```
In [1]: import numpy as np
        import pandas as pd
        import matplotlib.pyplot as plt
        import seaborn as sns
        from sklearn.model selection import train test split
        from sklearn.preprocessing import StandardScaler
        from sklearn import datasets
        boston = datasets.load_boston()
        features = pd.DataFrame(boston.data, columns=boston.feature_names)
```

In [2]: print(features.shape)

(506, 13)

## Out [2]:

	CRIM	ZN	INDUS	CHAS	NOX	RM	AGE	DIS	RAD	TAX	PTRATIO	В	LSTAT
0	0.00632	18.0	2.31	0.0	0.538	6.575	65.2	4.0900	1.0	296.0	15.3	396.90	4.98
1	0.02731	0.0	7.07	0.0	0.469	6.421	78.9	4.9671	2.0	242.0	17.8	396.90	9.14
2	0.02729	0.0	7.07	0.0	0.469	7.185	61.1	4.9671	2.0	242.0	17.8	392.83	4.03
3	0.03237	0.0	2.18	0.0	0.458	6.998	45.8	6.0622	3.0	222.0	18.7	394.63	2.94
4	0.06905	0.0	2.18	0.0	0.458	7.147	54.2	6.0622	3.0	222.0	18.7	396.90	5.33

## In [ ]:

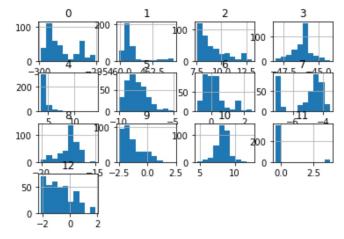
In [20]: # Now we have to standardize the data. We will use StandardScaler from preprocessing

X\_train, X\_test, y\_train, y\_test=train\_test\_split(features, targets, test\_size=.25, random\_s scaler= StandardScaler().fit(X train)

X train scaled=pd.DataFrame(scaler.transform(X train),index=X train.index.values,colum

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```
In [16]: # we can even check if we can reduce dimensionality of dataset using PCA, although her
%matplotlib inline
from sklearn.decomposition import PCA
pca=PCA()
pca.fit(X_train)
foo=pd.DataFrame(pca.transform(X_train))
x_axis=np.arange(1,pca.n_components_ + 1)
pca_scaled=PCA()
pca_scaled.fit(X_train_scaled)
foo_scaled=pd.DataFrame(pca.transform(X_train_scaled))
foo_scaled.hist()
```



```
In [18]: # import ,instantiate,fit
from sklearn.ensemble import RandomForestRegressor
forest=RandomForestRegressor(n_estimators=1000,oob_score=True,random_state=18)
# oob_True :whether to use out-of-bag samples to estimate
# the R^2 on unseen data.is kind of cross-validation
```

C:\Users\HP\Anaconda3\lib\site-packages\ipykernel\_launcher.py:6: DataConversionWar ning: A column-vector y was passed when a 1d array was expected. Please change the shape of y to (n\_samples,), for example using ravel().

```
Out[18]: RandomForestRegressor(bootstrap=True, criterion='mse', max_depth=None, max_features='auto', max_leaf_nodes=None, min_impurity_decrease=0.0, min_impurity_split=None, min_samples_leaf=1, min_samples_split=2, min_weight_fraction_leaf=0.0, n_estimators=1000, n_jobs=1, oob_score=True, random_state=18, verbose=0, warm_start=False)
```

```
In [19]: from sklearn.metrics import r2_score
    preds=forest.predict(X_test)
    test_score=r2_score(y_test,preds)
```

R squared score: 0.8683346228462773

In [ ]:

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