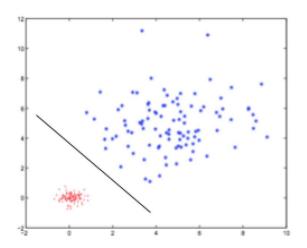
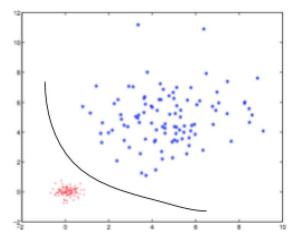
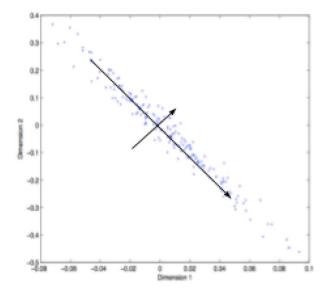
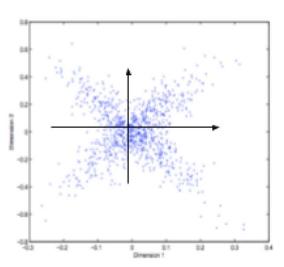
The left one is linear, the right one is quadratic boundary.





Question 2





Question 3

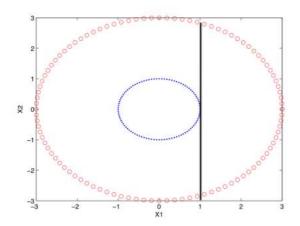
- a. Ture, given data D, the probability of Hypothesis is 1.
- b. False, given hypothesis H, the probability of data is not 1.
- c. False, the probability of data and probability is not 1.

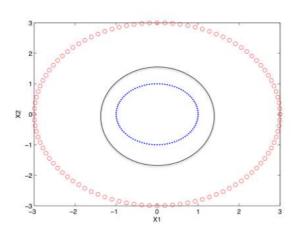
- a. The maximum number of leaf nodes is 2^{k-1} , the time complexity is O(k).
- b. The maximum number of leaf nodes is N, the maximum depth is N.

Question 5

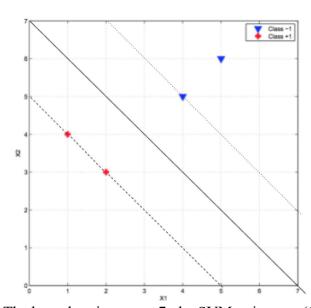
a. Correct is a, the first is $\exp(0) = 1$, the second is $\exp(-\infty) = 0$.

b.





c.



The boundary is y = -x+7, the SVM points are (1,4), (2,3) and (4,5).

- a. 100, it selects the variable that is significant in correlation analysis, thus the total 100 will be selected.
- b. Lower, because the training data will be more.
- c. Polynomial of degree 6, because it has more information that degree 4 and degree 5.

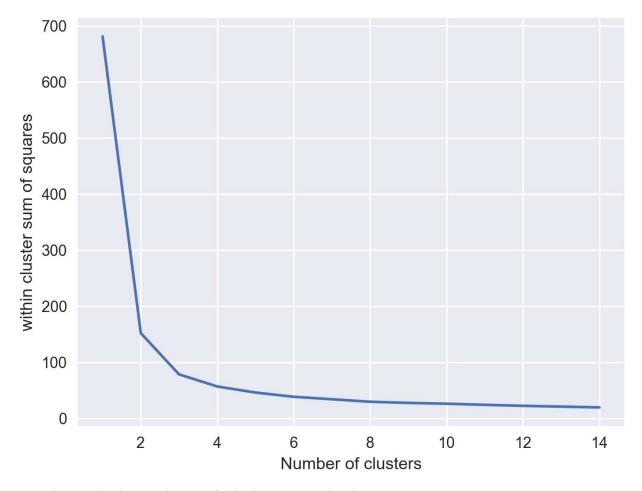
Question 7

No. It is not possible since the boundaries of decision trees are parallel to x or y axes.

Question 8

```
###Load libraries
import numpy as np
%matplotlib notebook
import matplotlib.pyplot as plt
# Though the following import is not directly being used, it is required
# for 3D projection to work
from mpl_toolkits.mplot3d import Axes3D
from sklearn.cluster import KMeans
from sklearn import datasets
import seaborn as sns; sns.set() # for plot styling
np.random.seed(5)
iris = datasets.load_iris()
X = iris.data
y = iris.target
```

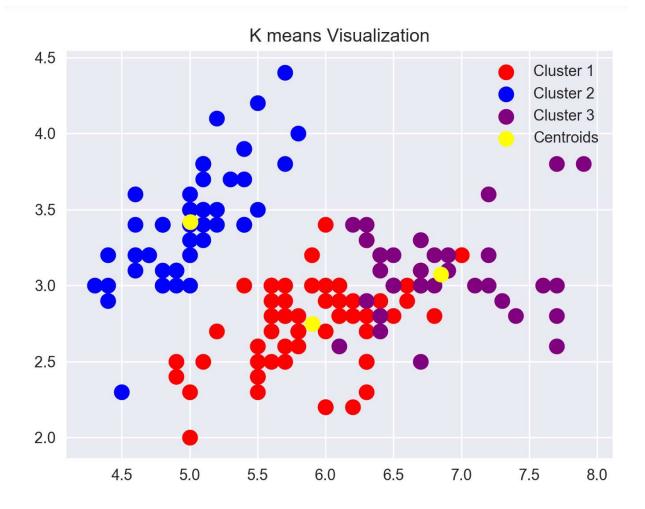
```
#Finding the optimum number of clusters for k-means classification
from sklearn.cluster import KMeans
ERR = []
for i in range(1, 15):
    kmeans = KMeans(n_clusters = i, init = 'k-means++', max_iter = 500, n_init = 10, random_state = 42)
    kmeans.fit(X)
    ERR.append(kmeans.inertia_)
#Plotting the results onto a line graph, allowing us to observe 'The elbow'
plt.plot(range(1, 15), ERR)
plt.xlabel('Number of clusters')
plt.ylabel('within cluster sum of squares')
plt.show()
```



From the graph above, cluster of 3 is the most optimal one.

```
#Applying kmeans to the dataset / Creating the kmeans classifier
kmeans = KMeans(n_clusters = 3, init = 'k-means++', max_iter = 300, n_init = 10, random_state = 0)
Y_kmeans = kmeans.fit_predict(X)

#Visualising the clusters
plt.scatter(X[Y_kmeans == 0, 0], X[Y_kmeans == 0, 1], s = 100, c = 'red', label = 'Cluster 1')
plt.scatter(X[Y_kmeans == 1, 0], X[Y_kmeans == 1, 1], s = 100, c = 'blue', label = 'Cluster 2')
plt.scatter(X[Y_kmeans == 2, 0], X[Y_kmeans == 2, 1], s = 100, c = 'purple', label = 'Cluster 3')
#Plotting the centroids of the clusters
plt.scatter(kmeans.cluster_centers_[:, 0], kmeans.cluster_centers_[:,1],s = 100, c = 'yellow', label = 'Centroids')
plt.title('K_means_Visualization')
plt.legend()
```



```
import numpy as np
from sklearn import datasets
from sklearn import svm
from sklearn.metrics import accuracy_score
from sklearn.model_selection import train_test_split
import matplotlib.pyplot as plt

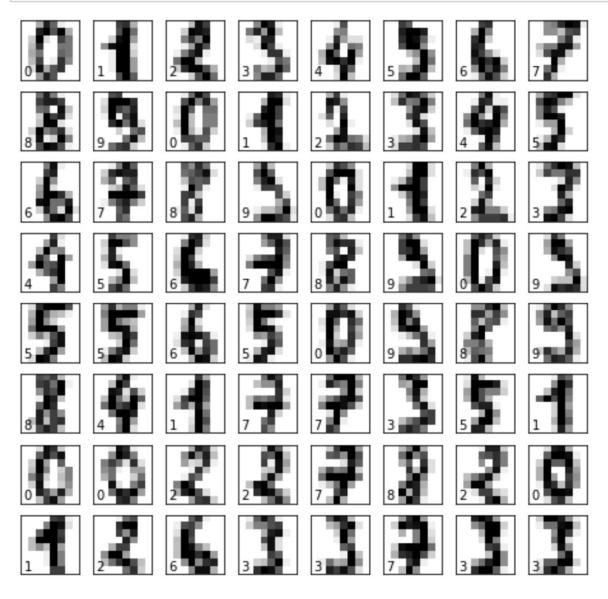
d = datasets.load_digits()
X = d.data
y = d.target
```

```
X.shape, y.shape
((1797, 64), (1797,))
```

```
###Split dataset into train/test
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.3,random_state=42)
```

```
##display the dataset
# set up the figure
fig = plt.figure(figsize=(8, 8))

for i in range(64):
    ax = fig.add_subplot(8, 8, i + 1, xticks=[], yticks=[])
    ax.imshow(d.images[i], cmap=plt.cm.binary,interpolation='nearest')
    ax.text(0, 7, str(y[i]))
```



```
import pandas as pd
def accuracy(ytest, ypred):
    pred=pd.DataFrame()
    pred['label']=ytest
    pred['Predicted']=ypred
    pred['Accuracy']=1*(pred.label==pred.Predicted)
    pred = pred.groupby('label').mean()
    return pred.drop(['Predicted'], axis=1)
 from sklearn.svm import SVC
 ac=[]
for c in [ 0.1, 1, 10, 100]:
    for q in [5e-7, 1e-6, 2e-6]:
        clf=SVC(C=c, gamma=g)
        clf.fit(Xtrain,ytrain)
        svm pred=clf.predict(Xtest)
        ac.append([c,q,accuracy(ytest, svm pred).mean()])
ac
[[0.1, 5e-07, Accuracy
                            0.1
  dtype: float64], [0.1, 1e-06, Accuracy
                                                0.1
  dtype: float64], [0.1, 2e-06, Accuracy
                                                0.1
  dtype: float64], [1, 5e-07, Accuracy
                                              0.1
  dtype: float64], [1, 1e-06, Accuracy
                                              0.1
  dtype: float64], [1, 2e-06, Accuracy
                                              0.1
  dtype: float64], [10, 5e-07, Accuracy
                                               0.718732
  dtype: float64], [10, 1e-06, Accuracy
                                               0.893875
  dtype: float64], [10, 2e-06, Accuracy
                                               0.933628
  dtype: float64], [100, 5e-07, Accuracy
                                                0.953352
  dtype: float64], [100, 1e-06, Accuracy
                                                0.966069
  dtype: float64], [100, 2e-06, Accuracy
                                                0.976435
  dtype: float64]]
# Now try all samples
svm=SVC(C=100, gamma=2e-6)
svm.fit(Xtrain,ytrain)
```

From the above we can find that, C = 100, gamma = 2e-6 can achieve the highest accuracy.

svm pred=svm.predict(Xtest)

```
svm_accuracy=accuracy(ytest, svm_pred)
print (svm_accuracy)
print (svm_accuracy.mean())
```

```
Accuracy
label
0
        1.000000
1
        1.000000
2
        1.000000
3
       0.962963
4
        1.000000
5
       0.969697
       0.981132
6
7
       0.981818
        0.953488
8
9
       0.915254
Accuracy
             0.976435
dtype: float64
```

```
from sklearn.metrics import confusion_matrix
import seaborn as sns; sns.set()
mat = confusion_matrix(ytest,svm_pred)
sns.heatmap(mat.T, square=True, annot=True, fmt='d', cbar=False)
plt.xlabel('true label')
plt.ylabel('predicted label');
```

