Machine Learning project Oil price prediction

Bogdana Kolic, Clemence Mottez, Matheo Le Masson

Motivations

- Crucial for industries and decision-makers
 - Impact on the global economy, investment and trading strategies, supply and demand dynamics, risk management and market forecasting
- Usual data: year and price
 - More comprehensive approach with 23 features
- Feature selection
 - determine what contribute to oil prices
 - deepen our understanding of the complex dynamics driving oil prices
 - provides us with enhanced predictive capabilities

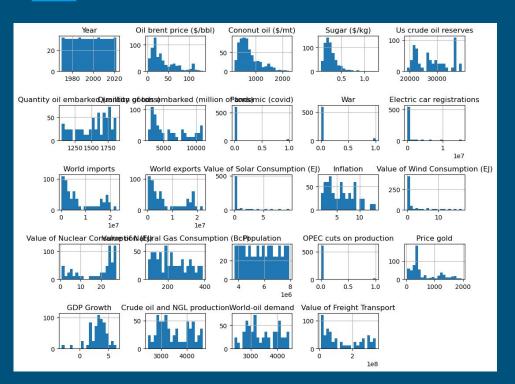
Data

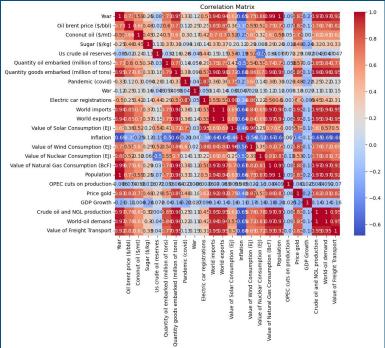
- How did we create our data set
 - Decided by ourselves what we thought could influence the price, asked domain experts
 - Sources: macrotrends.net, unctadstat.unctad.org, datasource.kapsarc.org, data.worldbank.org, tradingeconomics.com
 - Handle missing values
 - Normalization

$$x = \frac{x - x_{min}}{x_{max} - x_{min}}$$

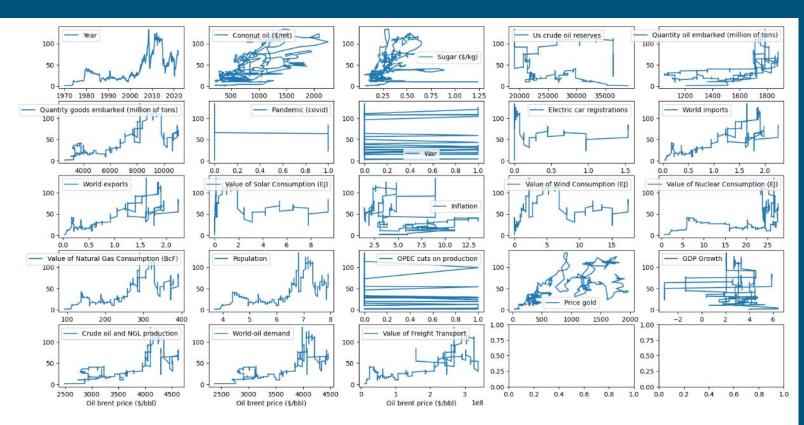
- Our dataset
 - 23 parameters, total of 24 columns
 - Monthly data from 1970 to 2022, total of 624 rows

Data visualization





Data visualization



Data preprocessing

- Removing instances with missing values
- Allow for the option of averaging the data over several months
 - 24th feature average price of oil in the past months?
- Manual feature selection
- Normalization
- Splitting the data into training and test sets

Manual feature selection and averaging

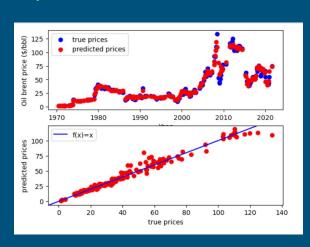
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In [6]: # discovered function rolling() at:
        # https://stackoverflow.com/questions/60274621/calculate-average-of-specified-range-of-values-in-pandas-column-and-store-as-ano
        # and function shift() at:
        # https://stackoverflow.com/questions/10982089/how-to-shift-a-column-in-pandas-dataframe
        # features that shouldn't ever be averaged
        no avg features = ['Year', 'War', 'OPEC cuts on production', 'Pandemic (covid)']
        # features we choose to use
        features = ['Year', 'Cononut oil ($/mt)', 'Sugar ($/kg)', 'Us crude oil reserves',
                     'Quantity oil embarked (million of tons)', 'Quantity goods embarked (million of tons)', 'Pandemic (covid)', 'War',
                    'Electric car registrations', 'World imports', 'World exports', 'Value of Solar Consumption (EJ)', 'Inflation',
                    'Value of Wind Consumption (EJ)', 'Value of Nuclear Consumption (EJ)', 'Value of Natural Gas Consumption (BCF)',
                    'Population', 'OPEC cuts on production', 'Price gold', 'GDP Growth', 'Crude oil and NGL production',
                    'World-oil demand', 'Value of Freight Transport']
        df = data.loc[:, features]
        num_months_avg = 1 # over how many months to take average
        take average = 0 # 1 or 0, indicates whether we do the averaging
        for feature in features:
            if feature not in no avg features:
                df[feature] = df[feature].rolling(num_months_avg).mean() # average
            elif num months avg > 1:
                df[feature] = df[feature].shift(-1) # if taking average, adjust the other features for forecasting
        df.dropna(inplace=True)
        df.describe()
```

Models

- Which models did we study?
 - Linear and Polynomial regression, decision trees, Random forest
- Hyperparameters
 - Ridge and Lasso penalty for Linear and Polynomial regression
 - Number of estimators, maximal depth, ... for decision trees
- Which one did we choose for feature selection.
 - xgbRegressor + random forest
 - The decision trees seem to work the best, different scores based on different metrics and features

Training

- Scoring and error metrics
 - Minimizing the *Mean squared error*
 - Maximizing R2 score
 - Minimizing the Cross-validation with mean squared error
 - Over the whole dataset
- Visualizing the results
 - two plots
 - ex: bagging regressor performance, all features



Hyperparameter tuning

- Initialize several options and test all the possibilities
- Evaluation done by cross-validation on the training set
- example on AdaBoost model:

AdaBoost

```
In [27]: # tuning the hyper-parameters:
         depths = [1, 2, 3, 4, 5, 10, 20, 30, 40, 50]
         estimators = [1, 2, 3, 4, 5, 10, 20, 30, 40, 50]
         min mse = math.inf
         max depth = 0
         num estimators = 0
         for d in depths:
              for e in estimators:
                     for sf in samples features:
                         adaboost model = AdaBoostRegressor(DecisionTreeRegressor(max depth=d),n estimators=e)
                         adaboost model.fit(x train, y train)
                         prediction = adaboost model.predict(x test)
                         mse = mean squared error(prediction, y test)
                         if mse<min mse:
                             min mse = mse
                             max depth = d
                             num estimators = e
                             max samples features = sf
         adaboost model = AdaBoostRegressor(DecisionTreeRegressor(max depth),n estimators=num estimators)
         adaboost model.fit(x train, v train)
         v predict adaboost = adaboost model.predict(x test)
         adaboost r2 score = r2 score(y test, y predict adaboost)
         adaboost_mse = mean_squared_error(y_test, y_predict_adaboost)
         adaboost cv = cross validate(adaboost model, df, oil prices, 4, mean squared error, np.mean, False)
         print("max depth: ", max depth)
         print("num estimators: ", num estimators)
         print("mse: ", adaboost mse)
         print("r2 score: ", adaboost_r2_score)
         print("cv: ", adaboost cv)
         model results["adaboost"] = (adaboost r2 score, adaboost mse, adaboost cv)
```

Interpretation of the results

- By sorting the final scores/ errors
 - example on the left: raw features
 - example on the right: with additional average oil price feature and averaging

Summing up the results In [35]: comparison r2 = sorted(model results.items(), key = lambda x:x[1], reverse = True) [('adaboost', (0.9829371845001577, 15.492112234042551, 713.3105756410255)), ('random forest', (0.9775000018764277, 20.428779541 06389, 698.034808000225)), ('polynomial_ridge', (0.9762947996355764, 21.52303789368814, 6762.762089856493)), ('bagging', (0.975 035427040476, 22.666479993730352, 640.6171660665955)), ('gradient_boosting', (0.9715275884435935, 25.851407430958524, 620.13537 79207891)), ('xgb_regressor', (0.9698048834023651, 27.415530294802856, 621.124659375191)), ('polynomial_lasso', (0.967417781600 923, 29.562889428608727, 2343.74914969669)), ('linear regression', (0.9361874946807662, 57.938298304294435, 1308536400, 38078 8)), ('linear ridge', (0.9359970842441161, 58.11118065682978, 67234304.81083053)), ('linear lasso', (0.9192127144482721, 73.350 47928418226, 2546211.101777478)), ('polynomial', (0.894844984516612, 95.47505813774859, 4.681250740188116e+23))] In [36]: comparison mse = sorted(model results.items(), key = lambda x:(x[1][1], x[1][0], x[1][2])) print(comparison mse) [('adaboost', (0.9829371845001577, 15.492112234042551, 713.3105756410255)), ('random forest', (0.9775000018764277, 20.428779541 . 66389, 698.034808000225)), ('polynomial_ridge', (0.9762947996355764, 21.52303789368814, 6762.762089856493)), ('bagging', (0.975 035427040476, 22.666479993730352, 640.6171660665955)), ('gradient_boosting', (0.9715275884435935, 25.851407430958524, 620.13537 79207891)), ('xgb_regressor', (0.9698048834023651, 27.415530294802856, 621.124659375191)), ('polynomial_lasso', (0.967417781600 923, 29.582889428608727, 2343.74914969669)), ('linear_regression', (0.9361874946807662, 57.338298304294435, 1308536400.38078 8)), ('linear ridge', (0, 9359970842441161, 58, 11118666682978, 67234304, 81083053)), ('linear lasso', (0, 9192127144482721, 73, 350 47928418226, 2546211.101777478)), ('polynomial', (0.894844984516612, 95.47505813774859, 4.681250740188116e+23))] In [37]: comparison_cv = sorted(model_results.items(), key = lambda x:(x[1][2], x[1][0], x[1][1])) print(comparison cv) [('gradient boosting', (0.9715275884435935, 25.851407430958524, 620.1353779207891)), ('xgb regressor', (0.9698048834023651, 27. 415530294802856, 621.124659375191)), ('bagging', (0.975035427040476, 22.666479993730352, 640.6171660665955)), ('random forest', (0.9775000018764277, 20.42877954106389, 698.034808000225)), ('adaboost', (0.9829371845001577, 15.492112234042551, 713.310575641 0255)), ('polynomial_lasso', (0.967417781600923, 29.582889428608727, 2343.74914969669)), ('polynomial_ridge', (0.97629479963557 64, 21.52303789368814, 6762,762089856493)), ('linear lasso', (0.9192127144482721, 73.35047928418226, 2546211,101777478)), ('linear lasso', (0.9192127144482721, 73.35047928418226, 2546211,101777478)), ear_ridge', (0.9359970842441161, 58.11118066682978, 67234304.81083053)), ('linear_regression', (0.9361874946807662, 57.93829830 4294435, 1308536400.380788)), ('polynomial', (0.894844984516612, 95.47505813774859, 4.681250740188116e+23))]

- Conclusion:

- similar results that change with each run
- decision trees perform well
- our choice is the XGBRegressor

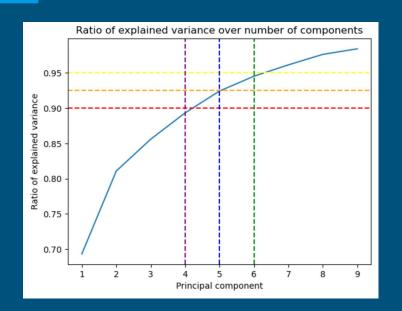
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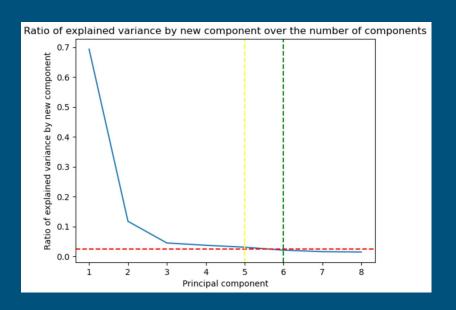
Feature selection

Why?

- Data 624x24!!!
 - Reduce computational cost
- Noise, redundant, irrelevant information have negative impact
 - Improve the performance of the model
- Capture essential dynamics of the oil market
 - Facilitate interpretation
 - What really predicts the price of oil?

PCA as an answer to: How many features should we keep?





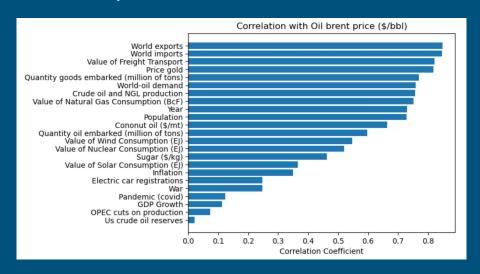
5 components explain 90% of variance Components after the 5th explain less than 2.5% of variation

Filter models

- Computationally efficient
- Involve statistical measures such as correlation

- Measures the linear relationship between each feature and the

target variable



Embedded models

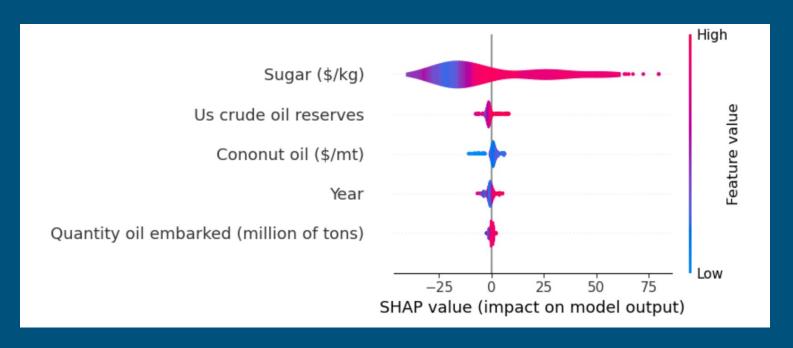
- Incorporates feature selection within the model training process
 - Optimizes both the model's performance and feature selection simultaneously
- Select relevant features based on their contribution to the model's accuracy

SHAP

- Interpretability tool
- Gain insights into feature importance and guide the feature selection process
- Show global contribution
 - Computed for each feature and used to rank the importance of features
- Show local feature contribution
 - for each instance

Embedded models

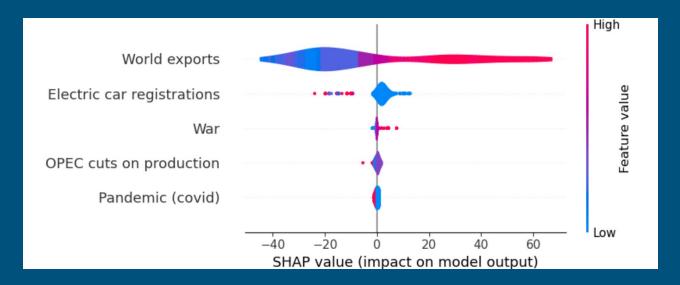
XGboost + SHAP



- Iteratively select and evaluate different subsets of features to find the subset that yields the best performance.
- Computationally expensive but accurate results
- Used Random Forest and XGB models

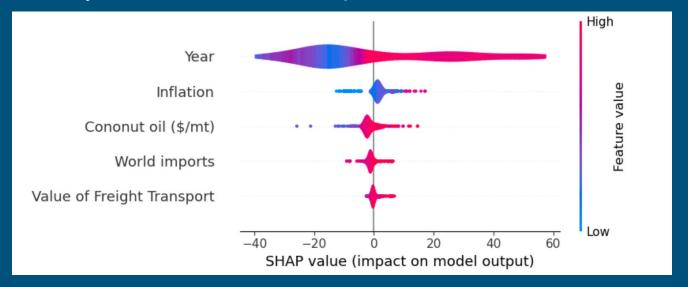
Sequential Forward Selection

- Starts with an empty set of features
- Iteratively adds features based on the best-performing subset



Recursive Feature Elimination (RFE)

- Starts with all features
- iteratively eliminates the least important feature



Boruta

- Compared importance of features with random shadow features
 - A feature is important if it can do better than the best randomized feature
- Used eBoruta (extension of Boruta that already uses the SHAP importance and that is mode agnostic)

Feature	Importance	
Year	20.353813	
Inflation	2.827759	
Cononut oil (\$/mt)	2.487869	
World imports	1.869873	
Sugar (\$/kg)	1.062382	

Selection

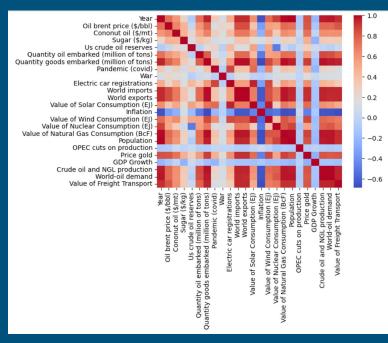
- Year
- World imports
- World exports
- Inflation
- Price of Gold
- War
- OPEC cuts on production

Correlation

- Small selection so we don't want too correlated variables
- Threshold at 0.95

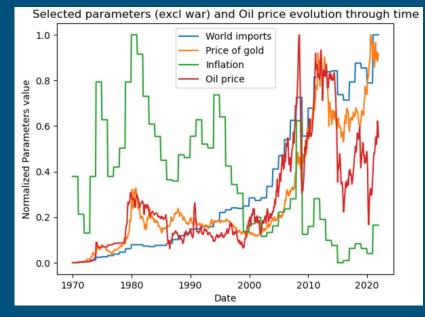
Result:

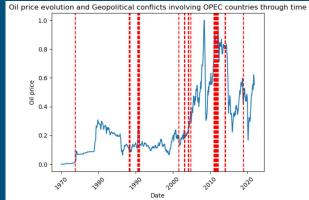
World export and world imports



Final selection

- Year
- World imports
- World exports <- too correlated
- Inflation
- Price of Gold
- War
- OPEC cuts on production <- only 5 features selected





MSE

Data / Model	XGB Regressor	Decision Tree	Random Forest
Raw Data	621	643	680
Data with PCA	253	390	318
Selected Data	372	474	555

Limits - What could do after?

- Limits:

 don't know if causation or correlation for example price of gold certainly a correlation

- After:

- Network model
- Future analysis and predictions
 make predictions for future oil prices
 Incorporate new data as it becomes available
- How to optimally compare models? With data modifications
- Optimize algorithm we used (for example SHAP is computationally very expensive)
- New model: RNNs with LSTM to avoid Vanishing Gradient problem

Network model

- Much better accuracy
 - MSE from cross-validation around 220 (compared to XGB around 400)
- Relatively simple:
 - 3 Hidden Dense layers with Relu activation
 - Dropout regularization to avoid overfitting

Other info for questions

- slides to explain the different models we used
- slides with term explanations, ...

SHAP values

