 14.909161967170652

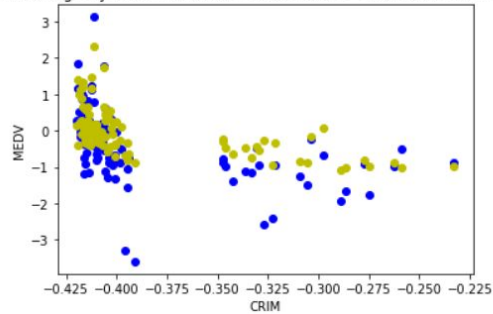


- For different  $C$  keeping  $d = 3$

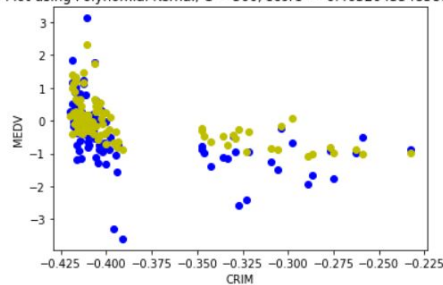
Plot using Polynomial Kernel,  $C = 100$ , score = 0.4632043548588621



Plot using Polynomial Kernel,  $C = 1000$ , score = 0.4632043548588621



Plot using Polynomial Kernel,  $C = 500$ , score = 0.4632043548588621





## SVR Implementation using sklearn

- Took 5 fold cross-validations.
- The following is the implementation:

```
svr_rbf = SVR(kernel='rbf', C=100, gamma=0.1, epsilon=1)
svr_lin = SVR(kernel='linear', C=100, gamma='auto')
svr_poly = SVR(kernel='poly', C=100, gamma='auto', degree=8, epsilon=0.1, coef0=1)
```

- Prediction:

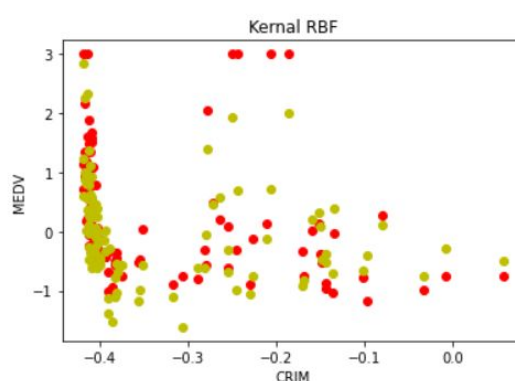
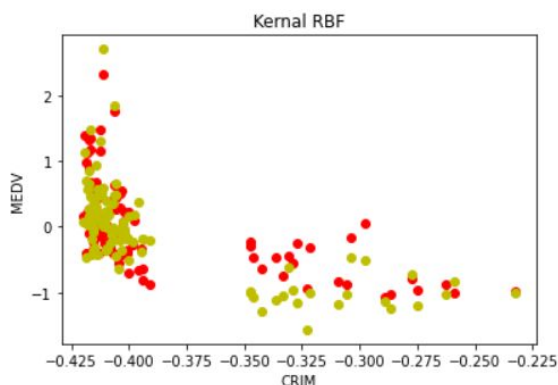
```
svr_poly.fit(X_train, Y_train)
poly_predict = svr_poly.predict(X_test)
scores_poly.append(svr_poly.score(X_test, Y_test))
```

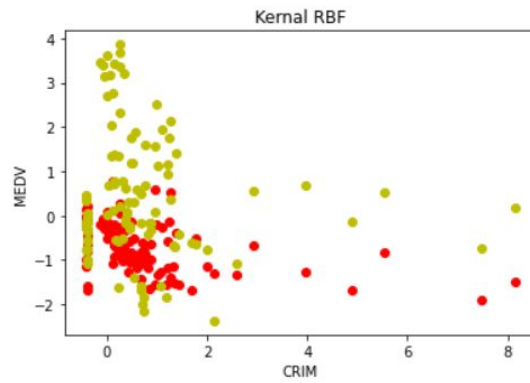
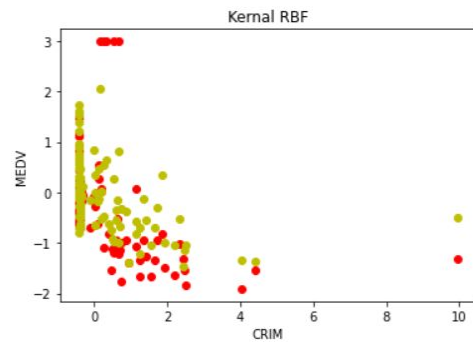
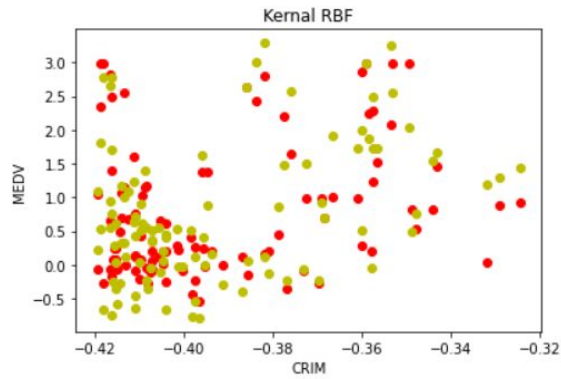
- Codes used for plotting, same plotting as above:

```
plt.title('Kernal Linear, eplison = 0.1, score = '+str(scores_lin[0]))
plt.xlabel('CRIM')
plt.ylabel('MEDV')
plt.scatter(X_test[:,0], Y_test, color = 'g', label="Y_test".format('r'))
plt.scatter(X_test[:,0], lin_predict, color = 'b', label="Y_predicted".format('y'))
plt.show()
```

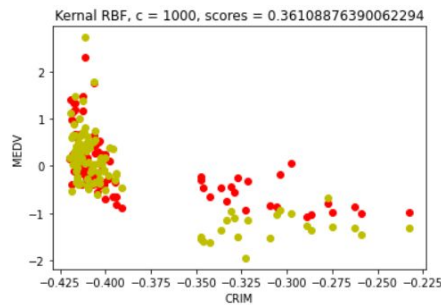
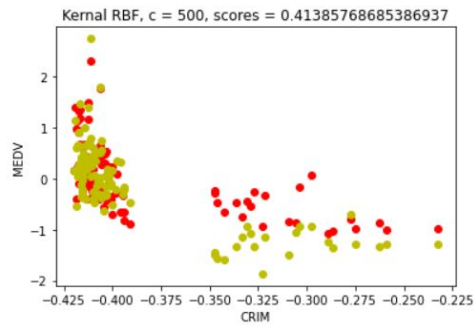
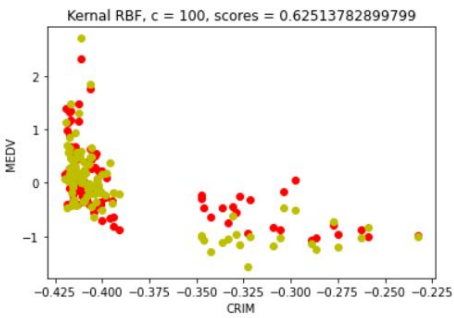
## Plotting of Graphs

- **RBF Kernal** for  $C = 100$ ,  $\gamma = 0.1$ ,  $\epsilon = 0.1$

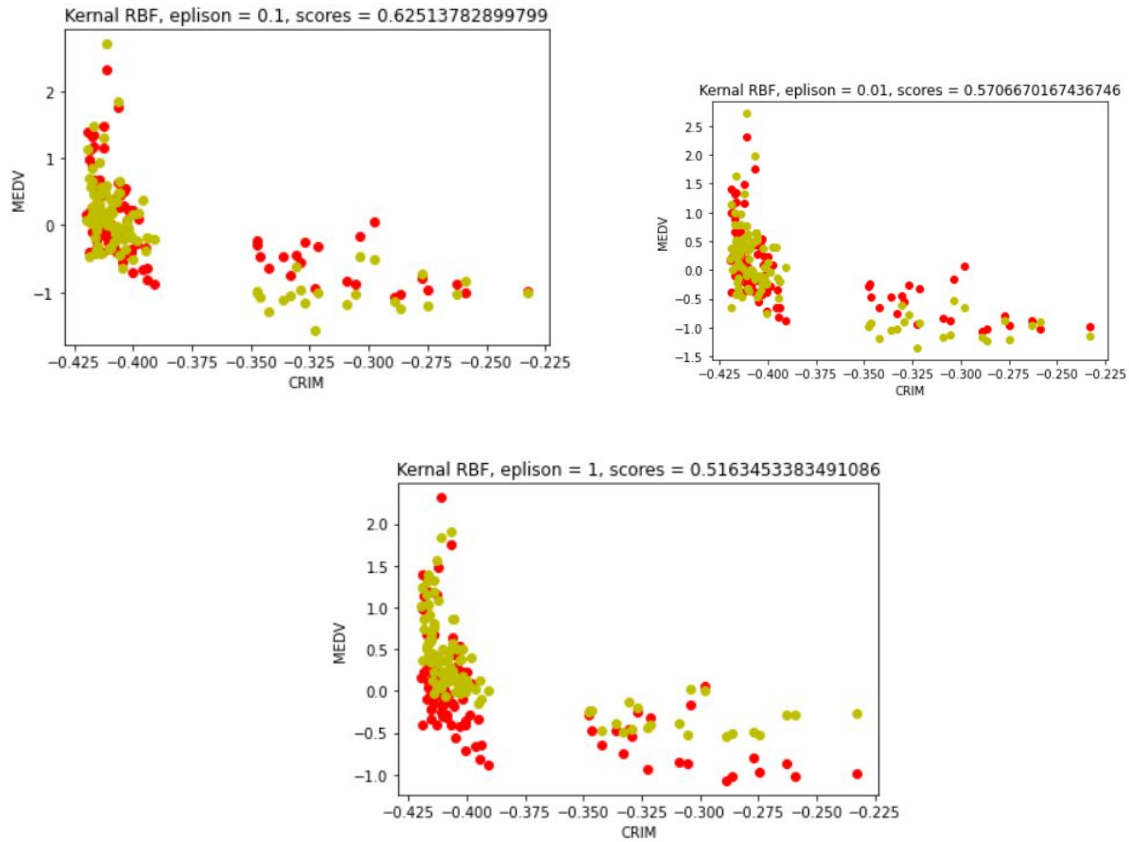




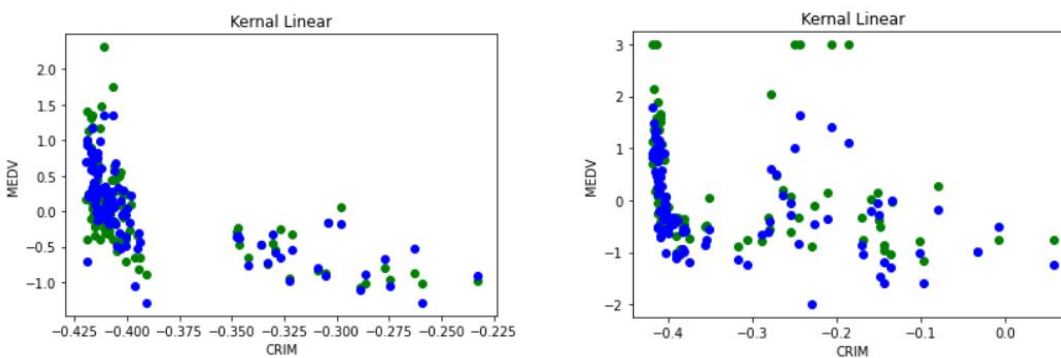
- For different C, keeping other parameters same:

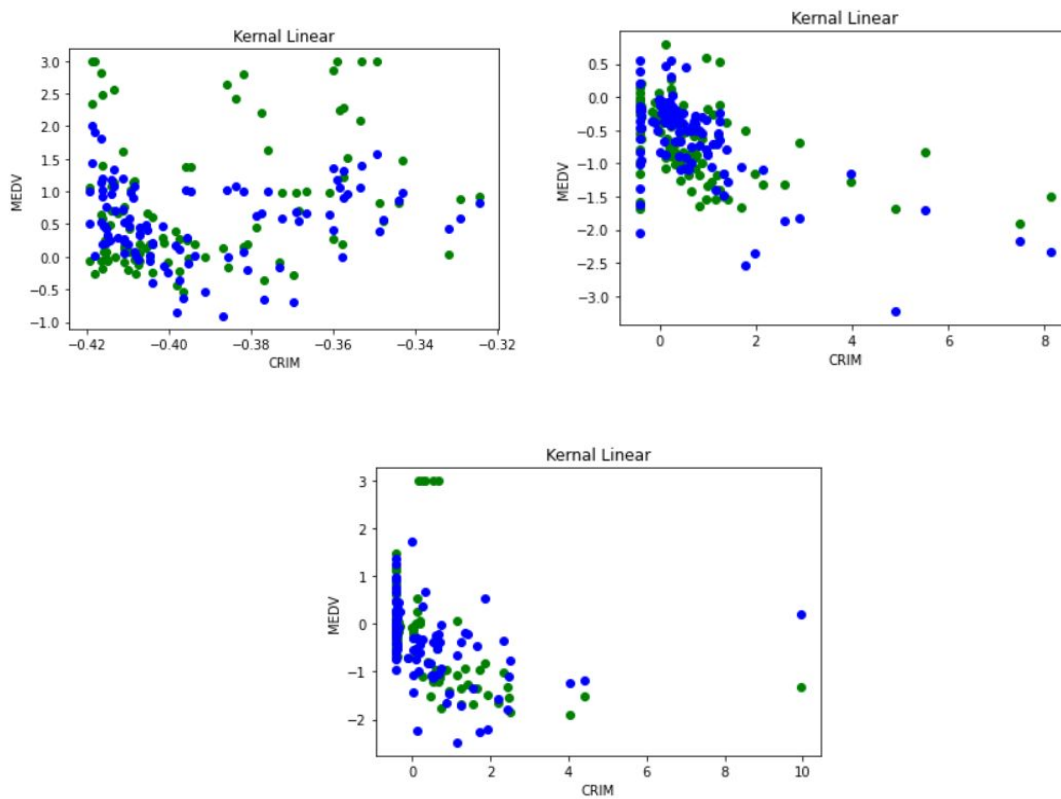


- For different epsilon, keeping  $C = 100$

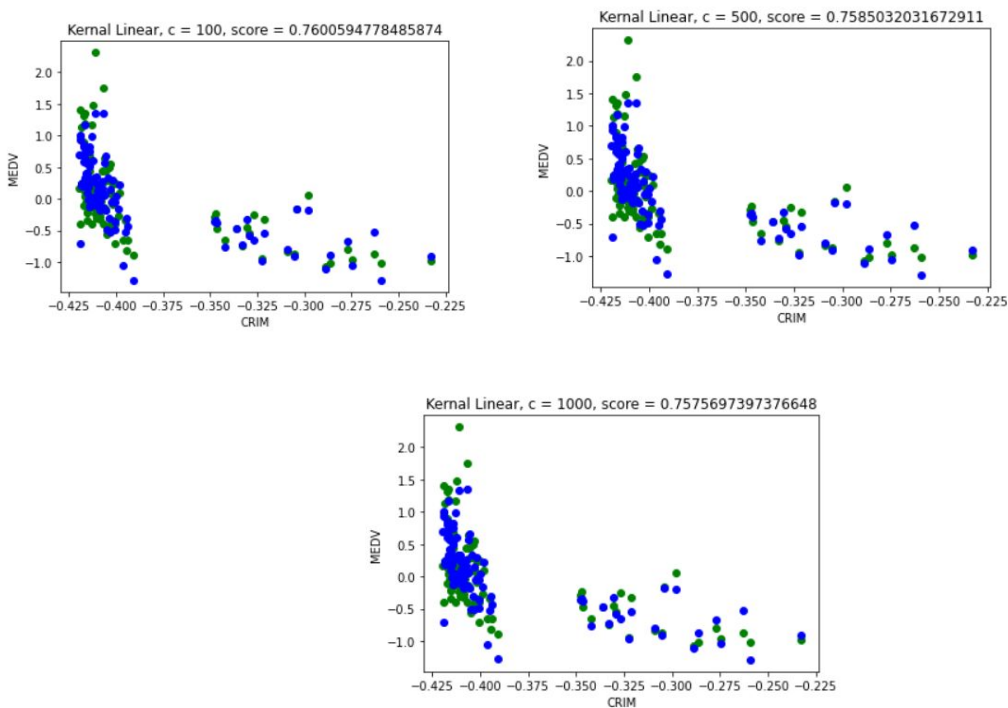


- Plot for **Linear Kernel** for  $C = 100$

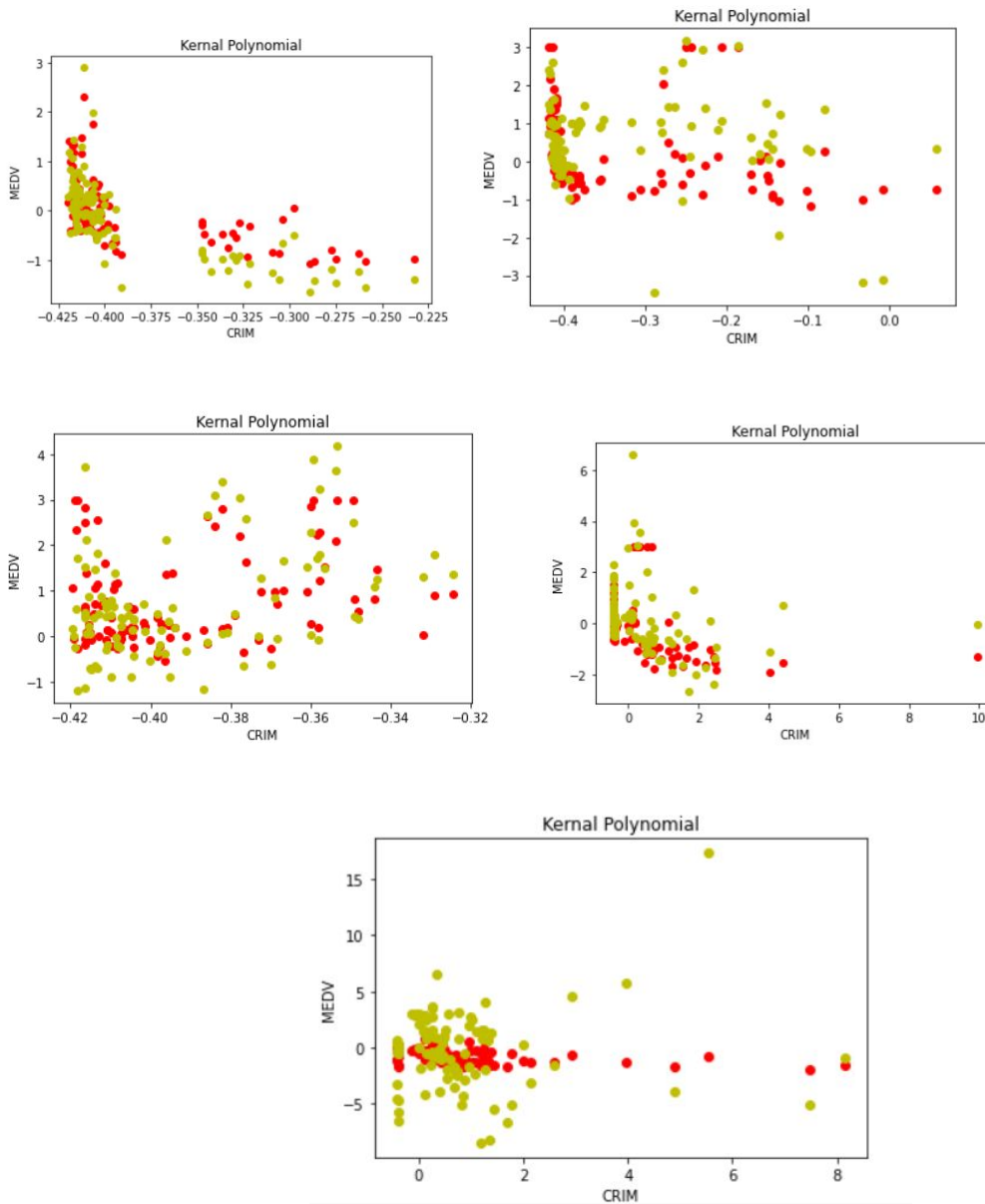




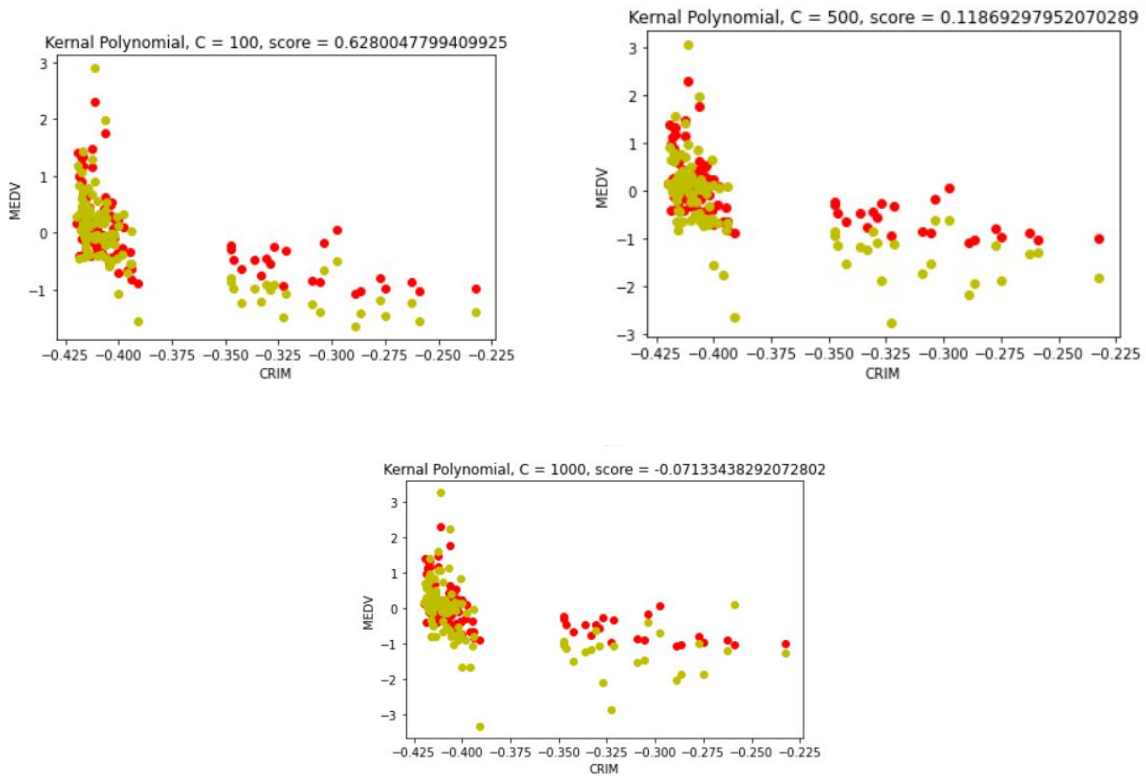
- Plot for Linear Kernel for different C



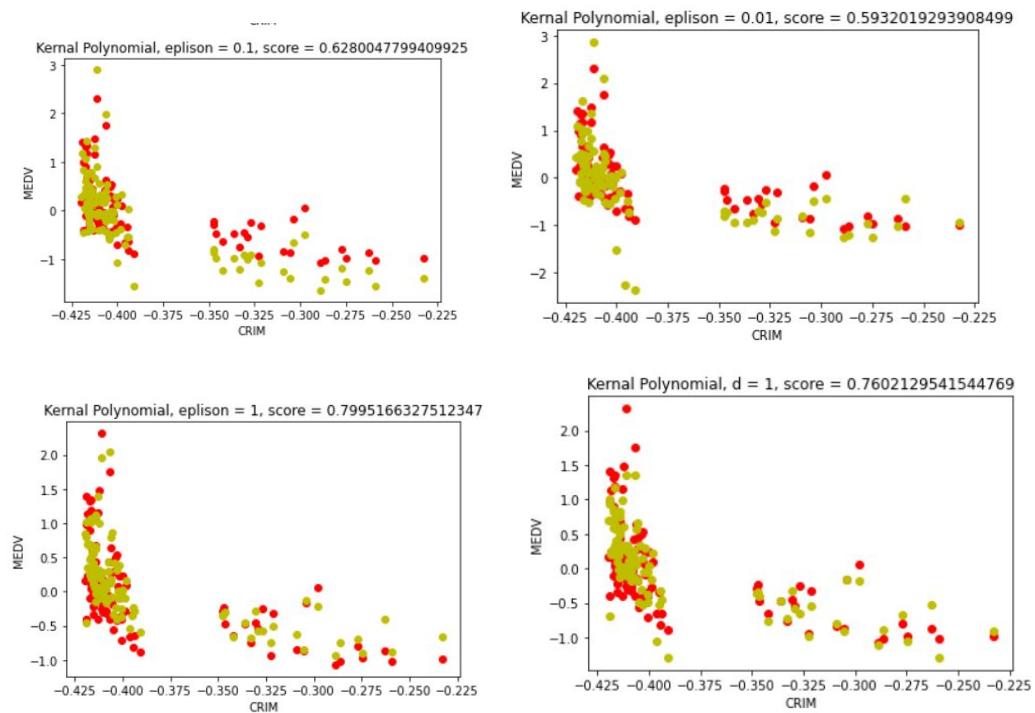
- Plot for **Polynomial Kernal** for  $C = 100$ , degree = 3, epsilon = 0.1, coeff0 = 1



- For different C, keeping other parameters same



- Plot for different epsilon





- Plot for different degrees of freedom(d)

