# UDACITY MACHINE LEARNING ENGINEER NANODEGREE

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**CAPSTONE REPORT** 

# 1. Data Exploration

When we load the azdias and customers dataset we raise a mixed type warning :

```
# Load in the data
azdias = pd.read_csv('azdias.csv', sep=',')

c:\users\lanth\appdata\local\programs\python\python37\lib\site-packages\IPython\core\interactiveshell.py:3057: DtypeWarning: Co
lumns (19,20) have mixed types. Specify dtype option on import or set low_memory=False.
    interactivity=interactivity, compiler=compiler, result=result)
```

Going forward we need to know what are the names of columns 19 and 20:

```
print(azdias.iloc[:,19:21].columns)
Index(['CAMEO_DEUG_2015', 'CAMEO_INTL_2015'], dtype='object')
```

Now we look at what they hold:

We can see that there are mainly numerical values: int, float and string But there are also some 'X' and 'XX', corresponding to missing values. So I have made a function cameo\_fix which simply replaces the X and XX by np.nan and sets all to float.

We know that customers dataset holds 3 columns that are not in azdias: 'PRODUCT\_GROUP', 'CUSTOMER\_GROUP' and 'ONLINE\_PURCHASE', therefore we will drop these columns as we want to have the same columns in each dataset.

Then we check if we have the same columns in azdias and customers:

```
list(set(azdias.columns) - set(customers.columns))
[]
list(set(customers.columns) - set(azdias.columns))
[]
```

# 2. Data Preprocessing

## 2.1 Data Cleaning

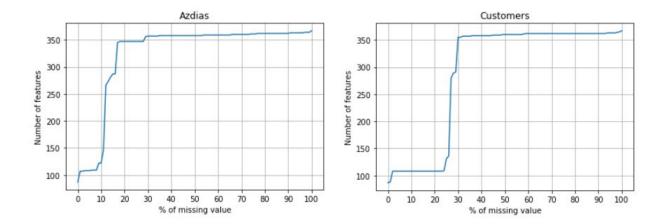
We will start the preprocessing process by implementing a missing\_pct function to calculate the proportion of missing values in each columns

Next we open DIAS Attributes - Values 2017.xlsx to see that it hold missing/unknown value for each column. So we put these into a dataframe to be able to replace the corresponding values by np.nan in each columns

	features	unknowns
0	AGER_TYP	-1
1	ALTERSKATEGORIE_GROB	-1,0,9
2	ALTER_HH	0
3	ANREDE_KZ	-1,0
4	BALLRAUM	-1
5	BIP_FLAG	-1
6	CAMEO_DEUG_2015	-1
7	CAMEO_DEUINTL_2015	-1
8	CJT_GESAMTTYP	0
9	D19_KK_KUNDENTYP	-1
10	EWDICHTE	-1
11	FINANZTYP	-1
12	FINANZ_ANLEGER	-1
13	FINANZ_HAUSBAUER	-1
14	FINANZ_MINIMALIST	-1
15	FINANZ_SPARER	-1
16	FINANZ_UNAUFFAELLIGER	-1
17	FINANZ_VORSORGER	-1
18	GEBAEUDETYP	-1,0
19	GEOSCORE_KLS7	-1,0
20	HAUSHALTSSTRUKTUR	-1,0
21	HEALTH_TYP	-1

To do so I write a missing\_to\_nans function and proceed to replace them and get the new missing\_pct values.

Then I implement a feature\_cap function that returns the list of features that have less missing value proportion than a desired number. This way I can plot the amount of features retained given a specific % of missing maximum:



We can see that we could have most of the azdias's columns with a cap fixed around 20%. But that would have a tremendous consequence on customers since we can see that at 20% less than a third of the customers's columns would be kept. So we need to choose the number based on customers and not azdias.

Therefore as 30% seems to fit with most of the columns in both dataset we will keep this number.

Once we create the features\_selected list for each dataset we compare to see which features are not in both:

```
list(set(azdias_features_selected) - set(customers_features_selected))
['REGIOTYP', 'KKK']

list(set(customers_features_selected) - set(azdias_features_selected))
```

So we drop REGIOTYP and KKK from azdias.

Then I use a features\_engineering to transform the features that are either of the wrong type. Those encoding form more than one thing, we split into 2 new features and drop the original one.

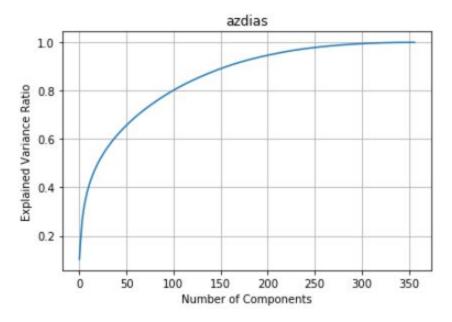
And finally it replaces every np.nan value by the most common value in the column.

### 2.2 Data Scaling/Dimensionality reduction

As there are very different ranges of values in each column, I implement a scaler\_tool function that simply performs a MinMaxScaler.

Then I use Principal component analysis to achieve dimensionality reduction.

As I need to choose a targeted number of component, I use a plot of the sum of the explained variance for each number of components to decide:



From this plot I choose to go for 150 because it explains 90% of the variance because the gap statistic that i plan to use next is very long to compute and 200 components would make it even longer.

# 3. Customers segmentation

### 3.1 Definitions

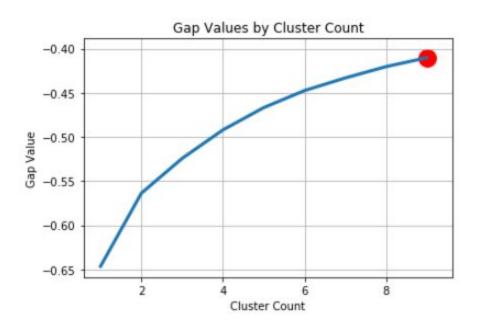
"Customer segmentation is the process of dividing customers into groups based on common characteristics so companies can market to each group effectively and appropriately."

From this definition I choose to use a clustering algorithm with the general demographic dataset (azdias), to then compare in which cluster the customers dataset will fit.

I use K Means as a clustering algorithm, but then I need to choose the optimal K.

## 3.2 Clustering

So I perform a gap statistic analysis (from Tibshirani, Walther, Hastie) with the function optimalK and then plot the gap value by cluster count to deduce that the optimal K is 9:



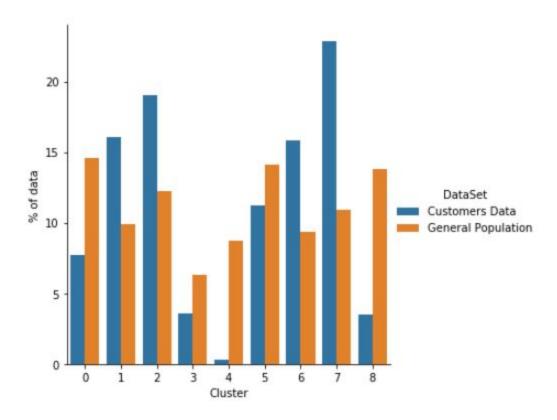
Then the azdias dataset that has been MinMaxed and then PCA(150) is now use to fit a KMeans model with K = 9

### 3.3 Segmentation

Now we have our model fitted we can transform the customers dataset through the same process to see where it ends :

Customers dataset -> preprocessing -> MinMaxScaler -> PCA(150)

Finally we can plot the resulting clusters proportion in each of the two datasets



### From that plot I can say:

- Cluster 7 is the very best segment for customers
- Cluster 1, 2 and 6 are very good also
- Cluster 0, 3, 4 and 8 are very bad.

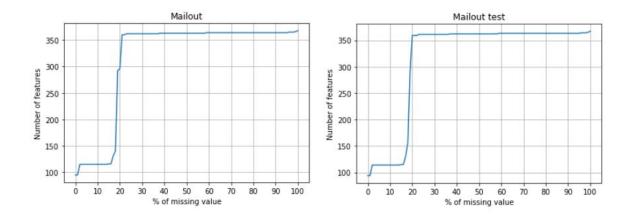
# 4. Supervised Learning

### 4.1 Cleaning/Preprocess

After loading mailout and mailout\_test we will first preprocess it :

- cameo\_fix
- missing\_to\_nans

As previously i will plot the amount of features kept by %of missing value cap :



We can see that about 22% would be enough but I will stay with 30% as the curve seems almost flat past 20%, it will not change much.

```
#Computing the list of columns in mailout_data that have less than 30% of missing values

mailout_data_missing = missing_pct(mailout_data)
mailout_data_features_selected = feature_cap(mailout_data_missing, 30)

#Computing the list of columns in mailout_test that have less than 30% of missing values

mailout_test_missing = missing_pct(mailout_test)
mailout_test_features_selected = feature_cap(mailout_test_missing, 30)

print(len(mailout_data_features_selected))
print(len(mailout_test_features_selected))

358
357
```

We need to check if the columns match:

```
list(set(mailout_data_features_selected) - set(mailout_test_features_selected))
['RESPONSE']

list(set(mailout_test_features_selected) - set(mailout_data_features_selected))
```

Since we know that by definition mailout\_test doesn't have the RESPONSE column, we can say that the columns will match after we separate RESPONSE from the mailout dataset.

- features\_engineering
- MinMaxScaler

At this point the datasets have the same number of columns, we just do a quick check of how imbalanced the class are :

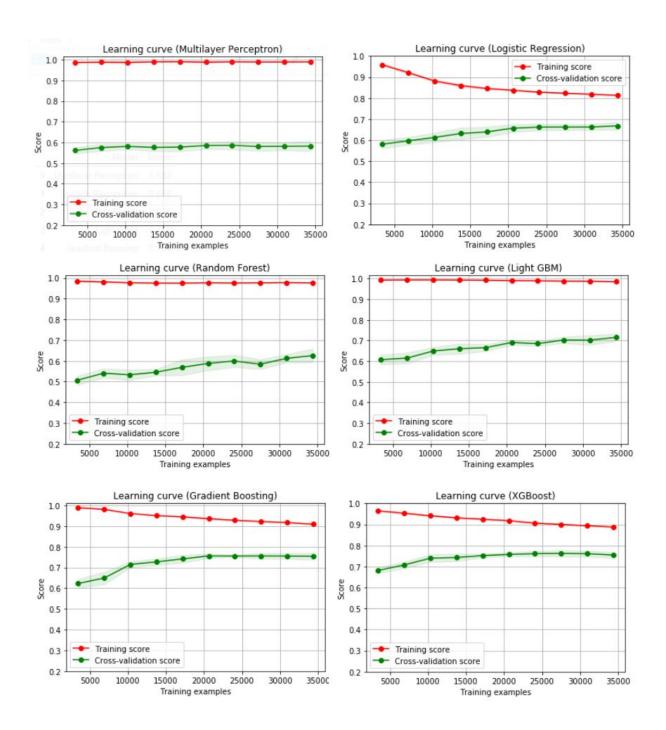
```
#How much in % is there of response
print(len(list(mailout_data.loc[mailout_data['RESPONSE'] == 1].index))/len(mailout_data)*100,"%")
1.2383036171500394 %
```

So the data is extremely imbalanced, we will need to address the issue. I have tried to resample the mailout\_data using imblearn TomekLinks, SMOTE, SMOTETomek and ClusterCentroids but it did not improve the quality of the model's prediction. So I've decided to not resample the data afterall.

### 4.2 Choosing model

To choose which model I will use, I have decided to plot the learning curve of the following models:

- Multi-layer Perceptron classifier
- Logistic Regression
- Random Forest
- Light GBM
- Gradient Boosting Classifier
- XGBoost



The plot\_learning\_curve function also return the average ROC\_AUC score for the tests set:

	Model	Score
0	Multilayer Perceptron	0.582
1	Logistic Regression	0.668
2	Random Forest	0.625
3	Light GBM	0.716
4	Gradient Boosting	0.754
5	XGBoost	0.755

As the a result we can see that XGBoost is the best performing model from scratch (though being closely followed up by Light GBM)
So I will go further with XGBoost.

### 4.3 Hyperparameters tuning

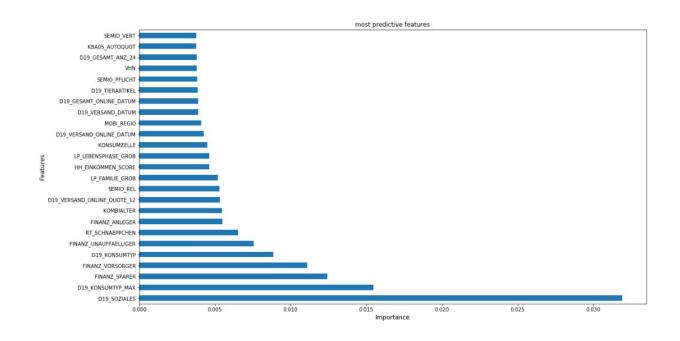
At this point we have only to find the hyperparameters to use for our XGBoost model, in order to have the best rank possible in the kaggle competition.

I have found a similar problem resolved on kaggle: <u>Bayesian Optimization of xgBoost</u> which suggests the use of the BayesSearchCV class from scikit-learn optimize. Though it is extremely time consuming, the results were very good.

The best classifier I have found with this optimization process :

```
XGBClassifier(base_score=0.5, booster='gbtree', colsample_bylevel=1, colsample_bynode=1, colsample_bytree=0.5, eval_metric='auc', gamma=1.0, learning_rate=0.0020276515169578386, max_delta_step=0, max_depth=5, min_child_weight=10, missing=None, n_estimators=251, n_jobs=-1, nthread=None, objective='binary:logistic', random_state=0, reg_alpha=1.0, reg_lambda=1, scale_pos_weight=34, seed=None, silent=None, subsample=0.5, verbosity=1)
```

After a long computation, the resulting model gives us the following features importance overview:



# 4.4 Kaggle competition

Overview	Data Notebooks Discussion Leaderboard Rules Team	My Submissions	Submit Predictions
17	Rahul Dixit	0.80561	7 1y
18	jxtrbtk	0.80557	163 20d
19	anacolada	0.80555	15 1mo
20	FC Su	0.80526	57 7mo
21	Maurizio Santamicone	0.80467	6 1y
22	yueureka	0.80461	1 6mo
23	Jahid Ahsan	0.80453	23 10d
24	NaomiNguyen	0.80444	18 5mo
25	rohan16	0.80389	6 7mo
26	Edu Burgoa	0.80374	14 24d
27	JPBedran	0.80370	5 1mo
28	yo Lanthroff	0.80336	5 8h

When I checked the leaderboard first, I saw that a lot of people achieved a score just above 0.80 and only one did break the 0.81 limit. So my primary goal was also to be in the same range of 0.80+

I am glad I could achieve that.