Applied Machine Learning

HW2:Task 1

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Disclaimer: For a reason I have not been able to identify, the pipelines I created in this task tended to generate ConvergenceWarnings. I have opted to ignore these warnings as the results output by the functions were good.

```
In [1]: import pandas as pd
        import matplotlib.pyplot as plt
        import numpy as np
        from numpy.random import RandomState
        from sklearn.datasets import fetch_openml
        from sklearn.model_selection import train_test_split, GridSearchCV, KFold
        from sklearn.preprocessing import OneHotEncoder, StandardScaler
        from sklearn.linear model import LogisticRegression, Ridge
        from sklearn.compose import make column transformer
        from sklearn.pipeline import make pipeline, Pipeline
        from sklearn.model selection import cross val score
        from sklearn.svm import LinearSVC
        from sklearn.neighbors import KNeighborsClassifier
        import warnings
In [2]: #collect the dataset
        credit g dict = fetch openml("credit-g")
```

1.1 Categorical Features vs. Continuous Features

```
In [4]: print(credit_g_dict.feature_names)

['checking_status', 'duration', 'credit_history', 'purpose', 'credit_amount',
    'savings_status', 'employment', 'installment_commitment', 'personal_status', '
    other_parties', 'residence_since', 'property_magnitude', 'age', 'other_payment
    _plans', 'housing', 'existing_credits', 'job', 'num_dependents', 'own_telephon
    e', 'foreign_worker']
```

In total, we have 20 features (excluding the target). Some of the categorical features are detailed by the "categories" key of the credit_g dictionary. Upon further analysis, we observe that there are additional categorical variables: installment_commitment, residence_since ,existing_credits ,num_dependents.

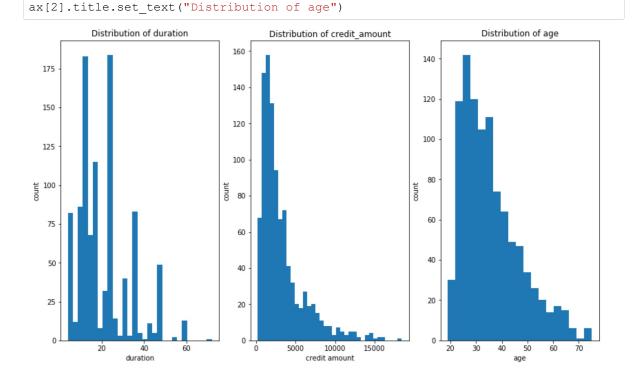
```
In [5]: print("The Categorical Features are : "+str(list(credit_g_dict["categories"].key
    s())+["installment_commitment", "residence_since" ,"existing_credits" ,"num_depe
    ndents"]))

The Categorical Features are : ['checking_status', 'credit_history', 'purpose
    ', 'savings_status', 'employment', 'personal_status', 'other_parties', 'proper
    ty_magnitude', 'other_payment_plans', 'housing', 'job', 'own_telephone', 'fore
    ign_worker', 'installment_commitment', 'residence_since', 'existing_credits',
    'num_dependents']
```

We can consider that the following features are continuous: duration, credit amount and age

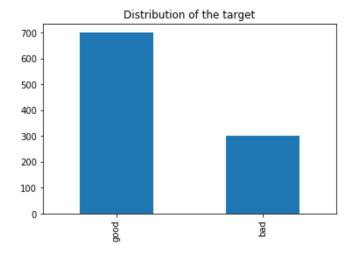
1.2 Distributions of Continuous Features and Target

```
In [6]: #create feature and target data frames
        df_X = pd.DataFrame(credit_g_dict.data,columns=credit_g_dict.feature_names)
        df_y = pd.DataFrame(credit_g_dict.target,columns=credit_g_dict.target_names)
In [7]: #distribution of the continuous variables
        fig, ax = plt.subplots(1,3,figsize=(14,8))
        ax[0].hist(df_X["duration"],bins="auto",)
        ax[1].hist(df_X["credit_amount"],bins="auto")
        ax[2].hist(df_X["age"],bins="auto")
        ax[0].set xlabel("duration")
        ax[1].set xlabel("credit amount")
        ax[2].set_xlabel("age")
        ax[0].set_ylabel("count")
        ax[1].set_ylabel("count")
        ax[2].set_ylabel("count")
        ax[0].title.set_text("Distribution of duration")
        ax[1].title.set text("Distribution of credit amount")
```



```
In [8]: #distribution of the target
    df_y['class'].value_counts().plot(kind='bar')
    plt.title("Distribution of the target")
```

```
Out[8]: Text(0.5, 1.0, 'Distribution of the target')
```



1.3 Preprocessing of the features

```
In [9]: #We encode the target to 0-1 values
         df y["IsGood"] = df y["class"].apply(lambda x: int(x=="good"))
         #We convert the categorical variables to objects
         #To do so we consider all columns other than age, duration and credit amount
         df = df X[df X.columns.difference(["age","duration","credit amount"])].astype("o
         bject")
         #We dummy encode the categorical variables and save them to df
         df = pd.get dummies(df)
         #We scale the continuous variables and append them to df
         scaler = StandardScaler()
         scaler.fit(df_X[["age","duration","credit_amount"]])
         df = pd.concat([df, pd.DataFrame(scaler.transform(df_X[["age","duration","credit
         amount"]]),columns=["age","duration","credit amount"])
         ], axis=1)
In [10]: np.shape(df)
Out[10]: (1000, 71)
```

Scaling the continuous variables and encoding the categorical variables now means our preprocessed dataframe, df, has 71 features.

Let us note that we also encoded the values of the target to be 1 if the class is "good" and 0 otherwise (stored in df_y["IsGood"]).

```
In [11]: #Initial Logistic Regression Model

X_train, X_test, y_train, y_test = train_test_split(df,df_y["IsGood"])

#We evaluate the model using a training/validation split
X_train_2, X_validate, y_train_2, y_validate = train_test_split(X_train,y_train)
model = LogisticRegression().fit(X_train_2,y_train_2)
model.score(X_validate,y_validate)
Out[11]: 0.7446808510638298
```

The logistic regression model without using pipelines yields 74.5% accuracy on the validation set.

1.4 Using Pipelines

We fit the first models without scaling the continuous features. We Dummy encode the categorical variables.

```
In [13]: | #We compare the classifiers without scaling the continuous features
         warnings.filterwarnings('ignore')
         df = df_X[df_X.columns.difference(["age","duration","credit_amount"])].astype("o
         bject")
         df = pd.concat([df,df X[["age","duration","credit amount"]]],axis=1)
         categorical = df.dtypes == object
         preprocess = make_column_transformer((OneHotEncoder(), categorical),("passthroug
        h",~categorical))
        X_train, X_test, y_train, y_test = train_test_split(df,df_y["IsGood"])
        model_log_reg = make_pipeline(preprocess, LogisticRegression())
        model_lin_SVC = make_pipeline(preprocess, LinearSVC())
        model KNN = make pipeline(preprocess, KNeighborsClassifier())
         scores log reg = np.mean(cross val score(model log reg, X train, y train))
         scores lin SVC = np.mean(cross val score(model lin SVC,X train,y train))
         scores KNN = np.mean(cross val score(model KNN, X train, y train))
        print("Logistic Regression with Pipeline Score : "+str(scores log reg))
        print("Linear SVC with Pipeline Score : "+str(scores lin SVC))
        print("K-Nearest Neighbors with Pipeline Score : "+str(scores KNN))
        Linear SVC with Pipeline Score: 0.6973333333333332
        K-Nearest Neighbors with Pipeline Score: 0.653333333333333333
```

In some cases, the returned scores are nan. This seems to indicate that we should scale our continuous variables if we want to obtain consistent results.

Let us see if the score on the dataset with scaled continuous features is better than for unscaled data(this would confirm our decision to scale).

We compare these results to those obtained when we scale the continuous features.

```
In [12]: warnings.filterwarnings('ignore')
         df = df_X[df_X.columns.difference(["age","duration","credit_amount"])].astype("o
         bject")
         df = pd.concat([df,df X[["age","duration","credit amount"]]],axis=1)
         categorical = df.dtypes == object
         preprocess = make column transformer((StandardScaler(), ~categorical),(OneHotEnc
         oder(), categorical))
         X train, X test, y train, y test = train test split(df,df y["IsGood"])
         X train 2, X validate, y train 2, y validate = train test split(X train, y train)
         model log reg = make pipeline(preprocess, LogisticRegression())
         model lin SVC = make pipeline(preprocess, LinearSVC())
         model KNN = make pipeline(preprocess, KNeighborsClassifier())
         scores log reg = np.mean(cross_val_score(model_log_reg, X_train, y_train))
         scores lin SVC = np.mean(cross val score(model lin SVC, X train, y train))
         scores KNN = np.mean(cross val score(model KNN, X train, y train))
         print("Logistic Regression with Pipeline Score : "+str(scores log reg))
         print("Linear SVC with Pipeline Score : "+str(scores lin SVC))
         print("K-Nearest Neighbors with Pipeline Score : "+str(scores KNN))
         Logistic Regression with Pipeline Score: 0.757333333333333334
         Linear SVC with Pipeline Score: 0.756
         K-Nearest Neighbors with Pipeline Score : 0.74
```

We observe that in general, scaling the continuous variables improves accuracy (a few percentage points improvement). From now on, we will consider the data where the continuous variables are scaled.

1.5 Parameter Tuning

```
In [14]: warnings.filterwarnings('ignore')
         pipe = Pipeline([("regressor", LogisticRegression())])
         param_grid = [{'regressor': [LogisticRegression()],
                        'regressor__C':np.logspace(-3,3,10)},
                       {'regressor': [LinearSVC()],
                        'regressor C':np.logspace(-3,3,10)},
                       {'regressor': [KNeighborsClassifier()],
                       'regressor__n_neighbors':range(1,10)
                      } ]
         grid = GridSearchCV(pipe, param grid,cv=5)
         grid.fit(X train,y train)
         print("Best model based on training : "+ str(grid.best params ))
         print("Score of best model on training : "+str(grid.score(X train,y train)))
         Best model based on training : {'regressor': LogisticRegression(C=2.1544346900
         31882, class weight=None, dual=False,
                            fit_intercept=True, intercept_scaling=1, l1_ratio=None,
                            max iter=100, multi class='auto', n jobs=None, penalty='12
                            random state=None, solver='lbfgs', tol=0.0001, verbose=0,
                            warm start=False), 'regressor C': 2.154434690031882}
         Score of best model on training: 0.7946666666666666
```

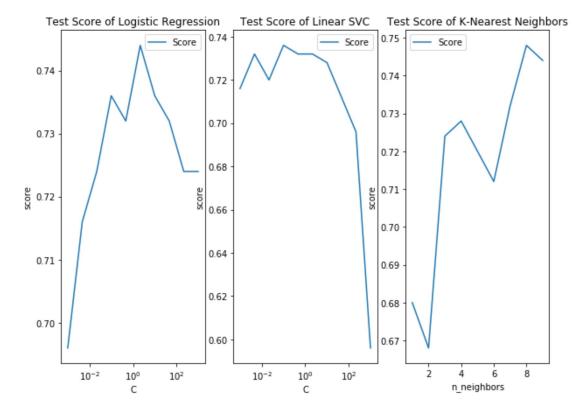
Parameter Tuning improves results by a few percentage points. The model which best performs on training data is sometimes Logistic Regression and sometimes LinearSVC. This seems to indicate that both methods reach similar optima. In this run, it was a Logistic Regression model with C=2.15. Let us see how this model performs on the test set.

```
In [15]: print("Test score : "+ str(grid.score(X_test,y_test)))
Test score : 0.744
```

The test score is similar to the validation scores of the previous models. This seems to indicate that the model has not overfitted.

```
In [16]: | #Performance as function of Parameters for Logistic Regression, LinearSVC and KN
         warnings.filterwarnings('ignore')
         C = np.logspace(-3,3,10)
         n neighbors = range(1,10)
         LR test score grid = [LogisticRegression(C=C[i]).fit(X train,y train).score(X te
         st,y test) for i in range(len(C))]
         SVC_test_score_grid = [LinearSVC(C=C[i]).fit(X_train,y_train).score(X_test,y_tes
         t) for i in range(len(C))]
         KNN test score grid = [KNeighborsClassifier(n neighbors=n neighbors[i]).fit(X tr
         ain,y_train).score(X_test,y_test) for i in range(len(n_neighbors))]
         fig, ax = plt.subplots(1,3,figsize=(10,7))
         ax[0].plot(C,LR test score grid,label="Score")
         ax[0].set xscale("log")
         ax[0].set xlabel("C")
         ax[0].set_ylabel("score")
         ax[0].title.set text("Test Score of Logistic Regression")
         ax[0].legend()
         ax[1].plot(C,SVC test score grid,label="Score")
         ax[1].set xscale("log")
         ax[1].set xlabel("C")
         ax[1].set ylabel("score")
         ax[1].title.set text("Test Score of Linear SVC")
         ax[2].plot(n_neighbors,KNN_test_score_grid,label="Score")
         ax[2].set_xlabel("n_neighbors")
         ax[2].set_ylabel("score")
         ax[2].title.set_text("Test Score of K-Nearest Neighbors")
         ax[2].legend()
```

Out[16]: <matplotlib.legend.Legend at 0x280742a9828>



The performance graphs confirm that Logistic Regression models are the most reliable of the three classification methods, as the test scores are consistently above 70% even when the parameters change. LinearSVC tends to have worse results for large values of C. K-Nearest Neighbors results are lower in general than for the other two methods.

1.6 Cross Validation without Stratification and with Shuffling

```
In [17]: # KFold Cross-validation with Shuffling
         warnings.filterwarnings('ignore')
         cv = KFold(5, shuffle=True)
         pipe = Pipeline([('regressor', LogisticRegression())])
         param grid = [{'regressor': [LogisticRegression()],
                        'regressor C':np.logspace(-3,3,10)},
                       {'regressor': [LinearSVC()],
                        'regressor__C':np.logspace(-3,3,10)},
                       { 'regressor': [KNeighborsClassifier()],
                       'regressor n neighbors':range(1,10)
         grid = GridSearchCV(pipe, param grid, cv=cv,)
         grid.fit(X train, y train)
         print("Best model based on training: "+ str(grid.best params))
         print("Score of best model on test : "+str(grid.score(X test,y test)))
         Best model based on training : {'regressor': LogisticRegression(C=0.4641588833
         6127775, class weight=None, dual=False,
                            fit intercept=True, intercept scaling=1, l1 ratio=None,
                            max iter=100, multi class='auto', n jobs=None, penalty='12
                            random_state=None, solver='lbfgs', tol=0.0001, verbose=0,
                            warm_start=False), 'regressor__C': 0.46415888336127775}
         Score of best model on test : 0.732
```

When we use shuffling and KFold, the best model on this run was LogisticRegression, but the value of the parameter found changed.

```
In [18]: # KFold Cross-Validation with random seed Shuffling
         warnings.filterwarnings('ignore')
         cv = KFold(5, shuffle=True, random_state=RandomState(1))
         pipe = Pipeline([('regressor', LogisticRegression())])
         param grid = [{'regressor': [LogisticRegression()],
                        'regressor C':np.logspace(-3,3,10)},
                       { 'regressor': [LinearSVC()],
                        'regressor__C':np.logspace(-3,3,10)},
                       { 'regressor': [KNeighborsClassifier()],
                       'regressor n neighbors':range(1,10)
                      } ]
         grid = GridSearchCV(pipe, param_grid,cv=cv,)
         grid.fit(X train, y train)
         print("Best model based on training : "+ str(grid.best params ))
         print("Score of best model on test : "+str(grid.score(X test,y test)))
         Best model based on training: {'regressor': LinearSVC(C=0.021544346900318832,
         class weight=None, dual=True,
                   fit intercept=True, intercept scaling=1, loss='squared hinge',
                   max iter=1000, multi class='ovr', penalty='12', random state=None,
                   tol=0.0001, verbose=0), 'regressor C': 0.021544346900318832}
         Score of best model on test : 0.72
```

Changing the Random Seed has once again changed the optimal estimator. The best estimator is now a LinearSVC model with C=0.022(results might change if you run the code again).

```
In [19]: | # KFold Cross-Validation with Shuffling, random seed and random state train test
         split
         warnings.filterwarnings('ignore')
         X_train, X_test, y_train, y_test = train_test_split(df,df_y["IsGood"],random_sta
         te=RandomState(1))
         \verb|#We convert X_train and X_test from sparse matrix to array|\\
         X train = X train.toarray()
         X_test = X_test.toarray()
         cv = KFold(5, shuffle=True, random state=RandomState(1))
         pipe = Pipeline([('regressor', LogisticRegression())])
         param grid = [{'regressor': [LogisticRegression()],
                         'regressor C':np.logspace(-3,3,10)},
                        {'regressor': [LinearSVC()],
                         'regressor__C':np.logspace(-3,3,10)},
                       { 'regressor': [KNeighborsClassifier()],
                       'regressor__n_neighbors':range(1,10)
                       } ]
         grid = GridSearchCV(pipe, param grid,cv=cv,)
         grid.fit(X train, y train)
         print("Best model based on training : "+ str(grid.best_params_))
         print("Score of best model on test : "+str(grid.score(X_test,y_test)))
         Best model based on training : {'regressor': LinearSVC(C=0.021544346900318832,
         class weight=None, dual=True,
                   fit_intercept=True, intercept_scaling=1, loss='squared_hinge',
                   max_iter=1000, multi_class='ovr', penalty='12', random_state=None,
                   tol=0.0001, verbose=0), 'regressor C': 0.021544346900318832}
         Score of best model on test : 0.76
```

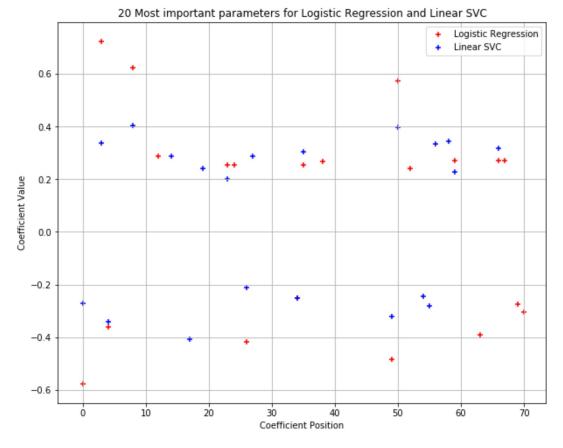
Changing the random state of the split into training and test and repeating this process for the train/validation split did not change the value of the parameter or the estimator in this instance. The best model is still a LinearSVC model. It is worth noting that for some runs, the best estimator is a LogisticRegression, whose parameters also change when we introduce shuffling, KFold and RandomStates.

1.7 Visualizing main coefficients

```
In [20]: | #We first select satisfactory parameters for LogisticRegression
         warnings.filterwarnings('ignore')
         pipe = Pipeline([('regressor', LogisticRegression())])
         param grid = [{'regressor': [LogisticRegression()],
                         'regressor__C':np.logspace(-3,3,10)}]
         grid = GridSearchCV(pipe, param grid)
         grid.fit(X_train,y_train)
         model_LR = LogisticRegression(C = grid.best_params_["regressor__C"]).fit(X_trai
         n,y_train)
         #We select satisfactory parameters for LinearSVC
         pipe = Pipeline([('regressor', LinearSVC())])
         param_grid = [{'regressor': [LinearSVC()],
                         'regressor__C':np.logspace(-3,3,10)}]
         grid = GridSearchCV(pipe, param_grid)
         grid.fit(X_train,y_train)
         model_SVC = LinearSVC(C = grid.best_params_["regressor__C"]).fit(X_train,y_trai
         n)
```

```
In [21]: \mid #We select the 20 most important coefficients (i.e those with largest absolute v
         alue) for Logistic Regression and LinearSVC
         coef_LR = np.absolute(model_LR.coef_)
         coef LR = np.argsort(coef LR) #sorts from smallest to largest
         coef LR = np.flip(coef LR)[0][:20] #selects 20 most important coefs
         coef SVC = np.absolute(model SVC.coef )
         coef_SVC = np.argsort(coef_SVC)
         coef SVC = np.flip(coef SVC)[0][:20]
         plt.figure(figsize=(10,8))
         plt.scatter(coef_LR,[model_LR.coef_[0][i] for i in coef_LR],marker="+",c="r",lab
         el="Logistic Regression")
         plt.scatter(coef_SVC,[model_SVC.coef_[0][i] for i in coef_SVC],marker="+",c="b",
         label="Linear SVC")
         plt.xlabel("Coefficient Position")
         plt.ylabel("Coefficient Value")
         plt.title("20 Most important parameters for Logistic Regression and Linear SVC")
         plt.grid()
         plt.legend(loc="upper right")
```

Out[21]: <matplotlib.legend.Legend at 0x280740128d0>



```
In [22]: print(np.sort(coef_LR))
    print(np.sort(coef_SVC))

[ 0  3  4  8 12 23 24 26 34 35 38 49 50 52 59 63 66 67 69 70]
    [ 0  3  4  8 14 17 19 23 26 27 34 35 49 50 54 55 56 58 59 66]
```

The 20 most important parameters for Linear Regression and for Linear SVC are almost identical (even though their magnitudes are different).