## 13-L1 Problem 1

We return to the plane-strain compression problem where the goal is to predict von Mises stress at a node given a set of its features.

Now, you will look at ensemble methods in sklearn to determine how well they perform in this context.

```
import numpy as np
import matplotlib.pyplot as plt
from sklearn.linear model import LinearRegression
from sklearn.tree import DecisionTreeRegressor
from sklearn.metrics import mean squared error
from sklearn.ensemble import StackingRegressor
def plot shape(dataset, index, model=None, lims=None):
    x = dataset["coordinates"][index][:,0]
    y = dataset["coordinates"][index][:,1]
    if model is None:
        c = dataset["stress"][index]
    else:
        c = model.predict(dataset["features"][index])
    if lims is None:
        lims = [min(c), max(c)]
    plt.scatter(x,y,s=5,c=c,cmap="jet",vmin=lims[0],vmax=lims[1])
    plt.colorbar(orientation="horizontal", shrink=.75,
pad=0,ticks=lims)
    plt.axis("off")
    plt.axis("equal")
def plot shape comparison(dataset, index, model, title=""):
    plt.figure(figsize=[6,3.2], dpi=120)
    plt.subplot(1,2,1)
    plot shape(dataset,index)
    plt.title("Ground Truth", fontsize=9, y=.96)
    plt.subplot(1,2,2)
    c = dataset["stress"][index]
    plot shape(dataset, index, model, lims = [min(c), max(c)])
    plt.title("Prediction", fontsize=9, y=.96)
    plt.suptitle(title)
    plt.show()
```

```
def load dataset(path):
    dataset = np.load(path)
    coordinates = []
    features = []
    stress = []
    N = np.max(dataset[:,0].astype(int)) + 1
    split = int(N*.8)
    for i in range(N):
        idx = dataset[:,0].astype(int) == i
        data = dataset[idx,:]
        coordinates.append(data[:,1:3])
        features.append(data[:,3:-1])
        stress.append(data[:,-1])
    dataset_train = dict(coordinates=coordinates[:split],
features=features[:split], stress=stress[:split])
    dataset test = dict(coordinates=coordinates[split:],
features=features[split:], stress=stress[split:])
    X train, X test = np.concatenate(features[:split], axis=0),
np.concatenate(features[split:], axis=0)
    y_train, y_test = np.concatenate(stress[:split], axis=0),
np.concatenate(stress[split:], axis=0)
    return dataset train, dataset test, X train, X test, y train,
y test
def get shape(dataset,index):
    X = dataset["features"][index]
    y = dataset["stress"][index]
    return X, y
```

### Loading the data

First, complete the code below to load the data and plot the von Mises stress fields for a few shapes.

You'll need to input the path of the data file, the rest is done for you.

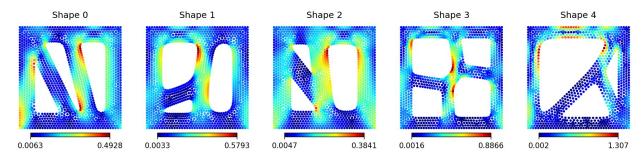
All training node features and outputs are in  $X_train}$  and  $y_train$ , respectively. Testing nodes are in  $X_test$ ,  $y_test$ .

dataset\_train and dataset\_test contain more detailed information such as node coordinates, and they are separated by shape.

Get features and outputs for a shape by calling get\_shape(dataset,index). N\_train and N test are the number of training and testing shapes in each of these datasets.

```
# YOU MAY NEED TO EDIT data_path
data_path = "stress_nodal_features.npy"
dataset_train, dataset_test, X_train, X_test, y_train, y_test =
load_dataset(data_path)
N_train = len(dataset_train["stress"])
N_test = len(dataset_test["stress"])
```

```
plt.figure(figsize=[15,3.2], dpi=150)
for i in range(5):
    plt.subplot(1,5,i+1)
    plot_shape(dataset_train,i)
    plt.title(f"Shape {i}")
plt.show()
```



## StackingRegressor

A StackingRegressor consists of N fitted regression models, which it evaluates to make N predictions. Then, these predictions are fed into another regression model, which makes a final prediction of the target.

### List of Regressors

First, initialize 4 regression models:

- 1. Linear Regression
- 2. Decision Tree regression, max depth 4
- 3. Decision Tree regression, max depth 8
- 4. Decision Tree regression, max depth 12

Then, store these in a list called models.

```
# YOUR CODE GOES HERE
# Define models, and put in a list called 'models'
model1 = LinearRegression()
model2 = DecisionTreeRegressor(max_depth=4)
model3 = DecisionTreeRegressor(max_depth=8)
model4 = DecisionTreeRegressor(max_depth=12)
models = [model1, model2, model3, model4]

named_models = [(f"Model {i+1}", model) for i, model in
enumerate(models)]
print(*named_models, sep="\n")

('Model 1', LinearRegression())
('Model 2', DecisionTreeRegressor(max_depth=4))
```

```
('Model 3', DecisionTreeRegressor(max_depth=8))
('Model 4', DecisionTreeRegressor(max_depth=12))
```

#### Final Regressor

Now make one more regressor, which will take as input the other four predictions, and combine them to make an improved prediction.

This can be another linear regression model. Call it final model.

```
# YOUR CODE GOES HERE
final_model = LinearRegression()
```

### Creating and training the StackingRegressor

Finally, we can combine all of our models into a StackingRegressor model. We fit this just as we would fit any sklearn model. Because of the size of the dataset, this may take a few minutes.

#### Performance of the model

Now we can investigate the performance of the model on test data, compared to constituent models. First, let's look at the performance of each individual model of our model srm. These can be accessed via srm.estimators\_.

```
for i, estimator in enumerate(srm.estimators_):
    print(f"\n{named_models[i][0]}:")
    train_err = mean_squared_error(y_train,
estimator.predict(X_train))
    test_err = mean_squared_error(y_test, estimator.predict(X_test))
    print(f"Training MSE: {train_err:.2e}")
    print(f" Testing MSE: {test_err:.2e}")
Model 1:
Training MSE: 8.11e-03
Testing MSE: 9.78e-03
```

```
Model 2:
Training MSE: 1.26e-02
Testing MSE: 1.51e-02

Model 3:
Training MSE: 7.56e-03
Testing MSE: 1.04e-02

Model 4:
Training MSE: 3.75e-03
Testing MSE: 8.33e-03
```

#### Stacking Regressor MSE on Test Data

Now compute the MSE of srm on training and testing data.

Note how the results, particularly on test data, compare to the individual models.

```
# YOUR CODE GOES HERE
train_mse = mean_squared_error(y_train, srm.predict(X_train))
test_mse = mean_squared_error(y_test, srm.predict(X_test))

print(f"\nStacking Regressor (final model):")
print(f"Training MSE: {train_mse:.2e}")
print(f"Testing MSE: {test_mse:.2e}")

Stacking Regressor (final model):
Training MSE: 4.10e-03
Testing MSE: 6.60e-03
```

## M13-L2 Problem 1

Once more, we will study the stress prediction problem, this time using XGBoost, a very powerful boosting method.

```
import numpy as np
import matplotlib.pyplot as plt
import xqboost as xqb
from xgboost import XGBRegressor
from sklearn.metrics import mean squared error
def plot shape(dataset, index, model=None, lims=None):
    x = dataset["coordinates"][index][:,0]
    y = dataset["coordinates"][index][:,1]
    if model is None:
        c = dataset["stress"][index]
    else:
        c = model.predict(dataset["features"][index])
    if lims is None:
        lims = [min(c), max(c)]
    plt.scatter(x,y,s=5,c=c,cmap="jet",vmin=lims[0],vmax=lims[1])
    plt.colorbar(orientation="horizontal", shrink=.75,
pad=0,ticks=lims)
    plt.axis("off")
    plt.axis("equal")
def plot shape comparison(dataset, index, model, title=""):
    plt.figure(figsize=[6,3.2], dpi=120)
    plt.subplot(1,2,1)
    plot_shape(dataset,index)
    plt.title("Ground Truth", fontsize=9, y=.96)
    plt.subplot(1,2,2)
    c = dataset["stress"][index]
    plot shape(dataset, index, model, lims = [min(c), max(c)])
    plt.title("Prediction", fontsize=9, y=.96)
    plt.suptitle(title)
    plt.show()
def load dataset(path):
    dataset = np.load(path)
    coordinates = []
    features = []
    stress = []
```

```
N = np.max(dataset[:,0].astype(int)) + 1
    split = int(N*.8)
    for i in range(N):
        idx = dataset[:,0].astype(int) == i
        data = dataset[idx,:]
        coordinates.append(data[:,1:3])
        features.append(data[:,3:-1])
        stress.append(data[:,-1])
    dataset_train = dict(coordinates=coordinates[:split],
features=features[:split], stress=stress[:split])
    dataset test = dict(coordinates=coordinates[split:],
features=features[split:], stress=stress[split:])
    X_train, X_test = np.concatenate(features[:split], axis=0),
np.concatenate(features[split:], axis=0)
    y_train, y_test = np.concatenate(stress[:split], axis=0),
np.concatenate(stress[split:], axis=0)
    return dataset train, dataset test, X train, X test, y train,
y_test
def get shape(dataset,index):
    X = dataset["features"][index]
    y = dataset["stress"][index]
    return X, v
def eval model(model, verbose=False):
    pred train = model.predict(X train)
    pred test = model.predict(X test)
    mse_train = mean_squared_error(y_train, pred_train)
    mse test = mean_squared_error(y_test, pred_test)
    if verbose:
        print(f"Train MSE = {mse train:.2e}")
        print(f"Test MSE = {mse_test:.2e}")
    return mse train, mse test
```

## Loading the data

First, complete the code below to load the data and plot the von Mises stress fields for a few shapes.

You'll need to input the path of the data file, the rest is done for you.

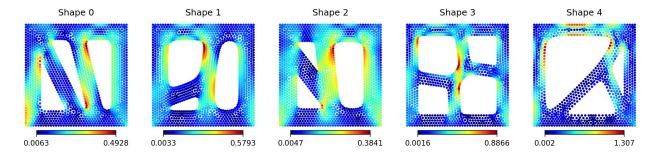
All training node features and outputs are in X\_train and y\_train, respectively. Testing nodes are in X\_test, y\_test.

dataset\_train and dataset\_test contain more detailed information such as node coordinates, and they are separated by shape.

Get features and outputs for a shape by calling get\_shape(dataset,index). N\_train and N\_test are the number of training and testing shapes in each of these datasets.

```
# YOU MAY NEED TO EDIT data_path
data_path = "stress_nodal_features.npy"
dataset_train, dataset_test, X_train, X_test, y_train, y_test =
load_dataset(data_path)
N_train = len(dataset_train["stress"])
N_test = len(dataset_test["stress"])

plt.figure(figsize=[15,3.2], dpi=150)
for i in range(5):
    plt.subplot(1,5,i+1)
    plot_shape(dataset_train,i)
    plt.title(f"Shape {i}")
plt.show()
```



# XGBoost Regressor

XGBoost models, like XGBRegressor here, can be used much like sklearn models.

First, define an instance of XGBRegressor with the desired parameters; then, fit the model with model.fit. You can evaluate a fitted model with model.predict.

The provided function mse\_train, mse\_test = eval\_model(model) to get MSE values on the train and test datasets.

```
eta = 0.8
depth = 9

params = dict(
    eta = eta,
    max_depth = depth,
)

model = XGBRegressor(objective ='reg:squarederror', seed = 123,
    n_estimators = 10, **params)
model.fit(X_train, y_train)

mse_train, mse_test = eval_model(model)
print(" eta depth | Train MSE Test MSE")
print("-------")
```

## Parametric study

Now let's examine the effects of varying the parameters eta and max\_depth, keeping n\_estimators as 10. For every combination of eta in [0.1, 0.3, 0.5, 0.7] and max\_depth in [5, 10, 15, 20], train an XGB regressor and report the train and test MSE values.

Which combination has the best performance on testing data?

```
# YOUR CODE GOES HERE
eta = [0.1, 0.3, 0.5, 0.7]
depth = [5, 10, 15, 20]
print(" eta depth | Train MSE Test MSE")
print("------
for e in eta:
   for d in depth:
      params = dict(
          eta = e,
          max depth = d,
      )
      model = XGBRegressor(objective ='reg:squarederror', seed =
123, n estimators = 10, **params)
      model.fit(X_train, y_train)
      mse train, mse test = eval model(model)
      # print(" eta depth | Train MSE Test MSE")
      # print("-----
      print(f" {e:.1f} {d:>2d} | {mse_train:.2e}
{mse test:.2e}")
      print("-----")
 eta
      depth |
                Train MSE Test MSE
 0.1 5
                 1.05e-02 1.28e-02
 0.1 10
                 5.93e-03 8.93e-03
                 3.86e-03 7.93e-03
 0.1 15
             3.35e-03 7.98e-03
 0.1 20
```

	0.3	5	5.43e-03	7.05e-03
	0.3	10	1.57e-03	4.58e-03
-	0.3	15	2.79e-04	4.59e-03
-	0.3	20	8.02e-05	4.73e-03
	0.5	5	4.60e-03	6.49e-03
	0.5	10	1.36e-03	4.80e-03
	0.5	15	1.47e-04	5.01e-03
-	0.5	20	8.30e-06	5.20e-03
	0.7	5	4.83e-03	7.03e-03
	0.7	10	1.44e-03	5.54e-03
	0.7	15	1.60e-04	5.82e-03
	0.7	20	9.05e-06	6.02e-03
		<b></b>	-1-3	

When eta = 0.3 and depth = 10, the model performed the best on the given testing data.