n [1]:	
1 [1].	<pre>import pandas as pu import numpy as np import matplotlib.pyplot as plt import seaborn as sns from sklearn.preprocessing import LabelEncoder from sklearn.utils import resample from xgboost import XGBClassifier from sklearn.metrics import roc_auc_score,precision_recall_curve,roc_curve from sklearn.model_selection import train_test_split from sklearn.linear_model import LogisticRegression from sklearn.manifold import TSNE from sklearn.decomposition import PCA, TruncatedSVD import time import matplotlib.patches as mpatches from sklearn.metrics import confusion_matrix</pre>
n [2]:	Loading Data %%time train_transactions=pd.read_csv('/input/train_transaction.csv') train_identity=pd.read_csv('/input/train_identity.csv') print('Train data set is loaded !') Train data set is loaded ! CPU times: user 18.7 s, sys: 2.66 s, total: 21.4 s Wall time: 21.6 s Getting basic Idea
n [3]: nt[3]:	TransactionID isFraud TransactionDT TransactionAmt ProductCD card1 card2 card3 card4 card5 card6 addr1 addr2 dis 1 2987000 0 86400 68.5 W 13926 NaN 150.0 discover 142.0 credit 315.0 87.0 19 2 2987001 0 86401 29.0 W 2755 404.0 150.0 mastercard 102.0 credit 325.0 87.0 N 2 2987002 0 86469 59.0 W 4663 490.0 150.0 visa 166.0 debit 330.0 87.0 283 3 2987003 0 86499 50.0 W 18132 567.0 150.0 mastercard 117.0 debit 476.0 87.0 N 4 2987004 0 86506 50.0 H 4497 514.0 150.0 mastercard 102.0 credit 420.0 87.0 N train_transactions.info() cclass 'pandas.core.frame.DataFrame'> RangeIndex: 590540 entries, 0 to 590539 Columns: 394 entries, TransactionID to V339 dtypes: float64(376), int64(4), object(14)
n [5]:	train_identity.info() cclass 'pandas.core.frame.DataFrame'> RangeIndex: 144233 entries, 0 to 144232 Data columns (total 41 columns): TransactionID
	id_19
n [6]:	Target variable x=train_transactions['isFraud'].value_counts().values sns.barplot([0,1],x) plt.title('Target variable count') Text(0.5, 1.0, 'Target variable count') Target variable count 500000
	The metric trap One of the major issues that novice users fall into when dealing with unbalanced datasets relates to the metrics used to evaluate their model. Using simpler metrics like accuracy_score can be misleading. In a dataset with highly unbalanced classes, if the classifier always "predicts" the most common class without performing any analysis of the features, it will still have a high accuracy rate, obviously illusory. False Positive. Predict an event when there was no event. False Negative. Predict no event when in fact there was an event. In the overview of the problem statement the organizers has described a situation where you stand at the queue for a long time and when your chance arrives, the transaction gets denied because it was interpreted as a Fraudulent transaction which many of us have faced. This is classical example of False Negative prediction. Change the performance metric As we saw above, accuracy is not the best metric to use when evaluating imbalanced datasets as it can be very misleading. Metrics that can provide better insight include:
	Confusion Matrix: a table showing correct predictions and types of incorrect predictions. Precision: the number of true positives divided by all positive predictions. Precision is also called Positive Predictive Value. It is a measur of a classifier's exactness. Low precision indicates a high number of false positives. Recall: the number of true positives divided by the number of positive values in the test data. Recall is also called Sensitivity or the True Positive Rate. It is a measure of a classifier's completeness. Low recall indicates a high number of false negatives. F1 Score: the weighted average of precision and recall. I don't understand why the competition hosts selected ROC_AUC as evaluation metric,I think ROC curves should be used when there are roughly equal numbers of observations for each class. Precision-Recall curves should be used when there is a moderate to large class imbalance. Merging transaction and identity dataset We will firt merge our transactions and identity datasets.
n [7]:	<pre>train=train_transactions.merge(train_identity, how='left', left_index=True, right_index=True) y_train=train['isFraud'].astype('uint8') print('Train shape', train.shape) del train_transactions, train_identity print("Data set merged ") Train shape (590540, 435) Data set merged</pre>
n [8]:	Reducing memory usage **time # From kernel https://www.kaggle.com/gemartin/load-data-reduce-memory-usage # WARNING! THIS CAN DAMAGE THE DATA def reduce_mem_usage2(df): """ iterate through all the columns of a dataframe and modify the data type to reduce memory usage. """ start_mem = df.memory_usage().sum() / 1024**2 print('Memory usage of dataframe is {:.2f} MB'.format(start_mem)) for col in df.columns: col_type = df[col].dtype if col_type != object: c_min = df[col].min() c_max = df[col].max() if str(col_type)[i3] == 'int': if c_min > np.iinfo(np.int6).min and c_max < np.iinfo(np.int8).max: df[col] = df[col].astype(np.int8) elif c_min > np.iinfo(np.int16).min and c_max < np.iinfo(np.int16).max: df[col] = df[col].astype(np.int16) elif c_min > np.iinfo(np.int32).min and c_max < np.iinfo(np.int32).max: df[col] = df[col].astype(np.int32) elif c_min > np.iinfo(np.int64).min and c_max < np.iinfo(np.int64).max: df[col] = df[col].astype(np.int64)
n [9]:	else: if c_min > np.finfo(np.float16).min and c_max < np.finfo(np.float16).max: df(col) = df(col).astype(np.float32).min and c_max < np.finfo(np.float32).max: df(col) = df(col).astype(np.float32) else: df(col) = df(col).astype(np.float32) else: df(col) = df(col).astype(inp.float64) else: df(col) = df(col).astype(inp.float64) else: df(col) = df(col).astype(inp.float64) end_mem = df.memory_usage().sum() / 1024**2 print('Memory usage after optimization is: {:.2f} MB'.format(end_mem)) print('Decreased by {:.1f}%'.format(100 * (start_mem - end_mem) / start_mem)) return df CPU times: user 0 ns, sys: 0 ns, total: 0 ns Wall time: 248 µs %%time train = reduce_mem_usage2(train) Memory usage after optimization is: 527.82 MB Decreased by 73.1% CPU times: user 52.3 s, sys: lmin 15s, total: 2min 8s Wall time: 2min 8s Splitting to train and validation
[10]:	We will now split the train dataset into train and validation set. We will keeep 20% of data for validation. X_train, X_test, y_train, y_test=train_test_split(train.drop('isFraud',axis=1), y_train,test_size=.2, rand_state=1) Resampling A widely adopted technique for dealing with highly unbalanced datasets is called resampling. It consists of removing samples from the majority class (under-sampling) and / or adding more examples from the minority class (over-sampling). Undersampling Oversampling Copies of the minority class
[11]:	Original dataset Resampling Techniques using sklearn 1. Oversample minority class Oversampling can be defined as adding more copies of the minority class. Oversampling can be a good choice when you don't have a to of data to work with. We will use the resampling module from Scikit-Learn to randomly replicate samples from the minority class. X=pd.concat([X_train, y_train], axis=1) not_fraud=X[X.isFraud==0] fraud=X[X.isFraud==1]
[12]: [12]:	0 455903 Name: isFraud, dtype: int64
	upsampled data class count 400000 200000 200000 100000 200000 200000 200000 100000 200000 200000 200000 200000 200000 200000 200000 200000 200000 2000000
[13]: [14]:	0 16529 Name: isFraud, dtype: int64
[15]:	We will review other resampling techniques. For ease of visualization, let's create a small unbalanced sample dataset using the make_classification method: from sklearn.datasets import make_classification X, y = make_classification(n_classes=2, class_sep=1.5, weights=[0.9, 0.1], n_informative=3, n_redundant=1, flip_y=0, n_features=20, n_clusters_per_class=1, n_samples=1000, random_state=10) df = pd.DataFrame(X)
:[15]:	df ['target'] = y df.target.value_counts().plot(kind='bar', title='Count (target)') <matplotlib.axessubplots.axessubplot 0x7fc68e72d550="" at=""> Count (target) 600 400 200 • We will do an experiment with this data without any resampling technique. • We will fit and predict the data on a Logistic regression model and observe the output scores.</matplotlib.axessubplots.axessubplot>
[17]:	<pre>def logistic(X,y): X_train,X_test,y_train,y_test=train_test_split(X,y,test_size=.2,random_state=1) lr=LogisticRegression() lr.fit(X_train,y_train) prob=lr.predict_proba(X_test) return (prob[:,1],y_test) probs,y_test=logistic(X,y) /opt/conda/lib/python3.6/site-packages/sklearn/linear_model/logistic.py:432: FutureWarning: Default s olver will be changed to 'lbfgs' in 0.22. Specify a solver to silence this warning. FutureWarning) • We will define two functions to plot precision_recall curve and roc curve def plot_pre_curve(y_test,probs): precision, recall, thresholds = precision_recall_curve(y_test, probs) plt.plot([0, 1], [0.5, 0.5], linestyle='') # plot the precision-recall curve for the model plt.plot(recall, precision, marker='.') plt.title("precision recall curve") plt.xlabel('Recall')</pre>
[19]:	<pre>plt.ylabel('Precision') # show the plot plt.show() def plot_roc(y_test,prob): fpr, tpr, thresholds = roc_curve(y_test, probs) # plot no skill plt.plot([0, 1], [0, 1], linestyle='') # plot the roc curve for the model plt.plot(fpr, tpr, marker='.') plt.title("ROC curve") plt.xlabel('false positive rate') # show the plot plt.show()</pre> plot_pre_curve(y_test,probs) precision recall curve 10 09
[20]:	Diot_roc (y_test, probs) ROC curve ROC curve
[21]:	We will also create a 2-dimensional plot function, plot_2d_space, to see the data distribution: def plot_2d_space(X_train, y_train, X=X, y=y ,label='Classes'): colors = ['#1F77B4', '#FF7F0E'] markers = ['o', 's'] fig, (ax1,ax2)=plt.subplots(1,2, figsize=(8,4)) for l, c, m in zip(np.unique(y), colors, markers): ax1.scatter(X_train[y_train==1, 0], X_train[y_train==1, 1],
	c=c, label=1, marker=m) for 1, c, m in zip(np.unique(y), colors, markers): ax2.scatter(X[y==1, 0], X[y==1, 1], c=c, label=1, marker=m) ax1.set_title(label) ax2.set_title('original data') plt.legend(loc='upper right') plt.show() Dimensionality Reduction and Clustering Understanding t-SNE: In order to understand this algorithm you have to understand the following terms: Euclidean Distance Conditional Probability Normal and T-Distribution Plots
[22]:	<pre>In the below section we will implement three major dimensionality reduction algorithms</pre>
[23]:	<pre>print("Truncated SVD took {:.2} s".format(t1 - t0)) T-SNE took 6.9 s PCA took 0.068 s Truncated SVD took 0.0028 s Now we will visualize the output of the above three algorithms in a 2D space. f, (ax1, ax2, ax3) = plt.subplots(1, 3, figsize=(24,6)) # labels = ['No Fraud', 'Fraud'] f.suptitle('Clusters using Dimensionality Reduction', fontsize=14) blue_patch = mpatches.Patch(color='#0A0AFF', label='No Fraud')</pre>
	<pre>red_patch = mpatches.Patch(color='#AF0000', label='Fraud') # t-SNE scatter plot ax1.scatter(X_reduced_tsne[:,0], X_reduced_tsne[:,1], c=(y == 0), cmap='coolwarm', label='No Fraud', newidths=2) ax1.scatter(X_reduced_tsne[:,0], X_reduced_tsne[:,1], c=(y == 1), cmap='coolwarm', label='Fraud', lin idths=2) ax1.seat_title('t-SNE', fontsize=14) ax1.grid(True) ax1.legend(handles=[blue_patch, red_patch]) # PCA scatter plot ax2.scatter(X_reduced_pca[:,0], X_reduced_pca[:,1], c=(y == 0), cmap='coolwarm', label='No Fraud', lin widths=2) ax2.scatter(X_reduced_pca[:,0], X_reduced_pca[:,1], c=(y == 1), cmap='coolwarm', label='Fraud', linew ths=2) ax2.grid(True) ax2.legend(handles=[blue_patch, red_patch]) # TruncatedSVD scatter plot ax3.scatter(X_reduced_svd[:,0], X_reduced_svd[:,1], c=(y == 0), cmap='coolwarm', label='No Fraud', lin widths=2) ax3.scatter(X_reduced_svd[:,0], X_reduced_svd[:,1], c=(y == 1), cmap='coolwarm', label='Fraud', linew ths=2) ax3.scatter(X_reduced_svd[:,0], X_reduced_svd[:,1], c=(y == 1), cmap='coolwarm', label='Fraud', linew ths=2) ax3.scatter(X_reduced_svd[:,0], X_reduced_svd[:,1], c=(y == 1), cmap='coolwarm', label='Fraud', linew ths=2) ax3.grid(True)</pre>
[24]:	Python imbalanced-learn module A number of more sophisticated resampling techniques have been proposed in the scientific literature. For example, we can cluster the records of the majority class, and do the under-sampling by removing records from each cluster, thus seeking to preserve information. In over-sampling, instead of creating exact copies of the minority class records, we can introduce small variations into those copies, creating more diverse synthetic samples. Let's apply some of these resampling techniques, using the Python library imbalanced-learn. It is compatible with scikit-learn and is part scikit-learn-contrib projects.
[25]:	Random under-sampling with imbalanced-learn from imblearn.under_sampling import RandomUnderSampler ran=RandomUnderSampler(return_indices=True) ##intialize to return indices of dropped rows X_rs,y_rs,dropped = ran.fit_sample(X,y) print("The number of removed indices are ",len(dropped)) plot_2d_space(X_rs,y_rs,X,y,'Random under sampling') The number of removed indices are 200 Random under sampling original data
[26]:	plot_pre_curve(y_test,probs) /opt/conda/lib/python3.6/site-packages/sklearn/linear_model/logistic.py:432: FutureWarning: Default s olver will be changed to 'lbfgs' in 0.22. Specify a solver to silence this warning. FutureWarning) precision recall curve 10
[27]:	plot_roc (y_test, probs) ROC curve 10 08 04 06 05 07 06 07 06 07 08 08 08 08 08 08 08 08 08 08 08 08 08
[28]:	Random over-sampling with imbalanced-learn from imblearn.over_sampling import RandomOverSampler ran=RandomOverSampler() X_ran,y_ran= ran.fit_resample(X,y) print('The new data contains {} rows '.format(X_ran.shape[0])) plot_2d_space(X_ran,y_ran,X,y,'over-sampled') The new data contains 1800 rows over-sampled original data original data

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Algorithmic Ensemble Techniques The above section, deals with handling imbalanced data by resampling original data to provide balanced classes. In this section, we give to look at an attentive approach i.e. Needlying existing classification algorithms to make them appropriate for imbalanced data and the stage classifiers. The approach involves constructive everal two stage classifiers from the original data and then aggregate their prediction Cache awareness and out-of-core computing using depth-first approach Tree pruning using depth-first approach Parallelized tree building Tree building Tree pruning using depth-first approach Regularization for avoiding overfitting In-built cross-validation capability XGBoost (Extreme Gradient Boosting) is an advanced and more efficient implementation of Gradient Boosting Algorithm discussed previous section. Advantages over Other Boosting Techniques It is 10 image faster than the normal Gradient Boosting as it implements parallel processing. It is highly flexible as users can define optimization objectives and evaluation orders. It has an inbuilt mechanism to handle missing values. Unlike gradient boosting which splitting a node as soon as it encounters a negative loss. XG Boost splits up to the maximum depth specified and prunes the treat and removes splits beyond which there is an only regalized loss. Extreme gradient boosting can be done using the XGBoost package in R and Python WORK IN PROGRESS If you like my work, please do upvote ^	1.0 - tune bositive rate 0.8 - 0.4 - 0.2 - 0.2 - 0.2 - 0.2 - 0.2 - 0.3 - 0.2 - 0.3 -	ROC curve	
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