ps02

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1 Data Science in Practice 2020

1.1 Problem Set 2

1.1.1 Descriptive report

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1.1.2 1. Initialization

Loading modules:

```
[1]: import pandas as pd
     import numpy as np
     import seaborn as sns
     from operator import itemgetter, attrgetter
     from scipy.stats import randint as sp_randint
     from sklearn.model_selection import train_test_split, GridSearchCV, __
     →RandomizedSearchCV, cross_val_score
     from sklearn.metrics import log_loss, classification_report, confusion_matrix,_
     →roc_curve, accuracy_score, roc_auc_score, precision_recall_fscore_support, u
     →precision_recall_curve, auc
     from sklearn.ensemble import RandomForestClassifier
     from sklearn.neighbors import KNeighborsClassifier,
     →NeighborhoodComponentsAnalysis
     from sklearn.preprocessing import StandardScaler
     # import time
     from matplotlib import pyplot as plt
     import argparse
     import logging
     import os
     import sys
     from ipyparallel import Client
     from ipyparallel.joblib import IPythonParallelBackend
```

```
from joblib import Parallel, parallel_backend, register_parallel_backend
```

Defining functions:

```
[2]: def df_pp():
         df_raw = pd.read_csv('customers.csv')
         # clean up lines
         total_length = len(df_raw)
         df_raw = df_raw[df_raw['TotalCharges'] != ' ']
         df_raw.TotalCharges = df_raw.TotalCharges.astype('float')
         cleanup_length = len(df_raw)
         print('%s lines were deleted.' %(total_length - cleanup_length))
         # output into dummies
         churn_dummy_dict = {'Yes': 1, 'No': 0}
         df_raw.Churn.replace(churn_dummy_dict, inplace = True)
         # X y splitting
         y = df_raw.Churn.copy()
         X_raw = df_raw.drop(columns = 'Churn').copy()
         # input into dummies
         X_raw_types = dict(X_raw.dtypes)
         features = list(X_raw.columns)
         categorical_features = [feat for feat in features if X_raw_types[feat] ==_u
     '0'1
         categorical_features.remove('customerID')
         X = pd.get_dummies(X_raw, columns = categorical_features,prefix_sep=':')
         X = X.drop(columns = 'customerID').copy()
         return df_raw, categorical_features, X, y, cleanup_length
[3]: def df_preprocessing_knn():
```

```
[3]: def df_preprocessing_knn() :
    df_raw,categorical_features, X, y, _ = df_pp()
    # train val splitting

X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.20)
```

```
y_train = y_train.tolist()
y_test = y_test.tolist()

return df_raw,categorical_features, X, X_train, X_test, y, y_train, y_test
```

```
[4]: def df_preprocessing_rf(test_size) :
    df_raw,categorical_features, X, y, cleanup_length = df_pp()

# train val splitting

train = np.random.rand(cleanup_length)> test_size

X_train = X[train]
X_test = X[~train]

y_train = y[train].tolist()
y_test = y[~train].tolist()

return df_raw,categorical_features, X, X_train, X_test, y, y_train, y_test
```

1.1.3 2. K-Nearest Neighbors Classifier

We start by scaling the data.

```
[7]: df_raw,categorical_features, X, X_train, X_test, y, y_train, y_test = df_preprocessing_knn()
```

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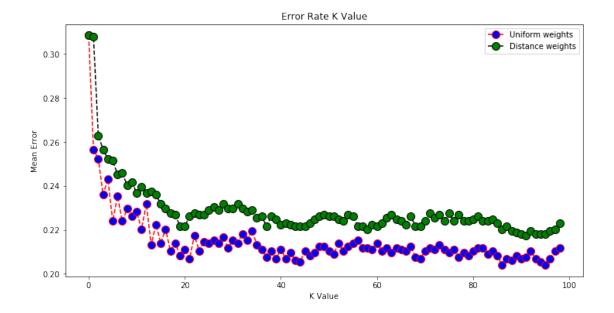
```
[8]: scaler = StandardScaler()
    scaler.fit(X_train)

X_train = scaler.transform(X_train)
    X_test = scaler.transform(X_test)
```

We run simulations for k-values between 1 and 100 to find the best fit parameters.

```
[9]: error_uni = np.array([0])
     error_dist = np.array([0])
     for i in range(1, 100):
         knn_uni = KNeighborsClassifier(n_neighbors=i, weights='uniform', n_jobs=-1)
         knn_uni.fit(X_train, y_train)
         pred_i_uni = knn_uni.predict(X_test)
         error_uni = np.append(error_uni, np.array(np.mean(pred_i_uni != y_test)))
         knn_dist = KNeighborsClassifier(n_neighbors=i, weights='distance',_
      \rightarrown jobs=-1)
         knn_dist.fit(X_train, y_train)
         pred_i_dist = knn_dist.predict(X_test)
         error_dist = np.append(error_dist, np.array(np.mean(pred_i_dist != y_test)))
     plt.figure(figsize=(12, 6))
     plt.plot(error_uni[1:], color='red', linestyle='dashed', marker='o',_
      →markerfacecolor='blue', markersize=10, label='Uniform weights')
     plt.plot(error_dist[1:], color='black', linestyle='dashed', marker='o', __
      →markerfacecolor='green', markersize=10, label='Distance weights')
     plt.title('Error Rate K Value')
     plt.xlabel('K Value')
     plt.ylabel('Mean Error')
     plt.legend(loc='best')
     plt.show()
     print('Minimum error rate with uniform weights: {:.2%} for k = {}'.
      →format(min(error_uni[1:]), np.argmin(error_uni[1:]) + 1))
     print('Minimum error rate with distance weights: {:.2%} for k = {}'.

→format(min(error dist[1:]), np.argmin(error dist[1:]) + 1))
```

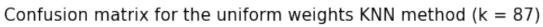


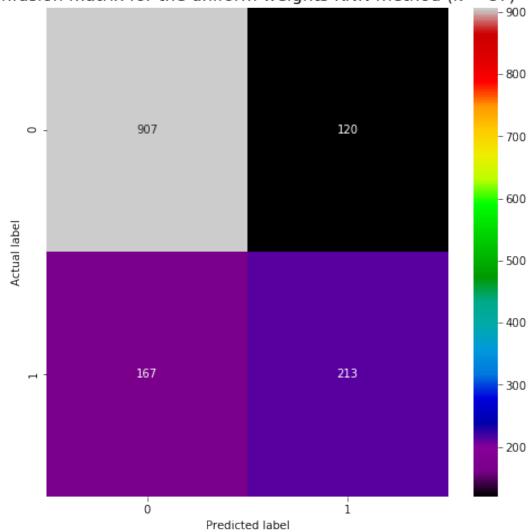
Minimum error rate with uniform weights: 20.40% for k = 87 Minimum error rate with distance weights: 21.75% for k = 92

```
UNIFORM WEIGHTS
Accuracy: 79.60%

precision recall f1-score support
```

0	0.84	0.88	0.86	1027
1	0.64	0.56	0.60	380
accuracy			0.80	1407
macro avg	0.74	0.72	0.73	1407
weighted avg	0.79	0.80	0.79	1407

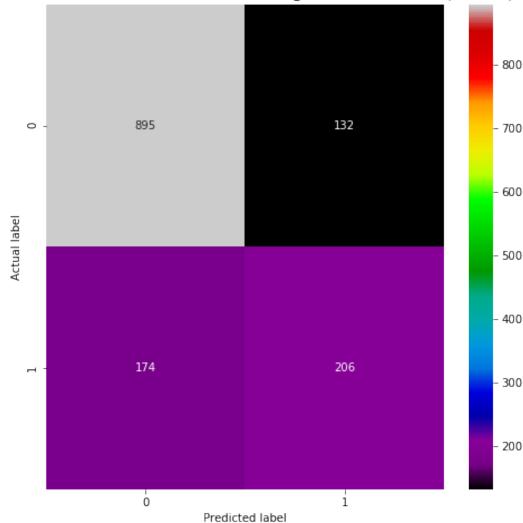




DISTANCE WEIGHTS Accuracy: 78.25%

	precision	recall	f1-score	support
0	0.84	0.87	0.85	1027
1	0.61	0.54	0.57	380
accuracy			0.78	1407
macro avg	0.72	0.71	0.71	1407
weighted avg	0.78	0.78	0.78	1407

Confusion matrix for the distance weights KNN method (k = 92)



1.1.4 3. Random Forest Classifier

We will compare two methods, which are grid search and random search.

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3.1 Grid search To avoid having too high a computational time, we will focus on 2 of the mot important parameters that are max depth and the number of estimators.

Max Depth This parameter is the depth of the trees, which is one of the most important. We range it between 4 (anything lower seems too low and imcreases computational time without much results) and 15.

Number of estimators This parameter is the number of trees that are going to be generated. Here the choice of the number of trees will mostly affect the computational time. Let's set the values between 10 and 500 and see the effects.

```
[15]: \max_{depth} = list(range(4,16))
[16]: n estimators = [10, 15, 20, 50, 100, 200, 500]
     Let's use the default 5 folds of cross validation.
[17]: grid_parameters = {'max_depth' : max_depth, 'n_estimators' : n_estimators }
[18]: grid_clf = GridSearchCV(clf, param_grid = grid_parameters,__
       →return_train_score=True, verbose = 3)
[19]: with parallel backend('ipyparallel'):
          grid_clf.fit(X_train, y_train);
     Fitting 5 folds for each of 84 candidates, totalling 420 fits
     [Parallel(n jobs=-1)]: Using backend IPythonParallelBackend with 8 concurrent
     workers.
     [Parallel(n_jobs=-1)]: Done 16 tasks
                                                 | elapsed:
                                                                5.2s
     [Parallel(n_jobs=-1)]: Done 112 tasks
                                                 | elapsed:
                                                               30.5s
     [Parallel(n_jobs=-1)]: Done 272 tasks
                                                 | elapsed: 1.3min
     [Parallel(n_jobs=-1)]: Done 420 out of 420 | elapsed:
                                                             2.3min finished
     Let's check which model is the best.
[20]: grid_best_score = grid_clf.best_score_
      grid_best_parameters = grid_clf.best_params_
      grid_best_max_depth = grid_best_parameters.get('max_depth')
      grid_best_n_estimators = grid_best_parameters.get('n_estimators')
      print('Grid search best_score: {:.5}'.format(grid_best_score))
      print('best_max_depth: {}'.format(grid_best_max_depth))
      print('best_n_estimators: {}'.format(grid_best_n_estimators))
     Grid search best_score: 0.80628
     best_max_depth: 7
     best_n_estimators: 20
[21]: |grid_clf_best = RandomForestClassifier(n_jobs = -1,max_depth = ___
       ⇒grid best max depth, n estimators = grid best n estimators )
```

```
[22]: grid_clf_best.fit(X_train, y_train);
```

Let's apply it to our validation set.

```
[23]: grid_y_pred = grid_clf_best.predict(X_test)
```

```
[24]: print("Accuracy: {:.2%}".format(accuracy_score(y_test, grid_y_pred)))
```

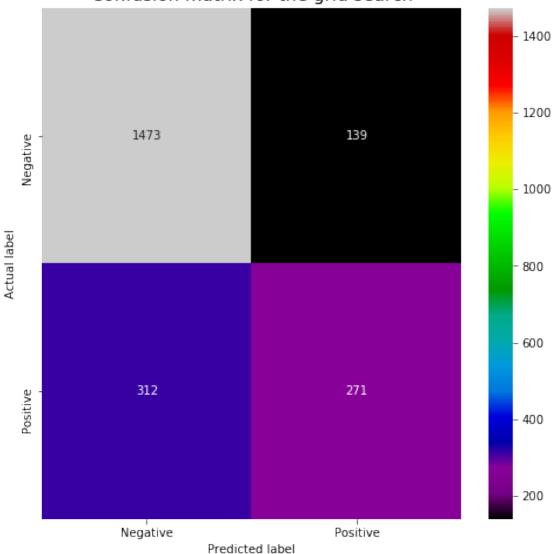
Accuracy: 79.45%

Given this accuracy, we can take a deeper look into the results.

```
[25]: grid_cm = confusion_matrix(y_test, grid_y_pred)
   index = ['Negative','Positive']
   columns = ['Negative','Positive']
   cm_df = pd.DataFrame(grid_cm,columns,index)

plt.figure(figsize=(8,8))
   sns.heatmap(cm_df, annot=True, fmt="d", cmap="nipy_spectral")
   plt.ylabel('Actual label');
   plt.xlabel('Predicted label');
   plt.xlabel('Predicted label');
   plt.title('Confusion matrix for the grid search', fontsize = 15);
```





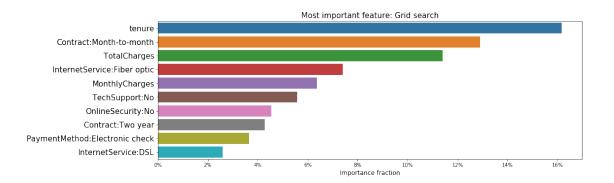
Feature importance

```
[26]: grid_feature_importances = [(list(X.columns)[i], grid_clf_best.

→feature_importances_[i]) for i in range(len(list(X.columns)))]

grid_feature_importances.sort(key=itemgetter(1), reverse = True)

plot_importance(grid_feature_importances, 10, 'Grid_search')
```



[27]: print(classification_report(y_test,grid_y_pred))

	precision	recall	f1-score	support
0	0.83	0.91	0.87	1612
1	0.66	0.46	0.55	583
accuracy			0.79	2195
macro avg	0.74	0.69	0.71	2195
weighted avg	0.78	0.79	0.78	2195

Feature selection Let's try to run the model again, but this time selecting only the most impacting features to save us some work and let's compare the results.

```
[28]: grid_selected_features = [grid_feature_importances[i][0] for i in range(15)] grid_X_train_sel = X_train[grid_selected_features] grid_X_test_sel = X_test[grid_selected_features]
```

```
[29]: with parallel_backend('ipyparallel'):
    grid_clf.fit(grid_X_train_sel, y_train);
```

Fitting 5 folds for each of 84 candidates, totalling 420 fits

 $[Parallel(n_jobs=-1)]$: Using backend IPythonParallelBackend with 8 concurrent workers.

[Parallel(n_jobs=-1)]: Done 420 out of 420 | elapsed: 2.2min finished

```
[30]: grid_best_score_sel = grid_clf.best_score_
    grid_best_parameters_sel = grid_clf.best_params_
    grid_best_max_depth_sel = grid_best_parameters.get('max_depth')
    grid_best_n_estimators_sel = grid_best_parameters.get('n_estimators')
```

Accuracy: 79.95%

3.2 Random Search After having explored a grid search, we can adopt another approach. Instead of searching for each value, let's give our model more parameters input, but instead let it choose randomly at each iteration one value for each parameter. It will then be evaluated again.

```
[36]: random_clf = RandomizedSearchCV(clf, param_distributions = random_parameters, 

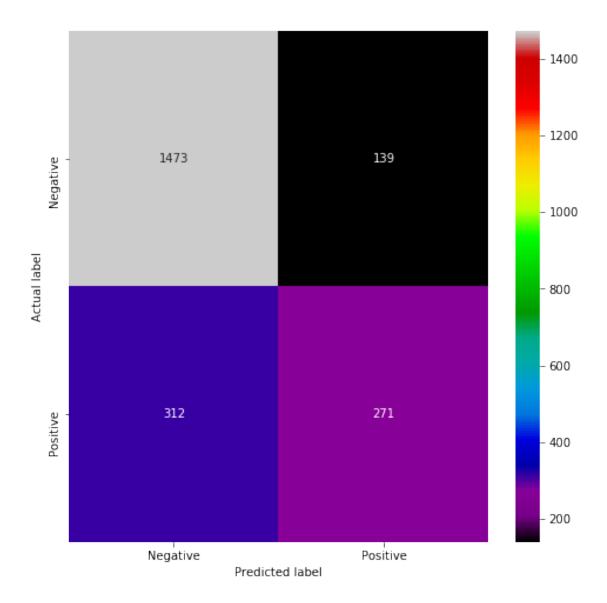
→n_iter = 20, verbose = 3)
```

```
[37]: with parallel_backend('ipyparallel'):
    random_clf.fit(X_train, y_train);
```

Fitting 5 folds for each of 20 candidates, totalling 100 fits

[Parallel($n_{jobs}=-1$)]: Using backend IPythonParallelBackend with 8 concurrent workers.

```
[Parallel(n_jobs=-1)]: Done 16 tasks | elapsed:
                                                             39.3s
     [Parallel(n_jobs=-1)]: Done 100 out of 100 | elapsed: 2.7min finished
[38]: random_best_score = random_clf.best_score_
      print('Random search best_score: {:.4}'.format(random_best_score))
      random_best_parameters = random_clf.best_params_
     Random search best_score: 0.8081
[39]: random_best_parameters
[39]: {'n_estimators': 410,
       'min_samples_leaf': 6,
       'max_features': 'sqrt',
       'max_depth': 45}
[40]: random_clf_best = RandomForestClassifier(max_depth = 35, max_features = 'sqrt', __
       →min_samples_leaf = 9, n_estimators = 250)
[41]: random_clf_best.fit(X_train, y_train);
[42]: random_y_pred = random_clf_best.predict(X_test)
[43]: random_cm = confusion_matrix(y_test, random_y_pred)
      annot_kws = {"ha": 'center', "va": 'center'}
      plt.figure(figsize=(8,8))
      sns.heatmap(cm_df, annot=True, fmt="d", cmap="nipy_spectral")
      plt.ylabel('Actual label');
      plt.xlabel('Predicted label');
```

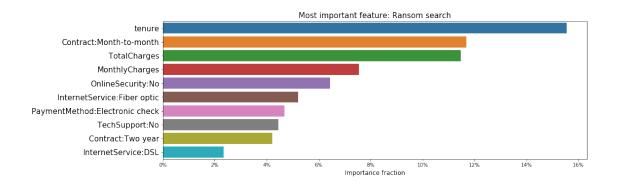


```
[44]: random_feature_importances = [(list(X.columns)[i], random_clf_best.

→feature_importances_[i]) for i in range(len(list(X.columns)))]

random_feature_importances.sort(key=itemgetter(1), reverse = True)

plot_importance(random_feature_importances, 10, 'Ransom search')
```



[45]: print(classification_report(y_test,random_y_pred))

	precision	recall	f1-score	support
0	0.84	0.90	0.87	1612
1	0.66	0.52	0.58	583
accuracy			0.80	2195
macro avg	0.75	0.71	0.73	2195
weighted avg	0.79	0.80	0.79	2195

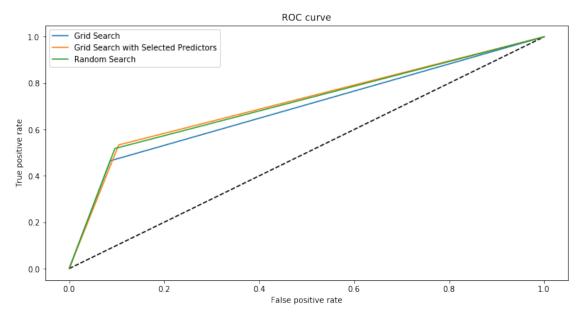
```
[46]: print("Accuracy: {:.2%}".format(accuracy_score(y_test, random_y_pred)))
```

Accuracy: 80.14%

3.3 Comparison

ROC Curve

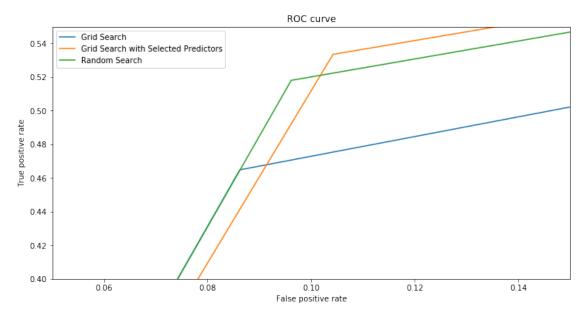
```
[47]: fpr_grid, tpr_grid, _ = roc_curve(y_test, grid_y_pred)
fpr_grid_sel, tpr_grid_sel, _ = roc_curve(y_test, grid_y_pred_sel)
fpr_random, tpr_random, _ = roc_curve(y_test, random_y_pred)
```



```
ROC AUC for Grid Search: 0.68930
ROC AUC for Grid Search with selected predictors: 0.71461
ROC AUC for Random Search: 0.71093
```

We observe that all three methods are better than a random prediction. The Grid Search with Selected Predictors has slightly better prediction than the Random Search, which has in turn slightly better prediction than the basic Grid Search.

```
plt.legend(loc='best')
plt.show()
```



Precision - Recall Curve

```
[50]: grid_precision, grid_recall, _ = precision_recall_curve(y_test, grid_y_pred)
grid_precision_sel, grid_recall_sel, _ = precision_recall_curve(y_test,

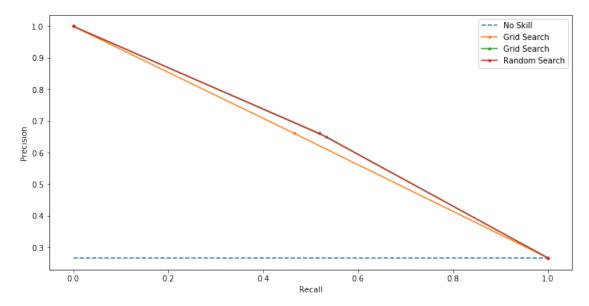
→grid_y_pred_sel)
random_precision, random_recall, _ = precision_recall_curve(y_test,

→random_y_pred)
```

```
[51]: y_test = np.array(y_test)
    no_skill = len(y_test[y_test == 1]) / len(y_test)

plt.figure(figsize=(12, 6))
    plt.plot([0, 1], [no_skill, no_skill], linestyle='--', label='No Skill')
    plt.plot(grid_recall, grid_precision, marker='.', label='Grid Search')
    plt.plot(grid_recall_sel, grid_precision_sel, marker='.', label='Grid Search')
    plt.plot(random_recall, random_precision, marker='.', label='Random Search')
    plt.xlabel('Recall')
    plt.ylabel('Precision')
    plt.legend(loc='best')
    plt.show()

pr_auc_grid = auc(grid_recall, grid_precision)
    print('Precision-Recall AUC for Grid Search: %.5f' % pr_auc_grid)
    pr_auc_grid_sel = auc(grid_recall_sel, grid_precision_sel)
```



Precision-Recall AUC for Grid Search: 0.63398

Precision-Recall AUC for Grid Search with selected predictors: 0.65332

Precision-Recall AUC for Random Search: 0.65343

1.1.5 4. Conclusion

We illustrated two main methods which are K Nearest Neighbors and Random Forest. Each of these methods can be used differently, given the different inputs we give them or the approach we want to have. In the long run, we see that all of them approach a 80% accuracy, with small differences among them.

Analyzing the feature importance can give the company some insights into what is affecting the rate of churn and how it could help them retain a higher number of customers.