Perform following Data Pre-processing tasks using python. Data reduction using variance threshold, univariate feature selection, recursive feature elimination, PCA.

```
import numpy as np
import matplotlib.pyplot as plt
from sklearn.model selection import train test split
from sklearn.feature selection import VarianceThreshold, RFE, SelectFromModel, SelectKBest
from sklearn.linear_model import LogisticRegression
from sklearn.ensemble import RandomForestClassifier
from sklearn.metrics import confusion_matrix, classification_report
from sklearn.datasets import load_iris
iris = load_iris()
X = iris.data
y = iris.target
print(X[1, :])
print(X.shape)
     [4.9 3. 1.4 0.2]
     (150, 4)
# In order to test the effectness of different feature selection methods, we add some nois
np.random.seed(100)
E = np.random.uniform(0, 1, size=(len(X), 10))
X = np.hstack((X, E))
X.shape
     (150, 14)
X_train, X_test, y_train, y_test = train_test_split(X, y, random_state=1, test_size=0.3)
X train.shape
     (105, 14)
```

Variance Thresholding

VarianceThreshold is a simple baseline approach to feature selection. It removes all features whose variance doesn't meet some threshold. By default, it removes all zero-variance features

```
sel_variance_threshold = VarianceThreshold()
X_train_remove_variance = sel_variance_threshold.fit_transform(X_train)
X_train_remove_variance.shape
# Variance threshold is applied but since the noise valued columns have non-zero variance,
```

(105, 14)

Univariate Feature Selection

Univariate feature selection works by selecting the best features based on univariate statistical tests. We compare each feature to the target variable, to see whether there is statistially significant relationship between them. When we analyze the relationship between one feature and the target variable we ignore the other features. That is why it is called 'univariate'. Each feature has its own test score. Finally, all the test scores are compared, and the features with top scores will be selected. These objects take as input a scoring function that returns univariate scores and p-values (or only scores for SelectKBest and SelectPercentile):

For regression: f_regression, mutual_info_regression > For classification: chi2, f_classif, mutual_info_classif

```
# Since this is iris dataset we are working on, we will use classfication techniques
# Analysis Of Variance (ANOVA)
sel_f = SelectKBest(f_classif, k=4)
X_train_f = sel_f.fit_transform(X_train, y_train)
print(sel_f.get_support())
print(sel_f.get_params())
               [ True True True False False False False False False False False
                 False False]
              {'k': 4, 'score_func': <function f_classif at 0x7fe618c788c0>}
# Chi2 Test
sel_chi2 = SelectKBest(chi2, k=4) # select 4 features
X train chi2 = sel chi2.fit transform(X train, y train)
print(sel_chi2.get_support())
print(sel chi2.get params())
               [ True True True False False False False False False False False
                 False False
              {'k': 4, 'score_func': <function chi2 at 0x7fe618c784d0>}
# mutual info classif Test
sel mutual = SelectKBest(mutual info classif, k=4)
X_train_mutual = sel_mutual.fit_transform(X_train, y_train)
print(sel_mutual.get_support())
print(sel_mutual.get_params())
               [ True True True False F
                 False False
               {'k': 4, 'score_func': <function mutual_info_classif at 0x7fe616705b00>}
```

Recursive Feature Elimination

Given an external estimator that assigns weights to features (e.g., the coefficients of a linear model), recursive feature elimination (RFE) is to select features by recursively considering smaller and smaller sets of features. First, the estimator is trained on the initial set of features and the importance of each feature is obtained either through a coef_ attribute or through a featureimportances attribute. Then, the least important features are pruned from current set of features. That procedure is recursively repeated on the pruned set until the desired number of features to select is eventually reached.

```
model_tree = RandomForestClassifier(random_state=100, n_estimators=50)
sel_rfe_tree = RFE(estimator=model_tree, n_features_to_select=4, step=1)
X_train_rfe_tree = sel_rfe_tree.fit_transform(X_train, y_train)
print(sel_rfe_tree.get_support())
```

[True False True True False False]

- PCA

Since PCA yields a feature subspace that maximizes the variance along the axes, it makes sense to standardize the data, especially, if it was measured on different scales. Although, all features in the Iris dataset were measured in centimeters, let us continue with the transformation of the data onto unit scale (mean=0 and variance=1), which is a requirement for the optimal performance of many machine learning algorithms

```
from sklearn.decomposition import PCA
from sklearn.preprocessing import StandardScaler
import pandas as pd

x = StandardScaler().fit_transform(X)
features = ['sepal length', 'sepal width', 'petal length', 'petal width']

#pd.DataFrame(data = x, columns = features).head()
pd.DataFrame(data = x).head()
```

principalDf.head(5)

4

principal component 1 principal component 2 0 -2.144349 1.254377 1 -2.226538 -1.647474 2 -2.238513 -0.668595 3 -2.501835 -2.506009

-2.550708

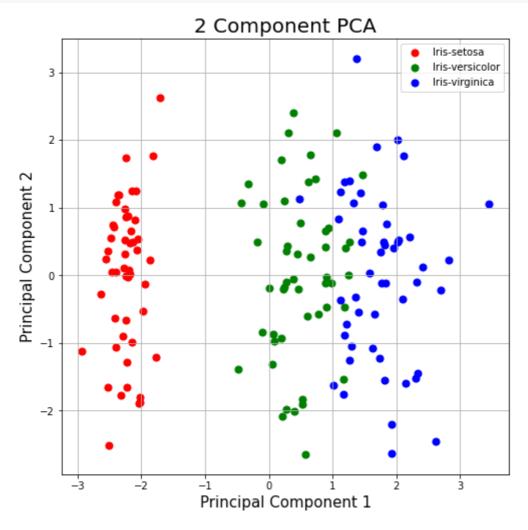
0.237216

```
finalDf = pd.concat([principalDf, df[['target']]], axis = 1)
finalDf.head(5)
```

	principal component 1	principal component 2	target
0	-2.144349	1.254377	Iris-setosa
1	-2.226538	-1.647474	Iris-setosa
2	-2.238513	-0.668595	Iris-setosa
3	-2.501835	-2.506009	Iris-setosa
4	-2.550708	0.237216	Iris-setosa

```
fig = plt.figure(figsize = (8,8))
ax = fig.add_subplot(1,1,1)
ax.set_xlabel('Principal Component 1', fontsize = 15)
ax.set_ylabel('Principal Component 2', fontsize = 15)
ax.set_title('2 Component PCA', fontsize = 20)

targets = ['Iris-setosa', 'Iris-versicolor', 'Iris-virginica']
colors = ['r', 'g', 'b']
```



₽		principal component 1	principal component 2	principal component 3	target
	0	-2.144349	1.254377	0.260207	Iris-setosa
	1	-2.226538	-1.647474	-2.488839	Iris-setosa
	2	-2.238513	-0.668595	2.283491	Iris-setosa
	3	-2.501835	-2.506009	-1.957746	Iris-setosa
	4	-2.550708	0.237216	-0.988219	Iris-setosa

```
import plotly.express as px
targets = ['Iris-setosa', 'Iris-versicolor', 'Iris-virginica']
colors = ['r', 'g', 'b']
fig = px.scatter_3d(finalDf1,x="principal component 1",y="principal component 2",z="princifig.show()
```

- target=Iris-setosa
- target=Iris-versicolor
- target=Iris-virginica

Differences between before and after using Feature Selection

```
#Before using Feature Selection
model_logistic = LogisticRegression(multi_class='multinomial', max_iter=1000)
model_logistic.fit(X_train, y_train)
predict = model_logistic.predict(X_test)
print(confusion_matrix(y_test, predict))

print(classification_report(y_test, predict))

[[14      0      0]
      [      0      16      2]
      [      0      1      12]]
```

precision recall f1-score support

	0	1.00	1.00	1.00	14
	1	0.94	0.89	0.91	18
	2	0.86	0.92	0.89	13
accur	acy			0.93	45
macro	avg	0.93	0.94	0.93	45
weighted	avg	0.94	0.93	0.93	45

```
#After using Feature Selection
model_logistic = LogisticRegression(solver='saga', multi_class='multinomial', max_iter=100
model_logistic.fit(X_train_f, y_train)
```

LogisticRegression(C=1.0, class_weight=None, dual=False, fit_intercept=True, intercept_scaling=1, l1_ratio=None, max_iter=10000, multi_class='multinomial', n_jobs=None, penalty='l2', random_state=None, solver='saga', tol=0.0001, verbose=0, warm_start=False)

```
X_test_f = sel_f.transform(X_test)
print(X_test.shape)
print(X_test_f.shape)
```

(45, 14) (45, 4)

[[14 0 0]

```
predict = model_logistic.predict(X_test_f)
print(confusion_matrix(y_test, predict))
print(classification_report(y_test, predict))
```

[0 17 1] [0 0 13]]				
	precision	recall	f1-score	support
0	1.00	1.00	1.00	14
1	1.00	0.94	0.97	18
2	0.93	1.00	0.96	13
accuracy			0.98	45
macro avg	0.98	0.98	0.98	45
weighted avg	0.98	0.98	0.98	45

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