Workshop: Introduction to Interpretability in ML

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Materials available at https://github.com/azsom/ODSC-East-2022

About me

- Born and raised in Hungary
- Astrophysics PhD at MPIA, Heidelberg, Germany
- Postdoctoral researcher at MIT (still in astrophysics at the time)
- Started at Brown in 2015 as a Data Scientist in CCV (Center for Computation and Visualization)
- Lead Data Scientist since 2017 at CCV
- Assistant Professor since summer of 2021 at the DSI

Data Science at AI+Training

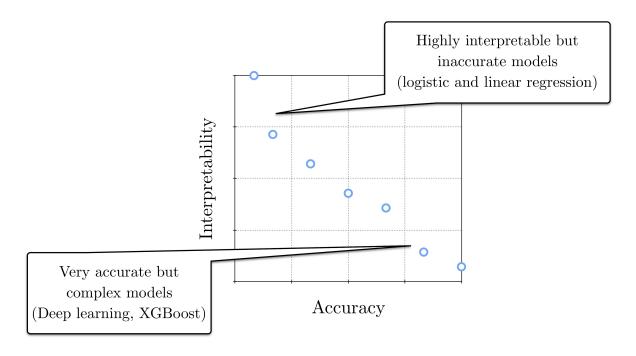
- Supervised Machine Learning Course Series https://app.aiplus.training/courses/supervised-machine-learning-series
- 6 courses that walk through the main steps of developing an ML pipeline
- github repo available here: https://github.com/azsom/Supervised-Learning
- Week 7 of the ODSC ML Certification https://aiplus.training/certificates/
- this workshop is based on course 6

Introduction to Interpretability in ML

By the end of this workshop, you will be able to

- calculate global explanations using the coefficients of linear models
- perform permutation importance using any ML model
- use the SHAP package to calculate local explanations

Motivation: simple vs. complex models



- local feature importance improves the interpretability of complex models
- check out this page for a good example

Motivation: build trust in your model

- · debugging ML models is tough
 - a model that runs without errors/warning is not necessarily correct
- model seems accurate but does it make predictions the way you expect it to be?
 - wolf vs. husky classifier
- model inspection is important
 - feature importance metrics measure useful they are at predicting the target variable
 - global: one score per feature, a vector of shape (1, n_features)
 - o general overview over a larger sample size
 - local: one score per feature and point, an array of shape (n_samples, n_features)
 - explanation specific to each sample
- · discuss the interpretability results with a subject matter expert
 - if they say something is surprising, you either
 - have a bug in your code (more likely),
 - you discovered something new which is exciting!

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Math recap

Linear regression:

$$f(X_i) = y_i' = heta_0 + X_{i1} heta_1 + X_{i2} heta_2 + ... = heta_0 + \sum_{j=1}^m heta_j X_{ij}$$
 ,

Logistic regression:

$$f(X_i)=y_i'=rac{1}{1+e^{-z}}$$
 ,

$$z = heta_0 + \sum_{j=1}^m heta_j x_{ij}$$

Here, y_i' is the prediction of the model for sample i using the X_i feature vector, and θ are coefficients.

What to do if you want to use the coefficients of your linear model as explanations

- Look at the absolute value of the coefficients!
 - a feature is important, if its coefficient's magnitude is large
 - doesn't matter if the coefficient is positive or negative
- Standardize ALL features!
 - most practitioners standarize the continuous features
 - but practicioners often forget to standarize the one-hot encoded or ordinal encoded features which is a mistake!
 - all features need to have the same mean and standard deviation if you want to use the coefficients to measure importance!

Simple ML pipeline

- we will use the diabetes dataset in this workshop
- https://www4.stat.ncsu.edu/~boos/var.select/diabetes.html
- "Ten baseline variables, age, sex, body mass index, average blood pressure, and six blood serum measurements were obtained for each of n = 442 diabetes patients, as well as the response of interest, a quantitative measure of disease progression one year after baseline."

```
df = pd.read csv('data/diabetes.csv')
        df.head()
        y = df['Y']
        X = df.loc[:, df.columns != 'Y']
        print(y.head())
        print(X.head())
        0
            151
        1
             75
        2
            141
        3
            206
            135
        Name: Y, dtype: int64
          AGE
                 SEX BMI BP S1
                                          S2 S3 S4 S5 S6
           59 male 32.1 101.0 157 93.2 38.0 4.0 4.8598 87
        0
           48 female 21.6 87.0 183 103.2 70.0 3.0 3.8918
        1
                                                                  69
        2
          72 male 30.5 93.0 156 93.6 41.0 4.0 4.6728 85
        3
          24 female 25.3 84.0 198 131.4 40.0 5.0 4.8903 89
           50 female 23.0 101.0 192 125.4 52.0 4.0 4.2905 80
In [2]: # our ML pipeline
        def MLpipe_KFold_RMSE(X, y, preprocessor, ML_algo, param_grid,random_state):
            This function splits the data to other/test (80/20) and then applies KFold
            The RMSE is minimized in cross-validation.
            1.1.1
            # split data to other/test 80/20, and the use KFold with 4 folds
            X_other, X_test, y_other, y_test = train_test_split(X, y, test_size=0.2, re
            kf = KFold(n splits=4, shuffle=True, random state=random state)
            # the sklearn pipeline
            if preprocessor == None:
                # no preprocessing
                pipe = Pipeline(steps=[('regressor', ML algo)])
            else:
               pipe = Pipeline(steps=[('preprocessor', preprocessor),
                                      ('regressor', ML_algo)])
            # loop through the hyperparameter combinations or use GridSearchCV
            grid = GridSearchCV(pipe, param grid=param grid,scoring='r2',
                               cv=kf, return_train_score=True)
            # for each combination, calculate the train and validation scores using the
            grid.fit(X other, y other)
            # find which hyperparameter combination gives the best validation score
            print(grid.best params )
            # calculate the test score
           test score = grid.score(X test,y test)
            print(test_score,r2_score(y_test,grid.predict(X_test)))
            return grid best_estimator_, test_score, (X_other, y_other), (X_test,y_test
In [3]: # let's try three different preprocessing techniques and collect the model coef
        from sklearn.linear model import Ridge
        from sklearn.pipeline import Pipeline
        from sklearn.model selection import ParameterGrid
        from sklearn.model selection import train test split
```

```
from sklearn.model selection import KFold
from sklearn.model selection import GridSearchCV
from sklearn.metrics import make_scorer
from sklearn.metrics import r2 score
from sklearn.preprocessing import StandardScaler, OrdinalEncoder
from sklearn.compose import ColumnTransformer
param_grid = {'regressor__alpha': [1e-3,1e-2,1e-1,1e0,1e1,1e2]}
num_ftrs = ['AGE', 'BMI', 'BP', 'S1', 'S2', 'S3', 'S4', 'S5', 'S6']
ord ftrs = ['SEX']
# ordinal encoder
ordinal_transformer = Pipeline(steps=[
    ('ordinal', OrdinalEncoder())])
# standard scaler
numeric transformer = Pipeline(steps=[
    ('scaler', StandardScaler())])
# ordinal encode SEX only!
preprocessor1 = ColumnTransformer(remainder = 'passthrough',
   transformers=[
        ('ord', ordinal_transformer, ord_ftrs)])
model, test_score, train_set, test_set = MLpipe_KFold_RMSE(X, y, preprocessor1,
print(test_score)
coefs no prep = model['regressor'].coef
print(coefs_no_prep)
# ordinal encode SEX, and standardize continuous features
preprocessor2 = ColumnTransformer(
   transformers=[
        ('num', numeric_transformer, num_ftrs),
        ('ord', ordinal transformer, ord ftrs)])
model, test_score, train_set, test_set = MLpipe_KFold_RMSE(X, y, preprocessor2,
print(test score)
coefs some prep = model['regressor'].coef
print(coefs_some_prep)
# ordinal encode SEX, standardize continuous features, and standarize all featu
preprocessor3 = Pipeline(steps=[
    ('prep2', preprocessor2),
    ('scaler', StandardScaler())])
model, test score, train set, test set = MLpipe KFold RMSE(X, y, preprocessor3,
print(test score)
coefs prep = model['regressor'].coef
print(coefs prep)
```

```
{'regressor alpha': 1.0}
        0.4520945451686046 0.4520945451686046
        0.4520945451686046
        [-22.80861461 0.14269511
                                     5.90541998 1.19867986 -1.07900835
           0.62662466 0.3774738
                                     9.77013169 60.79394666 0.21396887]
        {'regressor_alpha': 1.0}
        0.45418697357725524 0.45418697357725524
        0.45418697357725524
        [ \quad 1.79050083 \quad 25.75648091 \quad 16.71212465 \quad -34.65085803 \quad 17.05080545
                       11.71899321 31.39253301 2.45268588 -22.69228495]
           3.3870545
        {'regressor alpha': 1.0}
        0.4541465207069819 0.4541465207069819
        0.4541465207069819
        [ 1.80734179 25.73269892 16.73429974 -34.67195409 17.05307485
           3.36991411 11.76426044 31.3783838
                                                 2.45813922 -11.44818951]
In [4]: import matplotlib.pyplot as plt
        import numpy as np
        plt.rcParams.update({'font.size': 14})
        def plot_coefs(coefs,title,color,feature_names):
            feature_names = np.array(feature_names)
            sorted_indcs = np.argsort(np.abs(coefs))
            plt.barh(np.arange(10),coefs[sorted indcs])
            plt.yticks(np.arange(10),feature names[sorted indcs])
            plt.xlabel('coefficient')
            plt.title(title,color=color)
            return
        plt.figure(figsize=(10,5))
        plt.subplot(1,3,1)
        plot coefs(coefs no prep, 'no prep', 'r', ord ftrs + num ftrs)
        plt.annotate('BAD',xy=(10,10),xycoords = 'axes points',color='r')
        plt.subplot(1,3,2)
        plot_coefs(coefs_some_prep,'cont. ftrs. standardized','r', num_ftrs + ord_ftrs)
        plt.annotate('BAD',xy=(10,10),xycoords = 'axes points',color='r')
        plt.subplot(1,3,3)
        plot_coefs(coefs_prep, 'all ftrs. standardized', 'g', num_ftrs + ord_ftrs)
        plt.annotate('Correct!',xy=(10,10),xycoords = 'axes points',color='g')
        plt.tight_layout()
        plt.savefig('figures/coefs comparison.png',dpi=300)
        plt.show()
```



Introduction to Interpretability in ML

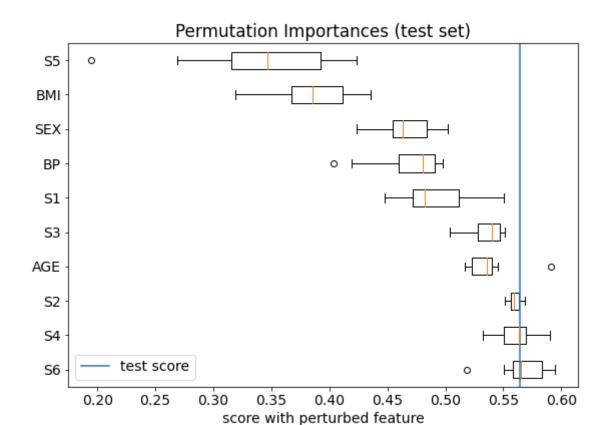
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- use the SHAP package to calculate local explanations

Permutation feature importance

- model agnostic, you can use it with any supervised ML model
- steps:
 - train a model and calculate a test score :)
 - randomly shuffle a single feature in the test set
 - recalculate the test score with the shuffled data
 - model score worsens because the shuffling breaks the relationship between feature and target
 - the larger the difference, the more important the feature is

```
y test = test set[1]
        ftr_names = X_test.columns
        nr runs = 10
        scores = np.zeros([len(ftr_names),nr_runs])
        print('test score = ',test score)
        # loop through the features
        for i in range(len(ftr_names)):
            print('shuffling '+str(ftr_names[i]))
            r2_scores = []
            for j in range(nr_runs):
                X_test_shuffled = X_test.copy()
                X_test_shuffled[ftr_names[i]] = np.random.permutation(X_test[ftr_names[
                r2 scores.append(model.score(X test shuffled,y test))
            print(' shuffled test score:',np.around(np.mean(r2_scores),3),'+/-',np.ar
            scores[i] = r2_scores
        test score = 0.5645801016038898
        shuffling AGE
           shuffled test score: 0.537 +/- 0.02
        shuffling SEX
           shuffled test score: 0.467 +/- 0.021
        shuffling BMI
           shuffled test score: 0.384 +/- 0.037
        shuffling BP
           shuffled test score: 0.468 +/- 0.031
        shuffling S1
           shuffled test score: 0.493 +/- 0.034
        shuffling S2
           shuffled test score: 0.56 +/- 0.005
        shuffling S3
           shuffled test score: 0.535 +/- 0.014
        shuffling S4
           shuffled test score: 0.561 +/- 0.016
        shuffling S5
           shuffled test score: 0.34 +/- 0.066
        shuffling S6
           shuffled test score: 0.567 +/- 0.021
In [7]: sorted_indcs = np.argsort(np.mean(scores,axis=1))[::-1]
        plt.rcParams.update({'font.size': 14})
        plt.figure(figsize=(8,6))
        plt.boxplot(scores[sorted_indcs].T,labels=ftr_names[sorted_indcs],vert=False)
        plt.axvline(test score, label='test score')
        plt.title("Permutation Importances (test set)")
        plt.xlabel('score with perturbed feature')
        plt.legend()
        plt.tight_layout()
        plt.show()
```



Check out sklearn's permutation importance!

https://scikit-

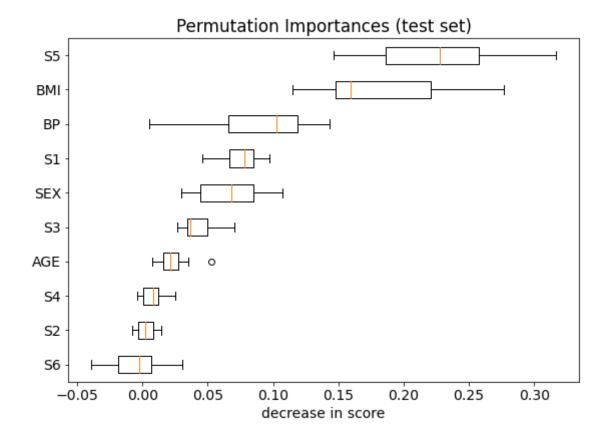
learn.org/stable/modules/generated/sklearn.inspection.permutation_importance.html

https://scikit-learn.org/stable/modules/permutation_importance.html#permutation-importance

```
In [8]: from sklearn.inspection import permutation_importance
    result = permutation_importance(model, X_test, y_test,n_repeats=10,random_state
    ftr_names = X_test.columns

sorted_indcs = result.importances_mean.argsort()

plt.rcParams.update({'font.size': 14})
    plt.figure(figsize=(8,6))
    plt.boxplot(result.importances[sorted_indcs].T,labels=ftr_names[sorted_indcs],v
    plt.title("Permutation Importances (test set)")
    plt.xlabel('decrease in score')
    plt.tight_layout()
    plt.show()
```



Cons of permutation feature importance

- strongly correlated features
 - if one of the features is shuffled, the model can still use the other correlated feature
 - both features appear to be less important but they might actually be important
 - solution:
 - o check the correlation matrix plot
 - remove all but one of the strongly correlated features
- · no feature interactions
 - one feature might appear unimportant but combined with another feature could be important
 - solution:
 - o permute two features to measure how important feature pairs are
 - this can be computationally expensive

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SHAP values

- one way to calculate local feature importances
- it is based on Shapely values from game theory
- read more here, here, and here

Cooperative game theory

- A set of *m* players in a coalition generate a surplus.
- Some players contribute more to the coalition than others (different bargaining powers).
- How important is each player to the coalition?
- How should the surplus be divided fairly amongst the players?

Cooperative game theory applied to feature attribution

- A set of *m* features in a model generate a prediction.
- Some features contribute more to the model than others (different predictive powers).
- How important is each feature to the model?
- How should the prediction be divided amongst the features?

How is it calculated?

$$\Phi_i = \sum_{S \subseteq M \setminus i} rac{|S|!(M-|S|-1)!}{M!} [f_x(S \cup i) - f_x(S)]$$

- Φ_i the contribution of feature i
- ullet M the number of features
- ullet S a set of features excluding i, a vector of 0s and 1s (0 if a feature is missing)
- |S| the number of features in S
- $f_x(S)$ the prediction of the model with features S

How is it calculated?

$$\Phi_i = \sum_{S \subset M \setminus i} rac{|S|!(M-|S|-1)!}{M!} [f_x(S \cup i) - f_x(S)]$$

- the difference feature i makes in the prediction:
 - $f_x(S \cup i)$ the prediction with feature i
 - $f_x(S)$ the prediction without feature i
- loop through all possible ways a set of S features can be selected from the M features excluding i
- ullet weight the contribution based on how many ways we can select |S| features

```
import shap
shap.initjs() # required for visualizations later on

explainer = shap.KernelExplainer(model[1].predict,model[0].transform(train_set[
# transform the test set
X_test_transformed = model[0].transform(X_test)
print(np.shape(X_test_transformed))
# calculate shap values on the first 1000 points in the test
shap_values = explainer.shap_values(X_test_transformed)
print(np.shape(shap_values))
ftr_names = num_ftrs + ord_ftrs
```

jς

Using 353 background data samples could cause slower run times. Consider using shap.sample(data, K) or shap.kmeans(data, K) to summarize the background as K samples.

```
(89, 10)

0% | | 0/89 [00:00<?, ?it/s]

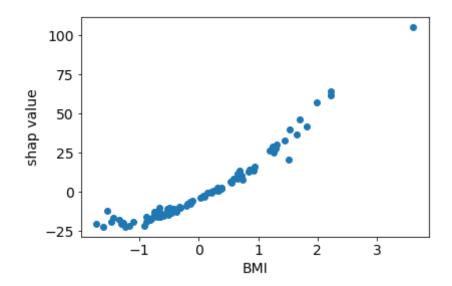
(89, 10)
```

Explain a point

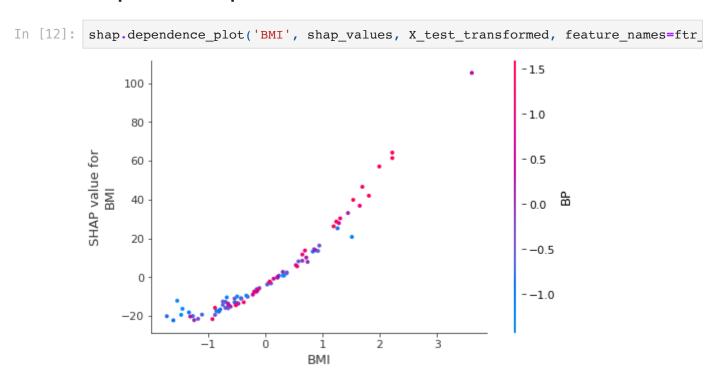
Feature value vs. shap value

```
In [11]: ftr = 'BMI'
   indx = np.argwhere(ftr_names==ftr)

plt.scatter(X_test_transformed[:,indx],shap_values[:,indx])
   plt.ylabel('shap value')
   plt.xlabel(ftr)
   plt.tight_layout()
   plt.show()
```

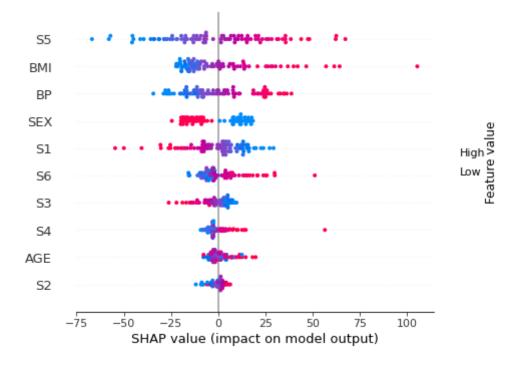


Dependence plot



It can also be used for global feature importance

In [13]: shap.summary_plot(shap_values, X_test_transformed,feature_names = ftr_names)



SHAP cons

- · it can be numerically expensive
 - an efficient shap method was developed for trees, see here
- how to estimate $f_x(S)$?
 - this is not trivial because models cannot change the number of features they use
 - usually the values of the dropped features are replaced with the mean or 0
 - this is an approximation

Recap: Introduction to Interpretability in ML

By now, you can

- calculate global explanations using the coefficients of linear models
- · perform permutation importance using any ML model
- use the SHAP package to calculate local explanations

Brown Data Science Master's Program

- if your company is looking to hire talented data scientists, reach out to me
- our students are looking for internships and full time positions!
- email: andras_zsom@brown.edu
- I'm also around tomorrow and until the early afternoon on Thursday