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BE-ETRX UID:2019110039 Sub-Minor ML

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Ise 2

<u>Aim:</u> To implement logistic regression, K nearest neighbor and support vector machine algorithms on the given credit card dataset.

Observation:

- To identify and deal with different samplings.
- To justify and compare the output obtained through different algorithms.

Code:

```
[4] import numpy as np
import pandas as pd

[5] from google.colab import drive
    drive.mount('/content/drive')

    Drive already mounted at /content/drive; to attemp

path = "/content/creditcard.csv"
    df = pd.read_csv(path)
```

df.isnull().sum() #checking for null values

Time	0
	0
V2	0
V 3	0
V4	0
V 5	0
V 6	0
V 7	0
V8	0
V 9	0
V1 V2 V3 V4 V5 V6 V7 V8 V9 V10 V11 V12 V13 V14 V15 V16 V17 V18 V19 V20 V21 V22 V23 V24 V25	0
V11	0
V12	0
V1 3	0
V14	0
V15	0
V16	0
V17	0
V18	1
V19	1 1
V20	
V21	1
V22	1
V23	1
V24	1
	1
V26	1
V27	1
V28	1
Amount	1
Class	
dtype:	1nt64

```
df['Class'].value_counts()[0] #number of normal transactions
 5970
 df['Class'].value_counts()[1] #number of fraud transactions
from sklearn.preprocessing import RobustScaler
rob scaler = RobustScaler()
df['scaled amount'] =
rob scaler.fit transform(df['Amount'].values.reshape(-1,1))
df['scaled time'] =
rob scaler.fit transform(df['Time'].values.reshape(-1,1))
df.drop(['Time','Amount'], axis=1, inplace=True)
df = df.sample(frac=1)
# amount of fraud classes 492 rows.
fraud df = df.loc[df['Class'] == 1]
non fraud df = df.loc[df['Class'] == 0][:492]
normal distributed df = pd.concat([fraud df, non fraud df])
```

new df = normal distributed df.sample(frac=1, random state=42)

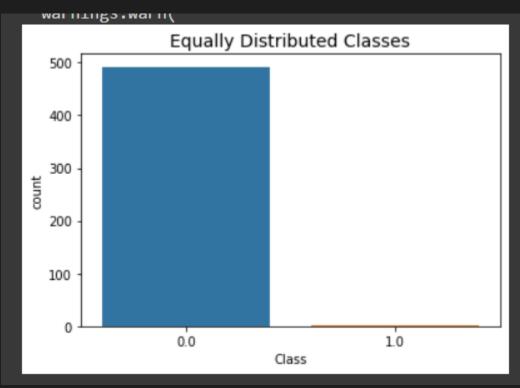
new df.head()

```
print('No Frauds', round(df['Class'].value_counts()[0]/len(df) * 100,2), '% of the dataset')
print('Frauds', round(df['Class'].value_counts()[1]/len(df) * 100,2), '% of the dataset')

X = df.drop('Class', axis=1)
y = df['Class']

No Frauds 99.93 % of the dataset
Frauds 0.05 % of the dataset

import seaborn as sns
import matplotlib.pyplot as plt
print('Distribution of the Classes in the subsample dataset')
print(new_df['Class'].value_counts()/len(new_df))
sns.countplot('Class', data=new_df)
plt.title('Equally Distributed Classes', fontsize=14)
plt.show()
```

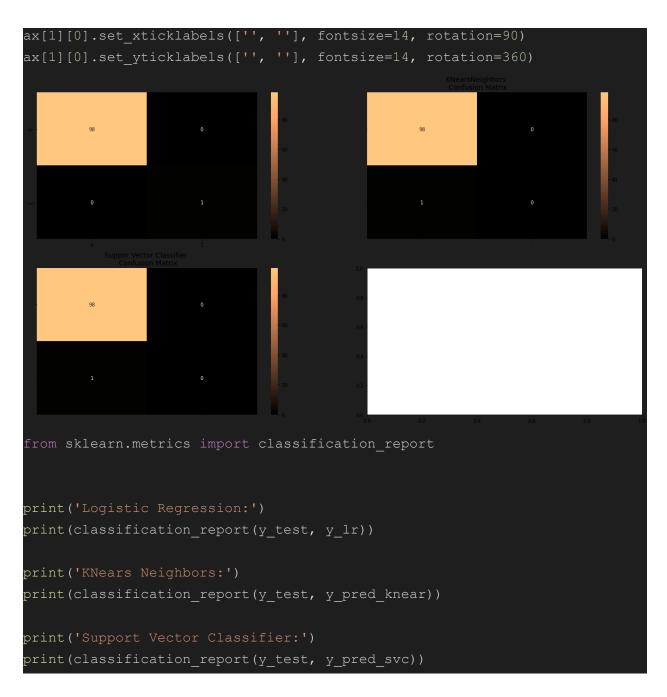


```
fig, ax = plt.subplots(figsize=(30,15))
sub_sample_corr = new_df.corr()
sns.heatmap(sub_sample_corr)
<matplotlib.axes._subplots.AxesSubplot at 0x7fd9f3cdf160>
       V1 -
       V2
       V3
       ٧4
       V5
       V6
       ٧7
       VΒ
       V9
      V10
      V11
      V12
      V13
      V14
      V15
      V16
      V17
      V18
      V19
      V20
      V21
      V22
      V23
      V24
```

```
X = new df.drop('Class', axis=1)
 y = new df['Class']
 from sklearn.model selection import train test split
 # This is explicitly used for undersampling.
 X train, X test, y train, y test = train test split(X, y, test size=0.2, random state=42)
 X train = X train.values
 X test = X test.values
 y train = y train.values
 y_test = y_test.values
 from sklearn.linear model import LogisticRegression
 from sklearn.svm import SVC
 from sklearn.neighbors import KNeighborsClassifier
 from sklearn.ensemble import RandomForestClassifier
 from sklearn.naive_bayes import GaussianNB
 classifiers = {
    "LogisiticRegression": LogisticRegression(),
    "KNearest": KNeighborsClassifier(),
    "Support Vector Classifier": SVC(),
    "RandomForestClassifier": RandomForestClassifier(),
    "Naive Bayes": GaussianNB()
from sklearn.model selection import cross val score
for key, classifier in classifiers.items():
    classifier.fit(X train, y train)
    training score = cross val score(classifier, X train, y train, cv=5)
score of", round(training score.mean(), 2) * 100, "% accuracy score")
from sklearn.model selection import GridSearchCV
log reg params = {"penalty": ['l1', 'l2'], 'C': [0.001, 0.01, 0.1, 1, 10,
100, 1000]}
grid log reg = GridSearchCV(LogisticRegression(), log reg params)
grid log reg.fit(X train, y train)
log reg = grid log reg.best estimator
```

```
knears params = {"n neighbors": list(range(2,5,1)),                           'algorithm': ['auto',
grid knears = GridSearchCV(KNeighborsClassifier(), knears params)
grid knears.fit(X train, y train)
knears neighbors = grid knears.best estimator
svc params = {'C': [0.5, 0.7, 0.9, 1], 'kernel': ['rbf', 'poly',
grid svc = GridSearchCV(SVC(), svc params)
grid svc.fit(X train, y train)
svc = grid svc.best estimator
log reg score = cross val score(log reg, X train, y train, cv=5)
print('Logistic Regression Cross Validation Score: ',
round(log reg score.mean() * 100, 2).astype(str) + '%')
knears score = cross val score(knears neighbors, X train, y train, cv=5)
print('Knears Neighbors Cross Validation Score', round(knears score.mean()
100, 2).astype(str) + '%')
svc score = cross val score(svc, X train, y train, cv=5)
print('Support Vector Classifier Cross Validation Score',
round(svc score.mean() * 100, 2).astype(str) + '%')
from sklearn.metrics import roc curve
from sklearn.model selection import cross val predict
log reg pred = cross val predict(log reg, X train, y train, cv=5,
                             method="decision function")
knears pred = cross val predict(knears neighbors, X train, y train, cv=5)
svc pred = cross val predict(svc, X train, y train, cv=5,
                             method="decision function")
from sklearn.metrics import roc auc score
print('Logistic Regression: ', roc auc score(y train, log reg pred))
print('KNears Neighbors: ', roc auc score(y train, knears pred))
print('Support Vector Classifier: ', roc auc score(y train, svc pred))
from sklearn.metrics import precision recall curve
```

```
precision, recall, threshold = precision recall curve(y train,
log reg pred)
from sklearn.metrics import recall score, precision score, f1 score,
accuracy score
y pred = log reg.predict(X train)
print('Recall Score: {:.2f}'.format(recall score(y train, y pred)))
print('Precision Score: {:.2f}'.format(precision score(y train, y pred)))
print('F1 Score: {:.2f}'.format(f1 score(y train, y pred)))
print('Accuracy Score: {:.2f}'.format(accuracy score(y train, y pred)))
print('---' * 45)
from sklearn.metrics import confusion matrix
clf = LogisticRegression(random state=0).fit(X, y)
y lr = clf.predict(X test)
y pred knear = knears neighbors.predict(X test)
y pred svc = svc.predict(X test)
lr cf = confusion matrix(y test, y lr)
kneighbors cf = confusion matrix(y test, y pred knear)
svc cf = confusion matrix(y test, y pred svc)
fig, ax = plt.subplots(2, 2, figsize=(22, 12))
sns.heatmap(lr cf, ax=ax[0][0], annot=True, cmap=plt.cm.copper)
ax[0][1].set title("L.R. \n Confusion Matrix", fontsize=14)
ax[0][1].set xticklabels(['', ''], fontsize=14, rotation=90)
ax[0][1].set yticklabels(['', ''], fontsize=14, rotation=360)
sns.heatmap(kneighbors cf, ax=ax[0][1], annot=True, cmap=plt.cm.copper)
ax[0][1].set title("KNearsNeighbors \n Confusion Matrix", fontsize=14)
ax[0][1].set xticklabels(['', ''], fontsize=14, rotation=90)
ax[0][1].set_yticklabels(['', ''], fontsize=14, rotation=360)
sns.heatmap(svc cf, ax=ax[1][0], annot=True, cmap=plt.cm.copper)
ax[1][0].set title("Suppor Vector Classifier \n Confusion Matrix",
fontsize=14)
```



Output:

			ression:	Logistic Reg
support	f1-score	recall	precision	
98	1.00	1.00	1.00	0.0
1	1.00	1.00	1.00	1.0
99	1.00			accuracy
99	1.00	1.00	1.00	macro avg
99	1.00	1.00	1.00	weighted avg

KNears Neighbors: recall 0.99 0.99 98 0.99 0.50 macro avg 0.49 99 0.98 weighted avg 0.99 0.98 Support Vector Classifier: recall f1-score support 0.99 0.99 0.00 0.00 0.00 0.99 accuracy 0.49 0.50 macro avg 0.98 0.99 0.98

Conclusion:

- We almost got the same results in both KNN and SVM.
- We got the maximum accuracy in the logistic regression algorithm.
- We successfully got the expected confusion matrix for all of the three algorithms.
- Logistic regression is a statistical method that is used for building machine learning models where the dependent variable is dichotomous: i.e. binary.
 Logistic regression is used to describe data and the relationship between one dependent variable and one or more independent variables.
- The k-nearest neighbors algorithm, also known as KNN or k-NN, is a non-parametric, supervised learning classifier, which uses proximity to make classifications or predictions about the grouping of an individual data point.
- Support Vector Machine or SVM is one of the most popular Supervised Learning algorithms, which is used for Classification as well as Regression problems. However, primarily, it is used for Classification problems in Machine Learning.
- The goal of the SVM algorithm is to create the best line or decision boundary that
 can segregate n-dimensional space into classes so that we can easily put the
 new data point in the correct category in the future. This best decision boundary
 is called a hyperplane.