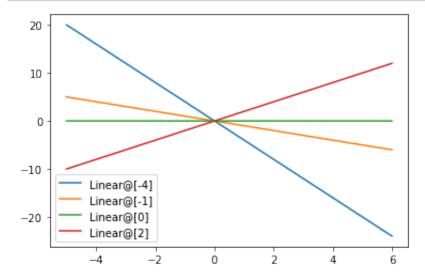
```
In [1]: import numpy as np
   import matplotlib.pyplot as plt
   import sklearn
   import scipy.spatial
   import functools
%matplotlib inline
```

```
In [2]:
        ### Kernel function generators
        def linear kernel(X1, X2):
             Computes the linear kernel between two sets of vectors.
             Args:
                 X1 - an \ n1xd \ matrix \ with \ vectors \ x1 \ 1,...,x1 \ n1 \ in \ the \ rows
                 X2 - an n2xd matrix with vectors x2_1,...,x2_n2 in the rows
             Returns:
                 matrix of size n1xn2, with x1_i^T x2_j in position i,j
             return np.dot(X1,np.transpose(X2))
        def RBF kernel(X1,X2,sigma):
             Computes the RBF kernel between two sets of vectors
             Args:
                 X1 - an \ n1xd \ matrix \ with \ vectors \ x1_1,...,x1_n1 \ in \ the \ rows
                 X2 - an n2xd matrix with vectors x2_1,...,x2_n2 in the rows
                 sigma - the bandwidth (i.e. standard deviation) for the RBF/Ga
             Returns:
                 matrix of size n1xn2, with exp(-||x1 i-x2 i||^2/(2 sigma^2)) i
             .....
             #TODO
             mat = scipy.spatial.distance.cdist(X1,X2, metric='sqeuclidean')
             return np.exp(-mat/(2*sigma**2))
        def polynomial kernel(X1, X2, offset, degree):
             Computes the inhomogeneous polynomial kernel between two sets of v
             Args:
                 X1 - an \ n1xd \ matrix \ with \ vectors \ x1_1,...,x1_n1 \ in \ the \ rows
                 X2 - an n2xd matrix with vectors x2_1,...,x2_n2 in the rows
                 offset, degree - two parameters for the kernel
             Returns:
                 matrix of size n1xn2, with (offset + <x1_i,x2_j>)^degree in po
             return (offset + linear_kernel(X1,X2))**degree
```

```
In [3]: # PLot kernel machine functions
plot_step = .01
xpts = np.arange(-5.0, 6, plot_step).reshape(-1,1)
prototypes = np.array([-4,-1,0,2]).reshape(-1,1)

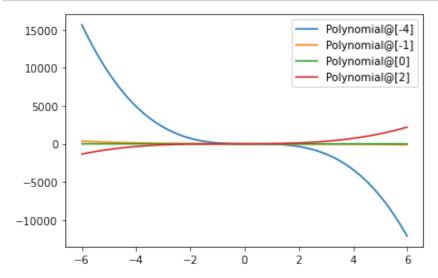
# Linear kernel
y = linear_kernel(prototypes, xpts)
for i in range(len(prototypes)):
    label = "Linear@"+str(prototypes[i,:])
    plt.plot(xpts, y[i,:], label=label)
plt.legend(loc = 'best')
plt.show()
```



## Part a)

```
In [4]: # PLot kernel machine functions
plot_step = .01
xpts = np.arange(-6.0, 6, plot_step).reshape(-1,1)
prototypes = np.array([-4,-1,0,2]).reshape(-1,1)

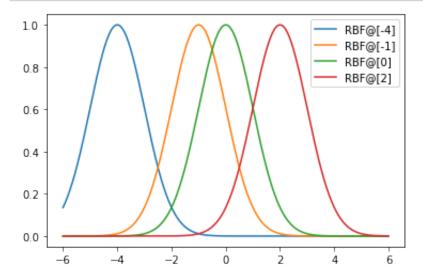
# Linear kernel
y = polynomial_kernel(prototypes, xpts,1,3)
for i in range(len(prototypes)):
    label = "Polynomial@"+str(prototypes[i,:])
    plt.plot(xpts, y[i,:], label=label)
plt.legend(loc = 'best')
plt.show()
```



## Part b

```
In [5]: # PLot kernel machine functions
plot_step = .01
xpts = np.arange(-6.0, 6, plot_step).reshape(-1,1)
prototypes = np.array([-4,-1,0,2]).reshape(-1,1)

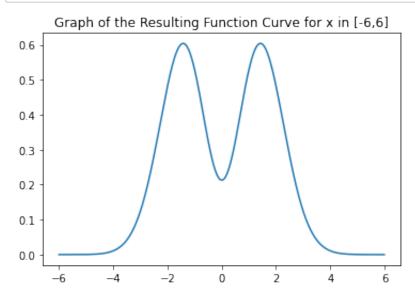
# Linear kernel
y = RBF_kernel(prototypes, xpts, 1)
for i in range(len(prototypes)):
    label = "RBF@"+str(prototypes[i,:])
    plt.plot(xpts, y[i,:], label=label)
plt.legend(loc = 'best')
plt.show()
```



```
In [6]: | class Kernel_Machine(object):
                __init__(self, kernel, training_points, weights):
                Aras:
                     kernel(X1,X2) - a function return the cross-kernel matrix
                    training points - an nxd matrix with rows x 1,..., x n
                    weights - a vector of length n with entries alpha_1,...,al
                .....
                self.kernel = kernel
                self.training_points = training_points
                self.weights = weights
            def predict(self, X):
                Evaluates the kernel machine on the points given by the rows of
                Args:
                    X – an nxd matrix with inputs x_1, \dots, x_n in the rows
                Returns:
                    Vector of kernel machine evaluations on the n points in X.
                         Sum \{i=1\}^R alpha i k(x j, mu i)
                .....
                kernel_matrix = self.kernel(X, self.training_points)
                 return (kernel matrix @ self.weights)
In [7]: k = functools.partial(RBF kernel, sigma=1)
        train_points = np.array([-1,0,1]).reshape(-1,1)
        weights = np.array([1,-1,1]).reshape(-1,1)
        X = np.array([-4,-1,0,1]).reshape(-1,1)
        test_kernel_object = Kernel_Machine(k, train_points, weights)
        test_kernel_object.predict(X)
Out[7]: array([[0.01077726],
                [0.52880462],
                [0.21306132],
                [0.5288046211)
```

```
In [8]: # PLot kernel machine functions
plot_step = .01
xpts = np.arange(-6.0, 6, plot_step).reshape(-1,1)
# Linear kernel
y = test_kernel_object.predict(xpts).reshape(-1,1)

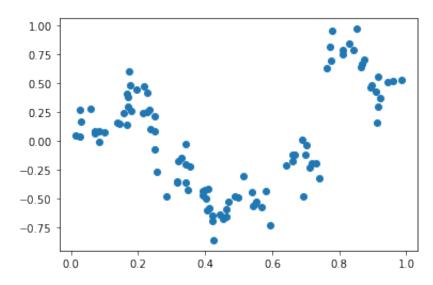
plt.plot(xpts, y)
plt.title('Graph of the Resulting Function Curve for x in [-6,6]')
plt.show()
```



Load train & test data; Convert to column vectors so it generalizes well to data in higher dimensions.

```
In [9]: data_train,data_test = np.loadtxt("krr-train.txt"),np.loadtxt("krr-tes
x_train, y_train = data_train[:,0].reshape(-1,1),data_train[:,1].resha
x_test, y_test = data_test[:,0].reshape(-1,1),data_test[:,1].reshape(-
plt.scatter(x_train,y_train)
```

Out[9]: <matplotlib.collections.PathCollection at 0x7fc8315edf10>

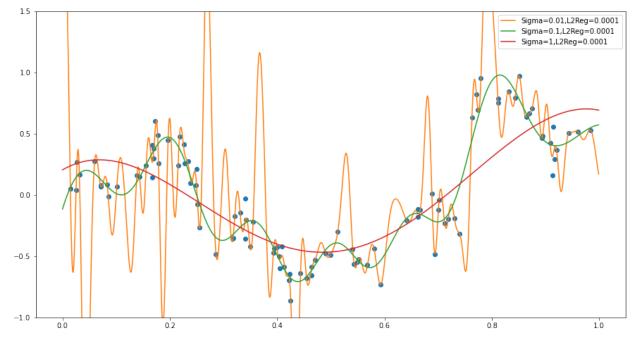


Looks like there is not a linear relationship, but rather a potential polynomial, or sinusoidal relationship, alternatively it could be a uniform + noise distribution.

## **Question 26**

```
In [10]: def train_kernel_ridge_regression(X, y, kernel, l2reg):
    kernel_matrix = kernel(X,X)
    matrix = ((np.identity(X.shape[0])*l2reg)+kernel_matrix)
    alpha = np.linalg.inv(matrix)@y
    return Kernel_Machine(kernel, X, alpha)
```

```
In [11]: plt.figure(figsize=(15,8))
    plot_step = .001
    xpts = np.arange(0 , 1, plot_step).reshape(-1,1)
    plt.plot(x_train,y_train,'o')
    l2reg = 0.0001
    for sigma in [.01,.1,1]:
        k = functools.partial(RBF_kernel, sigma=sigma)
        f = train_kernel_ridge_regression(x_train, y_train, k, l2reg=l2reg
        label = "Sigma="+str(sigma)+",L2Reg="+str(l2reg)
        plt.plot(xpts, f.predict(xpts), label=label)
    plt.legend(loc = 'best')
    plt.ylim(-1,1.5)
    plt.show()
```

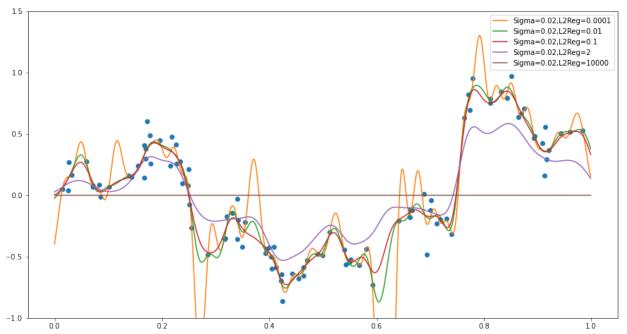


The smaller values of sigma,  $\sigma \in 0.01$  for instance, the curve overfits the training data, and does not generalize well (least bias). With a lower sigma, say  $\sigma \in 1$  the curve is much smoother with lower amplitude, and still does not fit the training data well (high bias). With  $\sigma = .1$  we have the best fit of the curve.

## **Problem 28**

```
In [12]: plt.figure(figsize=(15,8))
    plot_step = .001
    xpts = np.arange(0 , 1, plot_step).reshape(-1,1)
    plt.plot(x_train,y_train,'o')
    sigma= .02
    for l2reg in [.0001,.01,.1,2,10000]:
        k = functools.partial(RBF_kernel, sigma=sigma)
        f = train_kernel_ridge_regression(x_train, y_train, k, l2reg=l2reg label = "Sigma="+str(sigma)+",L2Reg="+str(l2reg)
        plt.plot(xpts, f.predict(xpts), label=label)

plt.legend(loc = 'best')
    plt.ylim(-1,1.5)
    plt.show()
```



As  $\lambda \to \infty$  the curve loses its amplitude, becomes less volatile, and becomes a line. Specifically, its the line that fits all points with least euclidean distance loss, which in this case approaches 0.

```
In [13]: from sklearn.base import BaseEstimator, RegressorMixin, ClassifierMixi
         class KernelRidgeRegression(BaseEstimator, RegressorMixin):
             """sklearn wrapper for our kernel ridge regression"""
             def init (self, kernel="RBF", sigma=1, degree=2, offset=1, l2re
                 self.kernel = kernel
                 self.sigma = sigma
                 self.degree = degree
                 self.offset = offset
                 self.l2reg = l2reg
             def fit(self, X, y=None):
                 This should fit classifier. All the "work" should be done here
                 if (self.kernel == "linear"):
                     self.k = linear kernel
                 elif (self.kernel == "RBF"):
                     self.k = functools.partial(RBF_kernel, sigma=self.sigma)
                 elif (self.kernel == "polynomial"):
                     self.k = functools.partial(polynomial_kernel,
                                                 offset=self.offset, degree=self
                 else:
                     raise ValueError('Unrecognized kernel type requested.')
                 self.kernel machine = train kernel ridge regression(X, y, sel
                 return self
             def predict(self, X, y=None):
                     getattr(self, "kernel_machine_")
                 except AttributeError:
                     raise RuntimeError("You must train classifer before predic
                 return(self.kernel_machine_.predict(X))
             def score(self, X, y=None):
                 # get the average square error
                 return(((self.predict(X)-y)**2).mean())
```

```
In [14]: from sklearn.model_selection import GridSearchCV, PredefinedSplit
         from sklearn.model selection import ParameterGrid
         from sklearn.metrics import mean squared error, make scorer
         import pandas as pd
         test fold = [-1]*len(x train) + [0]*len(x test)
                                                            #0 corresponds to te
         predefined split = PredefinedSplit(test fold=test fold)
In [15]: param_grid = [{'kernel': ['RBF'], 'sigma': [0.05, 0.055, 0.06],
                         'l2reg': [0.271, 0.27, 0.269]},
                       {'kernel':['polynomial'],'offset':[1.75, 1.8, 1.85],
                         'degree': [5,6,7], 'l2reg': [0.033, 0.034, 0.035]},
                       {'kernel':['linear'],'l2reg': [3.2, 4, 4.5]}]
         kernel_ridge_regression_estimator = KernelRidgeRegression()
         grid = GridSearchCV(kernel_ridge_regression_estimator,
                              param_grid,
                              cv = predefined_split,
                              scoring = make scorer(mean squared error, greater i
                              return train score=True
         grid.fit(np.vstack((x_train,x_test)),np.vstack((y_train,y_test)))
Out[15]: GridSearchCV(cv=PredefinedSplit(test_fold=array([-1, -1, ...,
                                                                              0]
         )),
                       estimator=KernelRidgeRegression(),
                       param_grid=[{'kernel': ['RBF'], 'l2reg': [0.271, 0.27, 0
         .269],
                                    'sigma': [0.05, 0.055, 0.06]},
                                   {'degree': [5, 6, 7], 'kernel': ['polynomial
         '],
                                    'l2reg': [0.033, 0.034, 0.035],
                                    'offset': [1.75, 1.8, 1.85]},
                                   {'kernel': ['linear'], 'l2reg': [3.2, 4, 4.5
         ]}],
                       return train score=True,
                       scoring=make_scorer(mean_squared_error, greater_is_bette
         r=False))
```

#### Out[16]:

	param_degree	param_kernel	param_l2reg	param_offset	param_sigma	mean_test_score ı
1	-	RBF	0.271	-	0.055	0.013821
4	-	RBF	0.270	-	0.055	0.013821
7	-	RBF	0.269	-	0.055	0.013821
8	-	RBF	0.269	-	0.06	0.013979
5	-	RBF	0.270	-	0.06	0.013982
12	5	polynomial	0.034	1.75	-	0.046631
15	5	polynomial	0.035	1.75	-	0.046974
37	-	linear	4.000	-	-	0.164510
38	-	linear	4.500	-	-	0.164511
36	-	linear	3.200	-	-	0.164511

39 rows × 7 columns

```
In [17]: rbf = df_toshow[df_toshow['param_kernel']=='RBF']
    rbf_min = rbf['mean_test_score'].min()
    rbf_min_row = rbf[rbf['mean_test_score']==rbf_min]
    print('Best parameters for RBF kernel:')
    rbf_min_row
```

Best parameters for RBF kernel:

#### Out[17]:

	param_degree	param_kernel	param_l2reg	param_offset	param_sigma	mean_test_score	m
1	-	RBF	0.271	-	0.055	0.013821	

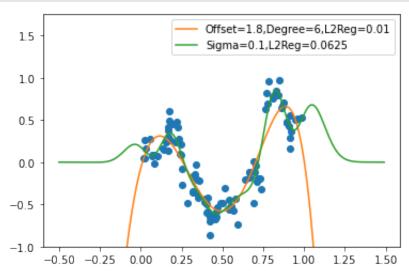
```
In [18]: polynomial = df_toshow[df_toshow['param_kernel']=='polynomial']
    poly_min = polynomial['mean_test_score'].min()
    poly_min_row = polynomial[polynomial['mean_test_score']==poly_min]
    print('Best parameters for polynomial kernel:')
    poly_min_row
```

Best parameters for polynomial kernel:

#### Out[18]:

	param_degree	param_kernel	param_l2reg	param_offset	param_sigma	mean_test_score	ı
22	6	polynomial	0.034	1.8	-	0.032405	

```
In [19]: ## Plot the best polynomial and RBF fits you found
         plot step = .01
         xpts = np.arange(-.5, 1.5, plot_step).reshape(-1,1)
         plt.plot(x_train,y_train,'o')
         #Plot best polynomial fit
         offset= 1.8
         degree = 6
         l2req = .01
         k = functools.partial(polynomial_kernel, offset=offset, degree=degree)
         f = train_kernel_ridge_regression(x_train, y_train, k, l2reg=l2reg)
         label = "Offset="+str(offset)+", Degree="+str(degree)+", L2Reg="+str(l2r
         plt.plot(xpts, f.predict(xpts), label=label)
         #Plot best RBF fit
         sigma = .1
         l2reg= .0625
         k = functools.partial(RBF_kernel, sigma=sigma)
         f = train_kernel_ridge_regression(x_train, y_train, k, l2reg=l2reg)
         label = "Sigma="+str(sigma)+",L2Reg="+str(l2reg)
         plt.plot(xpts, f.predict(xpts), label=label)
         plt.legend(loc = 'best')
         plt.ylim(-1,1.75)
         plt.show()
```



The best hyperparameters for the polynomial kernel were

 $Offset=1.8,\ Degree=6,\ l2reg=.034$  and for the RBF kernel the best parameters were  $\sigma=0.055,\ l2reg=0.271$ . Both curves seem to fit the graph reasonably well, though the RBF performed better with a lower min mean\_test\_score (

RBF = 0.013821, Polynomial = 0.032405). This could be because the best fitting polynomial curve could not capture the intricacies of the data without increasing its degree substantially, at the risk of over fitting on the train set, while the RBF kernel with its smoother curves could actually capture structure within the data that would generalize well.

Bayes function is defined in the following way:

$$E(y|x) = E(f(x) + \epsilon|x) \rightarrow E(f(x)|x) + E(\epsilon|x) = E(f(x))$$

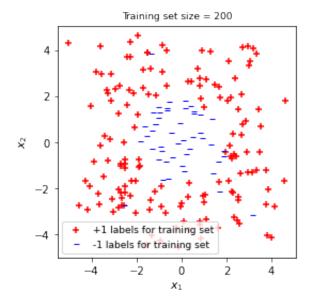
We can now calculate the risk of the bayes decision function:

$$E((f(x) - y)^2) = E((f(x) - f(x) + epsilon)^2) = E(\epsilon^2) = .1^2$$

As

$$Var(\epsilon) = E(\epsilon^2) - E^2(\epsilon) = E(\epsilon^2) - 0 = E(\epsilon^2)$$

```
In [20]: # Load and plot the SVM data
         #load the training and test sets
         data train,data test = np.loadtxt("svm-train.txt"),np.loadtxt("svm-tes
         x_train, y_train = data_train[:,0:2], data_train[:,2].reshape(-1,1)
         x_test, y_test = data_test[:,0:2], data_test[:,2].reshape(-1,1)
         #determine predictions for the training set
         yplus = np.ma.masked_where(y_train[:,0]<=0, y_train[:,0])</pre>
         xplus = x_train[~np.array(yplus.mask)]
         yminus = np.ma.masked_where(y_train[:,0]>0, y_train[:,0])
         xminus = x train[\simnp.array(yminus.mask)]
         #plot the predictions for the training set
         figsize = plt.figaspect(1)
         f, (ax) = plt.subplots(1, 1, figsize=figsize)
         pluses = ax.scatter (xplus[:,0], xplus[:,1], marker='+', c='r',
                               label = '+1 labels for training set')
         minuses = ax.scatter (xminus[:,0], xminus[:,1], marker=r'$-$',
                               c='b', label = '-1 labels for training set')
         ax.set_ylabel(r"$x_2$", fontsize=11)
         ax.set_xlabel(r"$x_1$", fontsize=11)
         ax.set_title('Training set size = %s'% len(data_train), fontsize=9)
         ax.axis('tight')
         ax.legend(handles=[pluses, minuses], fontsize=9)
         plt.show()
```



The data appears that it could be seperated by a quadratic boundary, (imagine a circle in the middle classifying the points). Additionally, a RBF boundary could work as well, as there is a circular distribution and you could imagine a guassian kernel coming out (as in the height extends to the z dimension) of the origin of the graph and decreasing otherwise.

```
In [21]: class train_soft_svm():
             def __init__(self, x_train, y_train, y_test, k, lambda_reg, epochs
                  self.x train = x train
                  self.y_train = y_train
                  self.y_test = y_test
                  self.k = k
                  self.lambda_reg = lambda_reg
                  self.epochs = epochs
                  self.alpha = None
             def predict(self, values):
                  return self.k(values, self.x_train) @ self.alpha
             def fit(self):
                  #Initialize helper variables
                  alpha = np.zeros(self.x_train.shape[0])
                  epoch, t = 0, 2
                 #Iterate over the epochs
                 while epoch < self.epochs:</pre>
                      for i in range(len(self.x train)):
                          alpha = alpha * (1-(1/t))
                          y_hat = self.k(self.x_train[i].reshape(1,2),self.x_tra
                          value = self.y_train[i] * y_hat
                          #If we have a missclassification, subtract the second
                          if value < 1:</pre>
                              step = y_train[i]/(t*self.lambda_reg)
                              alpha[i]+=1*step
                          t += 1
                      #Increment epoch counter variable
                      epoch += 1
                  self.alpha = alpha
                  return True
             def classification_error(self,y_hats):
                  error = 0
                  for i in range(len(y_hats)):
                      if y_hats[i] >= 0 and self.y_test[i] != 1:
                          error += 1
                      elif y_hats[i] < 0 and self.y_test[i] != -1:</pre>
                          error += 1
                  return error / len(y_hats)
```

Took a comprehensive iterative approach, started ide then honed in by restriciting the range I was searching over to get the best hyper parameters

```
In [22]: sigmas = np.arange(-3,0,.1)
         lambda_regs = np.logspace(-5, -2, num=40)
         rbf test accuracy = []
         rbf_lambda_reg = []
         rbf_sigma_list = []
         for sigma in sigmas:
             for lambda_reg in lambda_regs:
                 k = functools.partial(RBF kernel, sigma=sigma)
                 f = train_soft_svm(x_train, y_train, y_test, k, lambda_reg,20)
                 f.fit()
                 y_bar = f.predict(x_test)
                 rbf_test_accuracy.append(f.classification_error(y_bar))
                 rbf_lambda_reg.append(lambda_reg)
                 rbf_sigma_list.append(sigma)
         index = rbf_test_accuracy.index(min(rbf_test_accuracy))
         print("Best min mean test error at", rbf_test_accuracy[index]*100,
               "% with sigma = ",rbf_sigma_list[index],
               "lambda =", rbf_lambda_reg[index])
```

```
In [23]: degrees = np.arange(1,5,1)
         offsets = np.arange(1,10,20)
         lambda_regs = np.logspace(-5, -3, num=40)
         poly offset list = []
         poly_lambda_list = []
         poly degree list = []
         polynomial_test_accuracy = []
         for degree in degrees:
             for offset in offsets:
                 for lambda_reg in lambda_regs:
                     k = functools.partial(polynomial_kernel, offset=offset, de
                     f = train_soft_svm(x_train, y_train, y_test, k, lambda_reg
                     f.fit()
                     y_bar = f.predict(x_test)
                     polynomial_test_accuracy.append(f.classification_error(y_b
                     poly_lambda_list.append(lambda_reg)
                     poly_offset_list.append(offset)
                     poly_degree_list.append(degree)
         index = polynomial_test_accuracy.index(min(polynomial_test_accuracy))
         print("Best min mean test error at",polynomial_test_accuracy[index]*10
               "% with offset = ",poly_offset_list[index],
               "degree = ", poly_degree_list[index],
               "lambda =", poly_lambda_list[index])
```

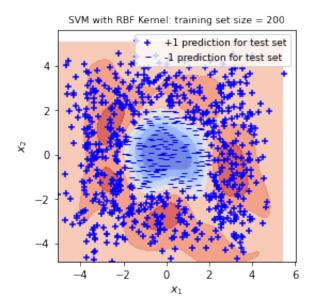
da = 0.0004375479375074184

## RBF Kernel, Optimal Fit, $\sigma = -.99999$ , $\lambda = .00412$

Best min mean test error at 5.625 % with offset = 1 degree = 2 lamb

```
xx, yy = np.meshgrid(np.arange(x1_min, x1_max, h),
                      np.arange(x2_min, x2_max, h))
Z = f.predict(np.c_[xx.ravel(), yy.ravel()])
Z = Z.reshape(xx.shape)
#determine the predictions for the test set
y_bar = f.predict(x_test)
vplus = np.ma.masked where(y bar<=0, y bar)</pre>
xplus = x_test[~np.array(yplus.mask)]
yminus = np.ma.masked_where(y_bar>0, y_bar)
xminus = x_test[~np.array(yminus.mask)]
print('Classification Error:',f.classification error(y bar)*100, '%')
#plot the learned boundary and the predictions for the test set
figsize = plt.figaspect(1)
f, (ax) = plt.subplots(1, 1, figsize=figsize)
decision =ax.contourf(xx, yy, Z, cmap=plt.cm.coolwarm, alpha=0.8)
pluses = ax.scatter (xplus[:,0], xplus[:,1], marker='+', c='b',
                      label = '+1 prediction for test set')
minuses = ax.scatter (xminus[:,0], xminus[:,1], marker=r'$-$',
                       c='b', label = '-1 prediction for test set')
ax.set_ylabel(r"$x_2$", fontsize=11)
ax.set_xlabel(r"$x_1$", fontsize=11)
ax.set_title('SVM with RBF Kernel: training set size = %s'% len(data_t
ax.axis('tight')
ax.legend(handles=[pluses, minuses], fontsize=9)
plt.show()
```

#### Classification Error: 3.375 %

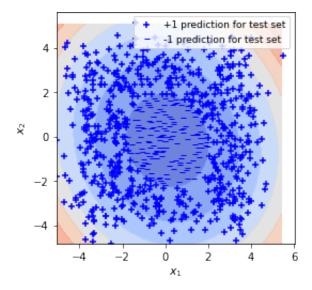


# Polynomial Kernel, Optimal Fit, Offset = 1, Degree = 2, $\lambda = 0.0004375479375074184$

```
In [25]: # Code to help plot the decision regions
         # (Note: This ode isn't necessarily entirely appropriate for
         the questions asked. So think about what you are doing.)
         k = functools.partial(polynomial_kernel, offset=1,degree=2)
         f = train_soft_svm(x_train, y_train, y_test, k, 0.0004375479375074184,
         f.fit()
         #determine the decision regions for the predictions
         x1_min = min(x_test[:,0])
         x1_max = max(x_test[:,0])
         x2_min = min(x_test[:,1])
         x2 max = max(x test[:,1])
         h=0.1
         xx, yy = np.meshgrid(np.arange(x1_min, x1_max, h),
                              np.arange(x2_min, x2_max, h))
         Z = f.predict(np.c_[xx.ravel(), yy.ravel()])
         Z = Z.reshape(xx.shape)
         #determine the predictions for the test set
         y bar = f.predict(x test)
         yplus = np.ma.masked where(y bar<=0, y bar)</pre>
         xplus = x test[~np.array(yplus.mask)]
         yminus = np.ma.masked_where(y_bar>0, y_bar)
         xminus = x_test[~np.array(yminus.mask)]
         print('Classification Error:',f.classification_error(y_bar)*100, '%')
         #plot the learned boundary and the predictions for the test set
         figsize = plt.figaspect(1)
         f, (ax) = plt.subplots(1, 1, figsize=figsize)
         decision =ax.contourf(xx, yy, Z, cmap=plt.cm.coolwarm, alpha=0.8)
         pluses = ax.scatter (xplus[:,0], xplus[:,1], marker='+', c='b',
                               label = '+1 prediction for test set')
         minuses = ax.scatter (xminus[:,0], xminus[:,1], marker=r'$-$',
                                c='b', label = '-1 prediction for test set')
         ax.set_ylabel(r"$x_2$", fontsize=11)
         ax.set_xlabel(r"$x_1$", fontsize=11)
         ax.set_title('SVM with RBF Kernel: training set size = %s'% len(data_t
         ax.axis('tight')
         ax.legend(handles=[pluses, minuses], fontsize=9)
         plt.show()
```

Classification Error: 5.625 %

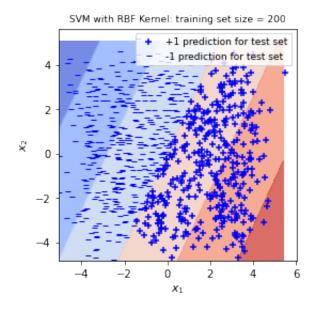
SVM with RBF Kernel: training set size = 200



### **Linear Kernel, Optimal Fit**

```
In [26]: # Code to help plot the decision regions
         # (Note: This ode isn't necessarily entirely appropriate for the quest
         sigma=1
         k = functools.partial(linear_kernel)
         f = train_soft_svm(x_train, y_train, y_test, k, 1e-05,50)
         f.fit()
         #determine the decision regions for the predictions
         x1 min = min(x test[:,0])
         x1_max= max(x_test[:,0])
         x2_min = min(x_test[:,1])
         x2_max = max(x_test[:,1])
         h=0.1
         xx, yy = np.meshgrid(np.arange(x1_min, x1_max, h),
                               np.arange(x2_min, x2_max, h))
         Z = f.predict(np.c_[xx.ravel(), yy.ravel()])
         Z = Z.reshape(xx.shape)
         #determine the predictions for the test set
         y_bar = f.predict(x_test)
         yplus = np.ma.masked_where(y_bar<=0, y_bar)</pre>
         xplus = x_test[~np.array(yplus.mask)]
         yminus = np.ma.masked where(y bar>0, y bar)
         xminus = x_test[~np.array(yminus.mask)]
         print('Classification Error:',f.classification_error(y_bar)*100, '%')
         #plot the learned boundary and the predictions for the test set
         figsize = plt.figaspect(1)
         f, (ax) = plt.subplots(1, 1, figsize=figsize)
```

#### Classification Error: 49.875 %



```
In [ ]:
```