iris\_data = load\_iris() In [0]: iris\_data.keys() Out[0]: dict\_keys(['data', 'target', 'target\_names', 'DESCR', 'feature\_names', 'filename']) In [0]: # print(iris\_data.DESCR) # for details of dataset In [0]: | iris\_data.feature\_names Out[0]: ['sepal length (cm)', 'sepal width (cm)', 'petal length (cm)' 'petal width (cm)'] In [0]: | from sklearn.model\_selection import train\_test\_split X\_train, X\_test, y\_train, y\_test = train\_test\_split(iris\_data['data'], iris\_data['target'], test\_siz e=0.25, random\_state=0) In [0]: y\_train Out[0]: array([1, 1, 2, 0, 2, 0, 0, 1, 2, 2, 2, 1, 2, 1, 1, 1, 0, 2, 1, 1, 1, 1, 2, 0, 0, 2, 1, 0, 0, 1, Θ, 2, 1, 0, 2, 2, 2, 0, 0, 2, 2, 0, 2, 0, 2, 0, 0, 2, 0, 0, 0, 1, 2, 2, 0, 0, 0, 1, 1, 0, 0, 1, 0, 2, 1, 2, 1, 0, 2, 0, 2, 0, 0, 2, 0, 2, 1, 1, 1, 2, 2, 1, 1, 0, 1, 2, 2, 0, 1, 1, 1, 1, 0, 0, 0, 2, 1, 2, 0]) In [0]: import pandas as pd from pandas import plotting as pdplt iris\_df = pd.DataFrame(X\_train,columns=iris\_data['feature\_names']) vis = pdplt.scatter\_matrix(iris\_df, c=y\_train, diagonal='hist', figsize=[10,10], marker='0', alpha= 0.5, s=40, hist\_kwds={'bins':20}) sepal length (cm) 4.0 sepal width (cm) 3.5 3.0 2.5 2.0 petal length (cm) 2.5 petal width 1.0 sepal length (cm) petal length (cm) In [0]: from sklearn.neighbors import KNeighborsClassifier knn = KNeighborsClassifier(n\_neighbors=1) knn.fit(X\_train,y\_train) perc\_accu = knn.score(X\_test,y\_test) \* 100 print(f"perc accu : {perc\_accu}") perc accu : 97.36842105263158 So using all 4 features we get 97% accuracy which is great but now let's reduce the feature space from 4D to 2D and let's see how well we do as the above scatter matrix clearly shows that if we just take petal's length and width then the model is very easily separable iris\_data = load\_iris() In [0]: iris\_data.data = iris\_data.data[:, 2:] iris\_data.feature\_names = iris\_data.feature\_names[2:] In [0]: iris\_data.feature\_names Out[0]: ['petal length (cm)', 'petal width (cm)'] In [0]: X\_train, X\_test, y\_train, y\_test = train\_test\_split(iris\_data.data,iris\_data.target,test\_size=0.25,r andom\_state=0) iris\_df = pd.DataFrame(X\_train, columns=iris\_data['feature\_names']) In [0]: vis = pdplt.scatter\_matrix(iris\_df, c=y\_train, diagonal='hist', figsize=[7,7], marker='0', alpha=0.5 s=40, hist\_kwds={'bins':20}) petal length (cm) 25 20 petal width (cm) 10 petal length (cm) petal width (cm) In [0]: from sklearn.neighbors import KNeighborsClassifier knn = KNeighborsClassifier(n\_neighbors=1) knn.fit(X\_train,y\_train) perc\_accu = knn.score(X\_test,y\_test) \* 100 print(f"perc accu : {perc\_accu}") perc accu : 97.36842105263158 So using just 2D feature space we get max accuracy of 97% which is exact same as 4D feature space gives us, so this tells us that the other two features are just redundant and adding just nothing to our model iris\_data = load\_iris() In [0]: In [0]: iris\_data.data = iris\_data.data[:, 1:3] iris\_data.feature\_names = iris\_data.feature\_names[1:3] iris\_data.feature\_names In [0]: Out[0]: ['sepal width (cm)', 'petal length (cm)'] In [0]: X\_train, X\_test, y\_train, y\_test = train\_test\_split(iris\_data.data,iris\_data.target,test\_size=0.25,r andom\_state=0) In [0]: iris\_df = pd.DataFrame(X\_train,columns=iris\_data['feature\_names']) vis = pdplt.scatter\_matrix(iris\_df, c=y\_train, diagonal='hist', figsize=[7,7], marker='0', alpha=0.5 s=40, hist\_kwds={'bins':20}) 4.0 sepal width (cm) 3.0 20 petal length (cm) 40 20 petal length (cm) sepal width (cm) In [0]: **from sklearn.neighbors import** KNeighborsClassifier knn = KNeighborsClassifier(n\_neighbors=1) knn.fit(X\_train,y\_train) perc\_accu = knn.score(X\_test,y\_test) \* 100 print(f"perc accu : {perc\_accu}") perc accu : 89.47368421052632 similarly we can also take petal length and sepal length In [0]: | iris\_data=load\_iris() iris\_data.data = iris\_data.data[:, [0,2]] In [0]: iris\_data.feature\_names = [iris\_data.feature\_names[i] for i in range(4) if i%2 == 0] iris\_data.feature\_names In [0]: Out[0]: ['sepal length (cm)', 'petal length (cm)'] In [0]: X\_train, X\_test, y\_train, y\_test = train\_test\_split(iris\_data.data,iris\_data.target,test\_size=0.25,r andom\_state=0) In [0]: | iris\_df = pd.DataFrame(X\_train, columns=iris\_data['feature\_names']) vis = pdplt.scatter\_matrix(iris\_df, c=y\_train, diagonal='hist', figsize=[7,7], marker='0', alpha=0.5 , s=40, hist\_kwds={'bins':20}) 7.5 7.0 sepal length (cm) 6.5 6.0 5.5 5.0 petal length (cm) sepal length (cm) petal length (cm) In [0]: from sklearn.neighbors import KNeighborsClassifier knn = KNeighborsClassifier(n\_neighbors=1) knn.fit(X\_train,y\_train) perc\_accu = knn.score(X\_test,y\_test) \* 100 print(f"perc accu : {perc\_accu}") perc accu : 94.73684210526315 similarly we can also take sepal's length and width In [0]: iris\_data = load\_iris() In [0]: | iris\_data.data = iris\_data.data[:, [0,1]] iris\_data.feature\_names = [iris\_data.feature\_names[i] for i in range(4) if i <= 1]</pre> iris\_data.feature\_names In [0]: Out[0]: ['sepal length (cm)', 'sepal width (cm)'] In [0]: X\_train, X\_test, y\_train, y\_test = train\_test\_split(iris\_data.data,iris\_data.target,test\_size=0.25,r andom\_state=0) In [0]: | iris\_df = pd.DataFrame(X\_train, columns=iris\_data['feature\_names']) vis = pdplt.scatter\_matrix(iris\_df, c=y\_train, diagonal='hist', figsize=[7,7], marker='0', alpha=0.5 , s=40, hist\_kwds={'bins':20}) 7.5 7.0 length (cm) 6.5 6.0 sepall 5.5 5.0 45 4.0 sepal width (cm) 3.5 3.0 25 20 40 sepal length (cm) sepal width (cm) In [0]: from sklearn.neighbors import KNeighborsClassifier knn = KNeighborsClassifier(n\_neighbors=1) knn.fit(X\_train,y\_train) perc\_accu = knn.score(X\_test,y\_test) \* 100 print(f"perc accu : {perc\_accu}") perc accu : 71.05263157894737 Now as from the top most scatter matrix it is clearly evident that sepal's length and sepal's width are the worst pair to classify the dataset so when we use that as the feature than we get worse results iris\_data = load\_iris() In [0]: In [0]: | iris\_data.data = iris\_data.data[:, :2] iris\_data.feature\_names = iris\_data.feature\_names[:2] iris\_data.feature\_names In [0]: Out[0]: ['sepal length (cm)', 'sepal width (cm)'] In [0]: X\_train, X\_test, y\_train, y\_test = train\_test\_split(iris\_data.data,iris\_data.target,test\_size=0.25,r andom\_state=0) In [0]: | iris\_df = pd.DataFrame(X\_train, columns=iris\_data['feature\_names']) vis = pdplt.scatter\_matrix(iris\_df, c=y\_train, diagonal='hist', figsize=[7,7], marker='0', alpha=0.5 , s=40, hist\_kwds={'bins':20}) 7.5 7.0 length (cm) 6.5 6.0 sepal 5.0 45 40 sepal width (cm) 3.5 25 20 40 sepal length (cm) sepal width (cm) In [0]: **from sklearn.neighbors import** KNeighborsClassifier knn = KNeighborsClassifier(n\_neighbors=1) knn.fit(X\_train,y\_train) perc\_accu = knn.score(X\_test,y\_test) \* 100 print(f"perc accu : {perc\_accu}") perc accu : 71.05263157894737 Until now we are dealing with 2D feature space and are succefully get great results but now try to reduce the feature space further i.e., further from 2D to 1D In [0]: iris\_data=load\_iris() In [0]: iris\_data.target\_names Out[0]: array(['setosa', 'versicolor', 'virginica'], dtype='<U10')</pre> In [0]: # %matplotlib qt %matplotlib inline import matplotlib.pyplot as plt import numpy as np fig, axes =plt.subplots(1,4, figsize=(14, 3)) # 1 column containing 4 figures of all 4 features setosa=iris\_data.data[iris\_data.target==0] versicolor=iris\_data.data[iris\_data.target==1] virginica=iris\_data.data[iris\_data.target==2] ax=axes.ravel()# flat axes with numpy ravel for i in range(4): \_,bins=np.histogram(iris\_data.data[:,i],bins=40) ax[i].hist(setosa[:,i],bins=bins,color='r',alpha=0.5) ax[i].hist(versicolor[:,i], bins=bins, color='g', alpha=0.5) ax[i].hist(virginica[:,i],bins=bins,color='b',alpha=0.5) ax[i].set\_title(iris\_data.feature\_names[i],fontsize=9) ax[i].axes.get\_xaxis().set\_visible(False) # the x-axis co-ordinates are not so useful, as we jus t want to look how well separated the histograms are ax[i].set\_yticks(()) ax[0].legend(['setosa', 'versicolor', 'virginica'], loc='best', fontsize=8) plt.tight\_layout() plt.show() Now from the above histograms it is very clear that petal's width and length are very good features to classify and sepal's both dimensions are very bad for classfying the flowers Let's first take petal's length In [0]: | iris\_data.data = iris\_data.data[:, 2:3] iris\_data.feature\_names = iris\_data.feature\_names[2:3] In [0]: | iris\_data.feature\_names Out[0]: ['petal length (cm)'] X\_train, X\_test, y\_train, y\_test = train\_test\_split(iris\_data.data,iris\_data.target,test\_size=0.25,r In [0]: andom\_state=0) In [0]: X\_train.shape Out[0]: (112, 1) In [0]: from sklearn.neighbors import KNeighborsClassifier knn = KNeighborsClassifier(n\_neighbors=1) knn.fit(X\_train,y\_train) perc\_accu = knn.score(X\_test,y\_test) \* 100 print(f"perc accu : {perc\_accu}") perc accu : 89.47368421052632 we got 90% accuracy which is not bad as we are dealing with just 1 feature Let's first take petal's width In [0]: iris\_data=load\_iris() In [0]: | iris\_data.data = iris\_data.data[:, 3:4] iris\_data.feature\_names = iris\_data.feature\_names[-1] In [0]: X\_train, X\_test, y\_train, y\_test = train\_test\_split(iris\_data.data,iris\_data.target,test\_size=0.25,r andom\_state=0) In [0]: X\_train.shape Out[0]: (112, 1) In [0]: from sklearn.neighbors import KNeighborsClassifier knn = KNeighborsClassifier(n\_neighbors=1) knn.fit(X\_train,y\_train) perc\_accu = knn.score(X\_test,y\_test) \* 100 print(f"perc accu : {perc\_accu}") perc accu : 78.94736842105263 So in this case we got around 80% accuracy although it is indeed low and we will not classify our model using this feauture because it is 10% lower than previous one but if we still think intiutively this feature still doing some good job in classfying the model because if we make a classifier which randomly picks a class from three classes then the accuracy will be 33.33%, 1/3 using basic probability, so as compared to random selection this classifer is way more than double of that accuracy so it's not as bad as it seems to Let's first take sepal's length iris\_data=load\_iris() In [0]: In [0]: iris\_data.data = iris\_data.data[:, 0:1] iris\_data.feature\_names = iris\_data.feature\_names[0:1] In [0]: X\_train, X\_test, y\_train, y\_test = train\_test\_split(iris\_data.data,iris\_data.target,test\_size=0.25,r andom\_state=0) In [0]: X\_train.shape Out[0]: (112, 1) In [0]: **from sklearn.neighbors import** KNeighborsClassifier knn = KNeighborsClassifier(n\_neighbors=1) knn.fit(X\_train,y\_train) perc\_accu = knn.score(X\_test,y\_test) \* 100 print(f"perc accu : {perc\_accu}") perc accu: 47.368421052631575 Now this is extremly bad and nothing more than a random selection of 1 from three (whose accuracy is 33.33%) and this is what we expect as the histogram clearly shows that this feature is bad in classfying Let's first take sepal's width In [0]: iris\_data=load\_iris() In [0]: iris\_data.data = iris\_data.data[:, 1:2] iris\_data.feature\_names = iris\_data.feature\_names[1] iris\_data.feature\_names In [0]: Out[0]: 'sepal width (cm)' In [0]: X\_train, X\_test, y\_train, y\_test = train\_test\_split(iris\_data.data,iris\_data.target,test\_size=0.25,r andom\_state=0) In [0]: X\_train.shape Out[0]: (112, 1) In [0]: from sklearn.neighbors import KNeighborsClassifier knn = KNeighborsClassifier(n\_neighbors=1) knn.fit(X\_train,y\_train) perc\_accu = knn.score(X\_test,y\_test) \* 100 print(f"perc accu : {perc\_accu}") perc accu : 50.0 and in this case also we get expected results Until now we are just making our hands dirty on our own visualization using scatter matrix and histograms but now it's time for using some magic and the magic we our going to use is PCA and SVD LET'S FIRST LOOK AT PCA In [0]: iris\_data = load\_iris() In [0]: # Variance ratio of the four features ex\_variance=np.var(iris\_data.data,axis=0) ex\_variance\_ratio = ex\_variance/np.sum(ex\_variance) print( ex\_variance\_ratio ) [0.14994532 0.04154411 0.68145793 0.12705264] In [0]: X\_train, X\_test, y\_train, y\_test = train\_test\_split(iris\_data.data,iris\_data.target,test\_size=0.25,r andom\_state=0) In [0]: from sklearn.decomposition import PCA initial 4 features will be combined to make the best 3 principal components these are not features but are principal compnents generated from the original 4 features these principal components are some complex mixture of those 4 features pca = PCA(n\_components=3) pca.fit(X\_train)  $X_pca = pca.transform(X_train)$ X\_test = pca.transform(X\_test) print("shape of X\_pca", X\_pca.shape) shape of  $X_pca$  (112, 3) In [0]: pca.explained\_variance\_ratio\_ Out[0]: array([0.92743115, 0.05245721, 0.01557205]) the following matrix is the exact composition of features to form the PCs In [0]: pca.components\_ Out[0]: array([[ 0.37649644, -0.06637905, 0.85134571, 0.35924188], [ 0.6240207 , 0.75538031, -0.18479376, -0.07648543], [-0.60667794, 0.57674603, 0.08522779, 0.54040922]])df = pd.DataFrame(pca.components\_, columns=iris\_data.feature\_names) In [0]: df Out[0]: sepal length (cm) sepal width (cm) petal length (cm) petal width (cm) 0 0.376496 -0.066379 0.359242 0.851346 -0.076485 0.624021 0.755380 1 -0.184794 0.576746 2 -0.606678 0.085228 0.540409 The following plot shows that how the PC's are composed with the given features as shown in above table, this is a magical composition of all features in different proportion which gives us these PC's which suited the model best In [0]: these 3 principal components are some complex mixture of those 4 features heat-plot to see how the features mixed up to create the components. plt.matshow(pca.components\_, cmap='viridis') plt.yticks([0,1,2],['PC1','PC2','PC3'],fontsize=10) plt.colorbar() plt.xticks(range(len(iris\_data.feature\_names)), iris\_data.feature\_names, rotation=65, ha='left') plt.tight\_layout() plt.show() /usr/local/lib/python3.6/dist-packages/matplotlib/figure.py:2366: UserWarning: This figure includes A xes that are not compatible with tight\_layout, so results might be incorrect. warnings.warn("This figure includes Axes that are not compatible " 0.8 0.6 PC1 0.4 0.2 PC2 0.0 -0.2PC3 -0.4-0.6 In [0]: pca.singular\_values\_ Out[0]: array([22.23515693, 5.28812784, 2.88119225]) Below is a covariance heatmap, diagonals are the variance of that indexed feature more this variance is the more powerful is that individual feature for classifying data and the off-digonal are the relation/variance between the features i.e., how one feature vary with other feautre lower this covariance is better those features are as those feature are more likely to be orthogonal and both contributes individually In [0]: import seaborn as sns the corvariation heatmap of features df\_iris = pd.DataFrame(iris\_data.data,columns=iris\_data.feature\_names) s=sns.heatmap(df\_iris.cov(),cmap='coolwarm') s.set\_yticklabels(s.get\_yticklabels(),rotation=30,fontsize=7) s.set\_xticklabels(s.get\_xticklabels(),rotation=30,fontsize=7) plt.show() 3.0 15 0.0 -1.5In [0]: pca.get\_covariance() Out[0]: array([[ 0.75918838, -0.02083172, 1.39139318, 0.57112452], [-0.02083172, 0.19026705, -0.27999356, -0.10247426], [ 1.39139318, -0.27999356, 3.24250965, 1.36125483], [ 0.57112452, -0.10247426, 1.36125483, 0.61062741]]) As the most darkest spot on the digonal is the petal's length hence this is the feature which dominates our model i.e., most of the the data is varied due to petal's length and it also being verified by our previous results because using petal's length as the only predictor for the model we get around 90% accuracy. Then in the off-diagonal petal length and petal width covar is highest which tells that they are highly related i.e., if one increases then other increases and same for decrease and also this pair has the highest covar hence classify data better than any other pair and we also saw that at the very begining last PC contributes less so we take first two In [0]:  $X_pca = X_pca[:, 0:2]$  $X_{test} = X_{test}[:, 0:2]$ In [0]: **from sklearn.neighbors import** KNeighborsClassifier knn = KNeighborsClassifier(n\_neighbors=4) knn.fit(X\_pca,y\_train) perc\_accu = knn.score(X\_test,y\_test) \* 100 print(f"perc accu : {perc\_accu}") perc accu : 97.36842105263158 The below plot clearly shows that how the first two PC's seperates the three classes in different slots which makes a clear decision boundary between them leads to an easy and efficient classification In [0]: fig = plt.figure() ax1 = fig.add\_subplot(111) f1x=list() f1y=list() f2x=list() f2y=list() f3x=list() f3y=list() Xax=X\_pca[:,0] Yax=X\_pca[:,1] # Yax = [i for i in range(len(Xax))]for (i,j) in zip(Xax,Yax): var = knn.predict([[i,j]]) **if** var==0: f1x.append(i) fly.append(j) elif var==1: f2x.append(i) f2y.append(j) else: f3x.append(i) f3y.append(j) ax1.scatter(f1y, f1x, s=40, alpha=0.4, c='b', marker="\*", label='setosa') ax1.scatter(f2y, f2x, s=40, alpha=0.4, c='r', marker="o", label='versicolor') ax1.scatter(f3y, f3x, s=40, alpha=0.4, c='g', marker="+", label='virginica') plt.show() 4 3 2 1 0 -1 -2 -3 -1.0In [0]: import seaborn as sns the correlation heatmap of features df\_iris = pd.DataFrame(iris\_data.data,columns=iris\_data.feature\_names) s=sns.heatmap(df\_iris.corr(),cmap='coolwarm') s.set\_yticklabels(s.get\_yticklabels(),rotation=30,fontsize=7) s.set\_xticklabels(s.get\_xticklabels(),rotation=30,fontsize=7) plt.show() 0.8 0.4 0.0 -0.4-0.8The heatmap shows that petal length and petal width are the highest correlated pair of distinct features and this was also the case in above cells and petal length and sepal width is the least correlated or closest to orthogonal pair **USING SVD** In [0]: | iris\_data = load\_iris()  $A = USV^T$ In [0]: A = iris\_data.data

In [0]: from sklearn.datasets import load\_iris

