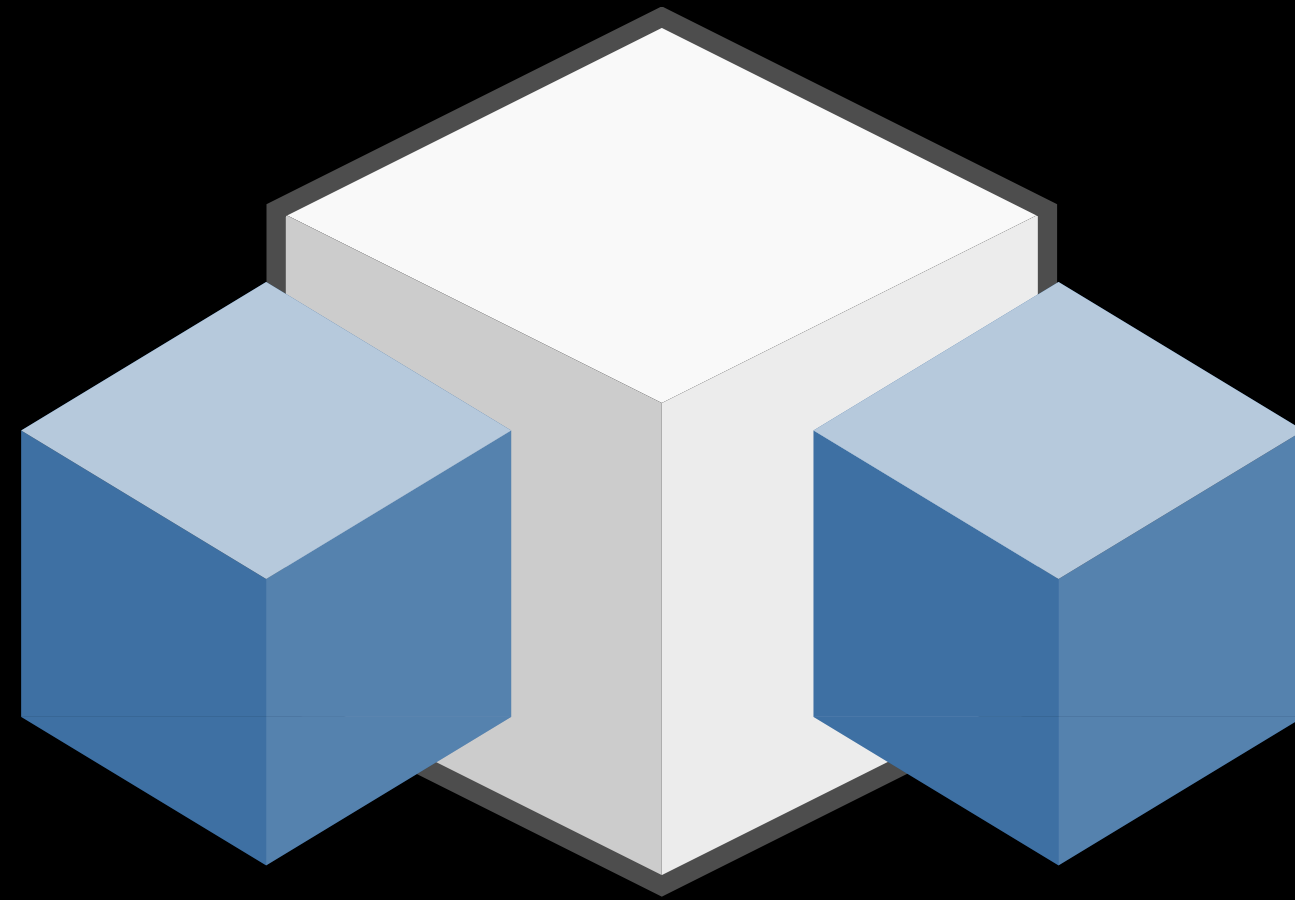


# Prediction of Orthorhombic Lattice Constants

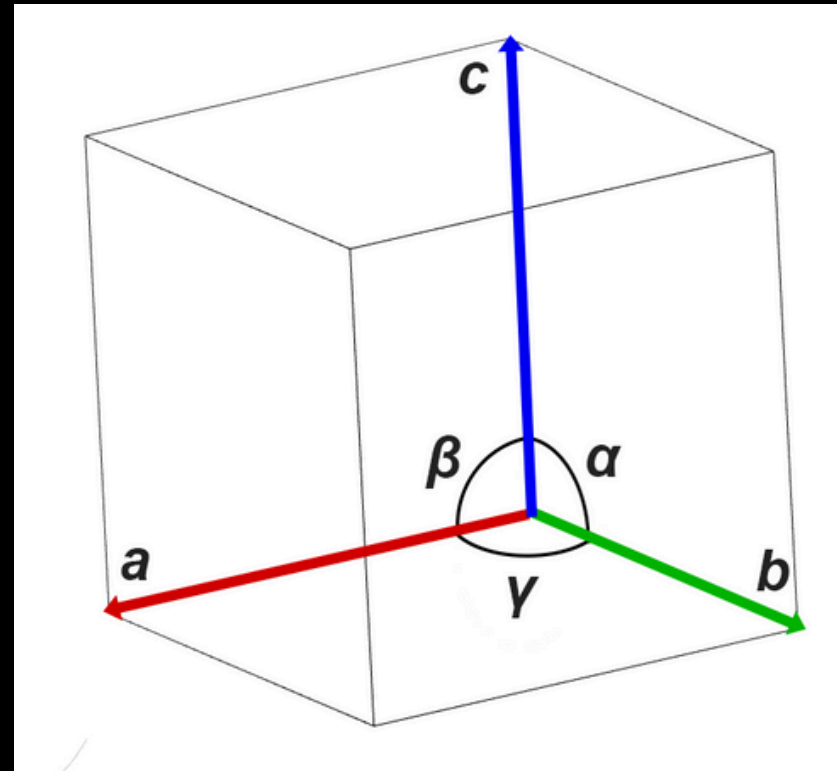


But why?

The structure of crystal materials  
is denoted by the lattice structure.

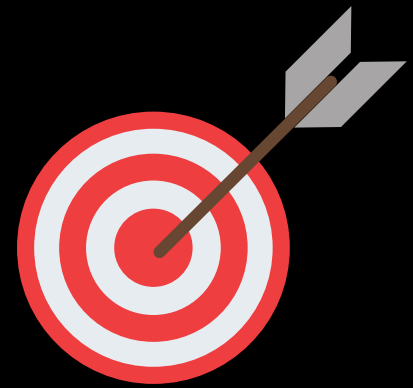
The lattice structure is  
a repeating pattern.  
Its basic unit is the unit cell.

The unit cell is fully  
described by the six  
lattice constants.



These are:  
a, b, c;  
alpha, beta,  
gamma.

Many properties of the crystal material depend on the **lattice structure**, which depends on the **unit cell**, which depends on the **lattice constants**.



## Objectives:

- predict orthorhombic lattice constants
- examine feature importance

# Method to the Madness



THOU SHALT PREDICT.

# Method to the Madness

Dataset: Materials Project

Pre-processing: pandas, numpy

Model: random forest regression (from sklearn)

Visualization: matplotlib



# Method to the Madness

The dataset is taken from the Materials Project database  
and modified by Li et al.

It has been featurized using MAGPIE and the matminer library.  
(insert database shape)

# Method to the Madness

The dataset was already clean (NaN values, etc.) so I proceeded to filtering out the orthorhombic crystal materials.  
(I then perform further cleaning by removing the outliers.  
More on this later.)

# Method to the Madness

I passed the dataset through a function to  
separate features and targets.

I performed an 80–20 train–test–split for the  
features and targets.

# Method to the Madness

I used sklearn's random forest regressor class to create three RFR models with `n_estimators=300` and `max_features=300, 200, 100` respectively.

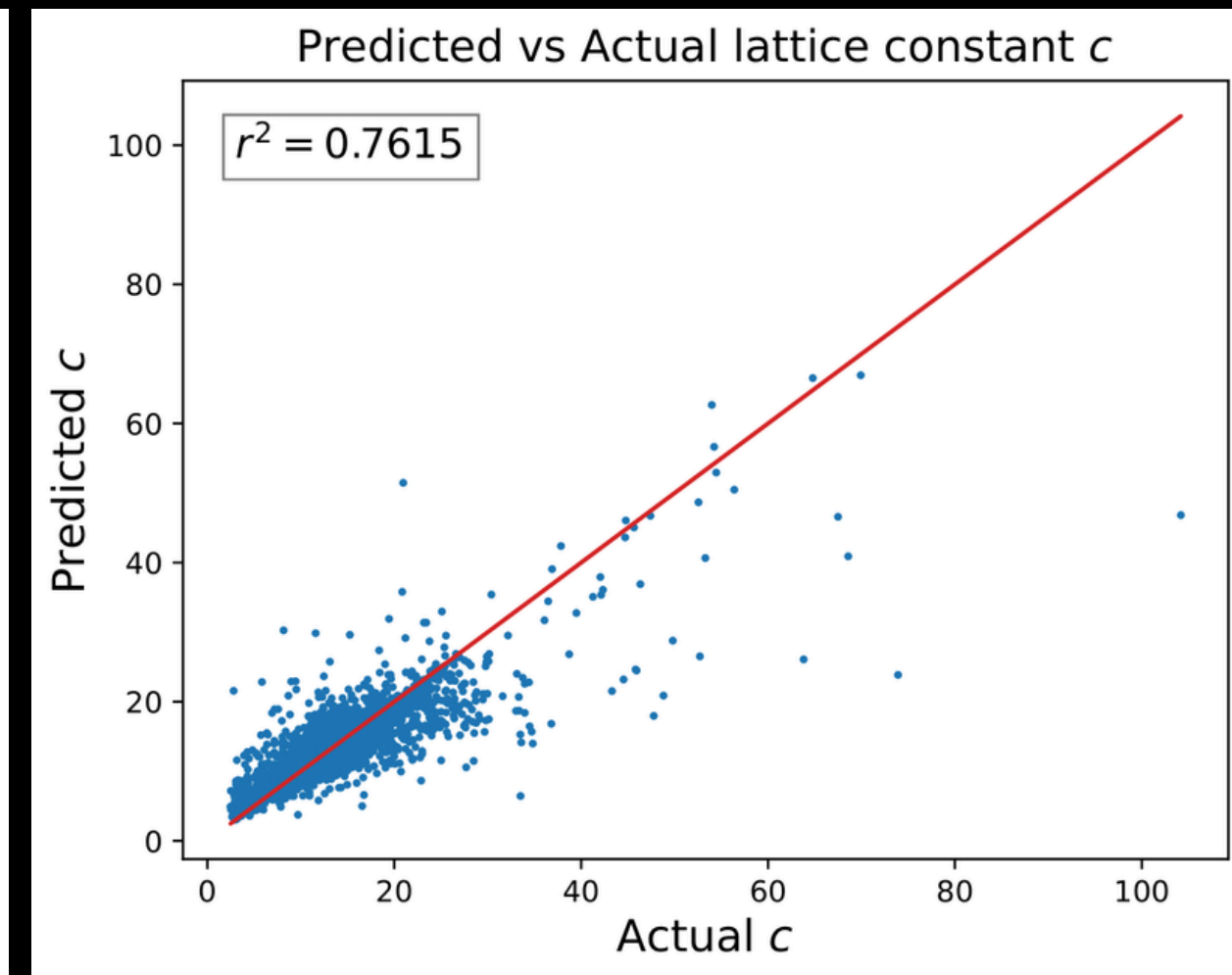
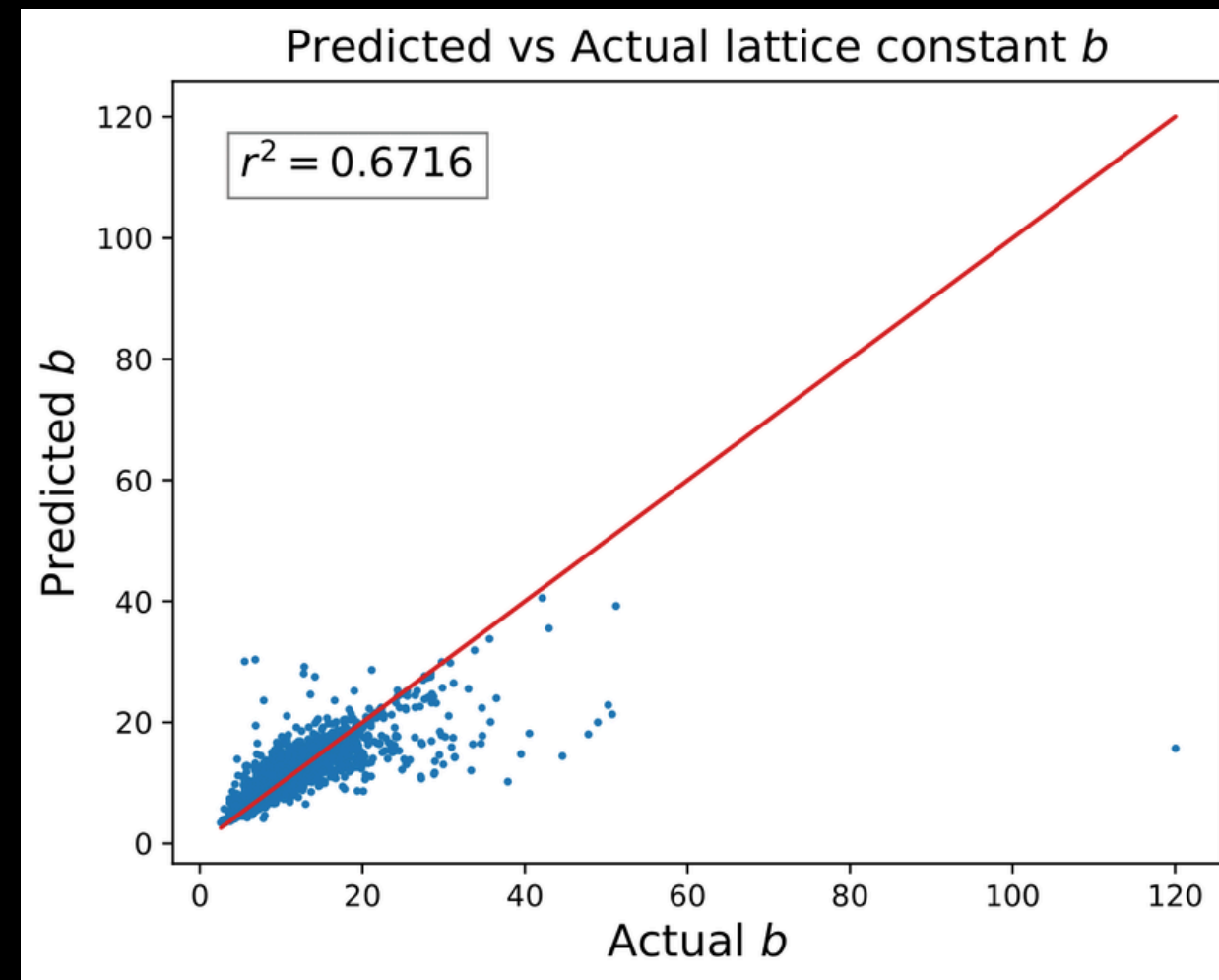
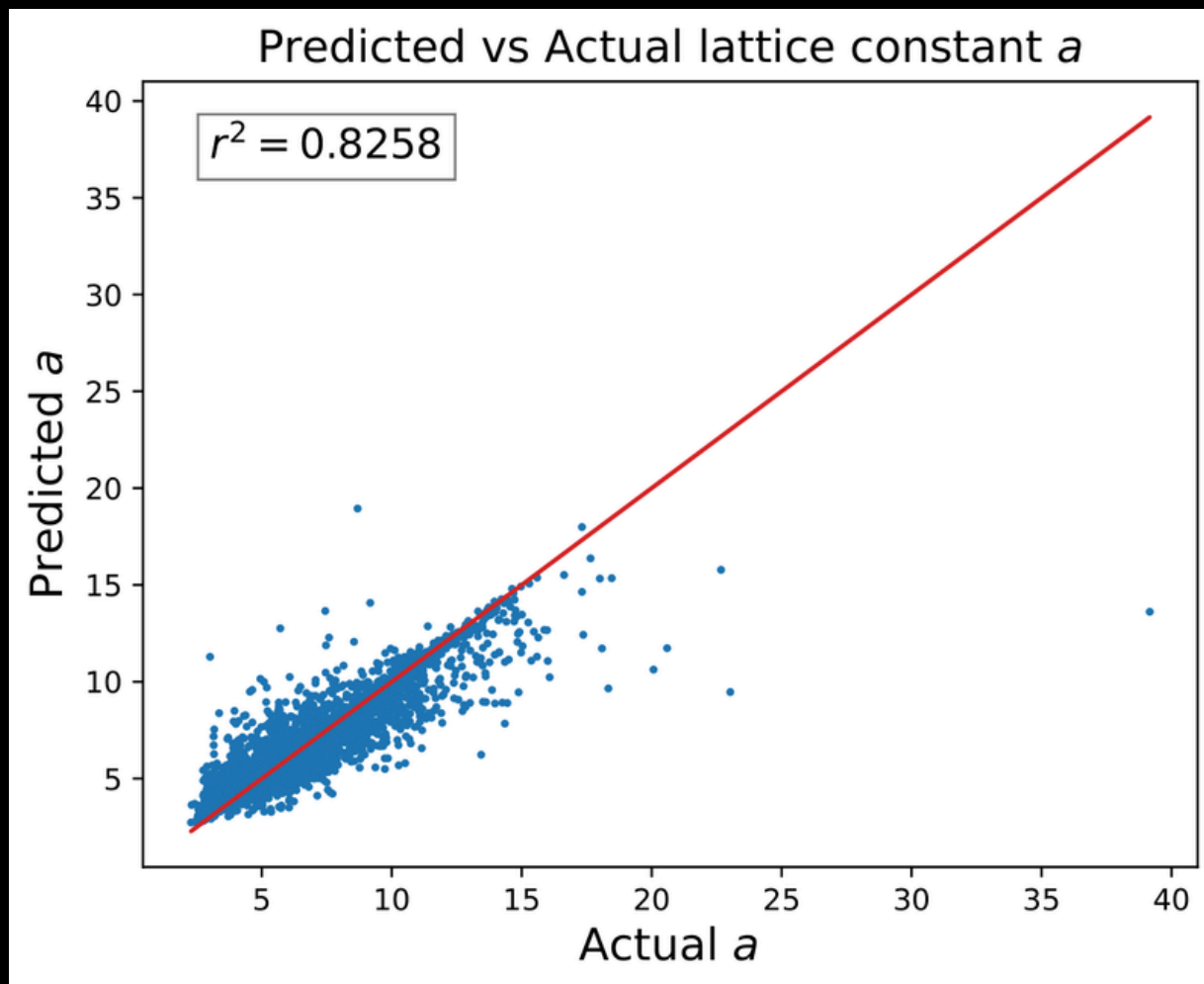
# Method to the Madness

I took the mean of the result of each RFR model  
as my final prediction.

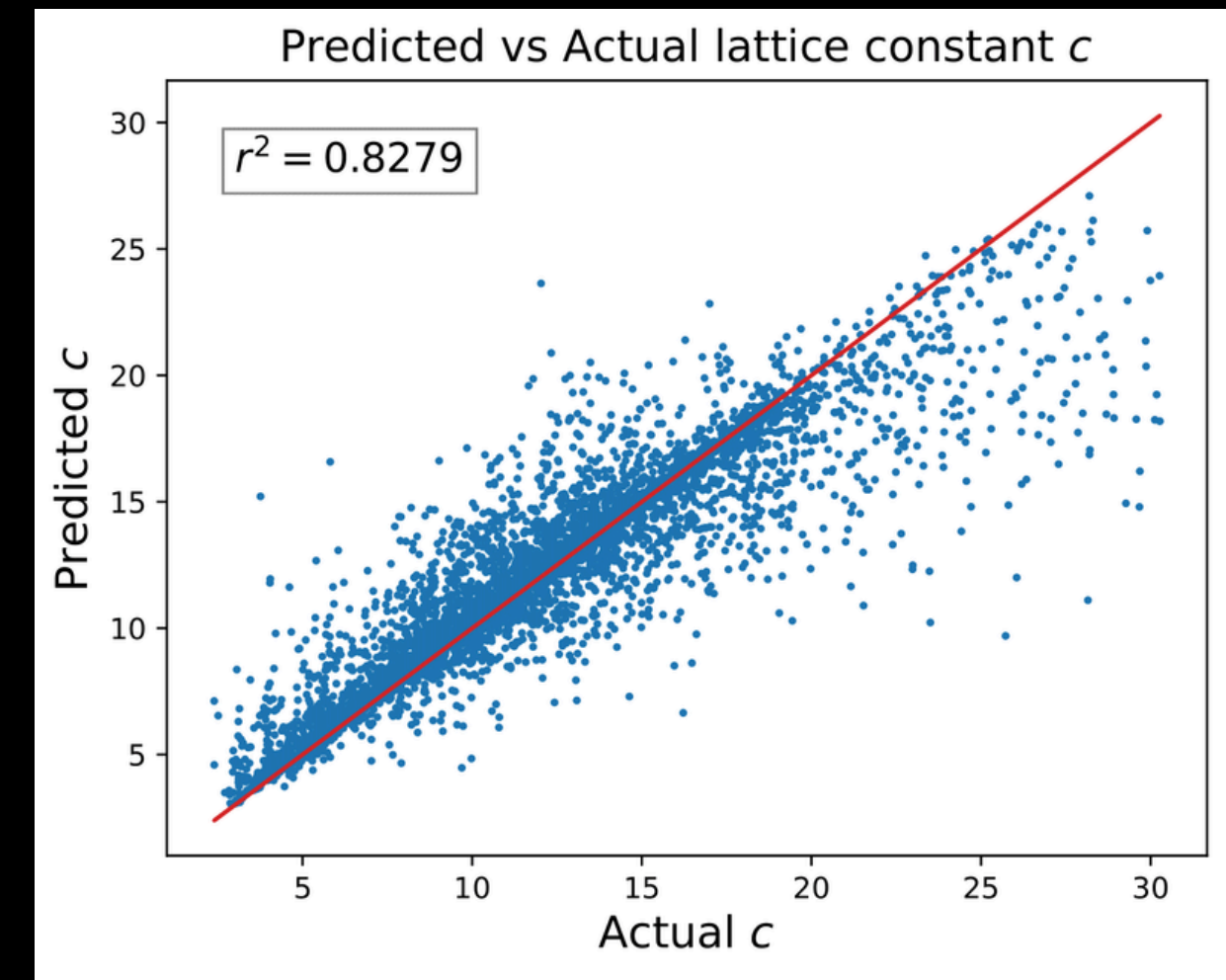
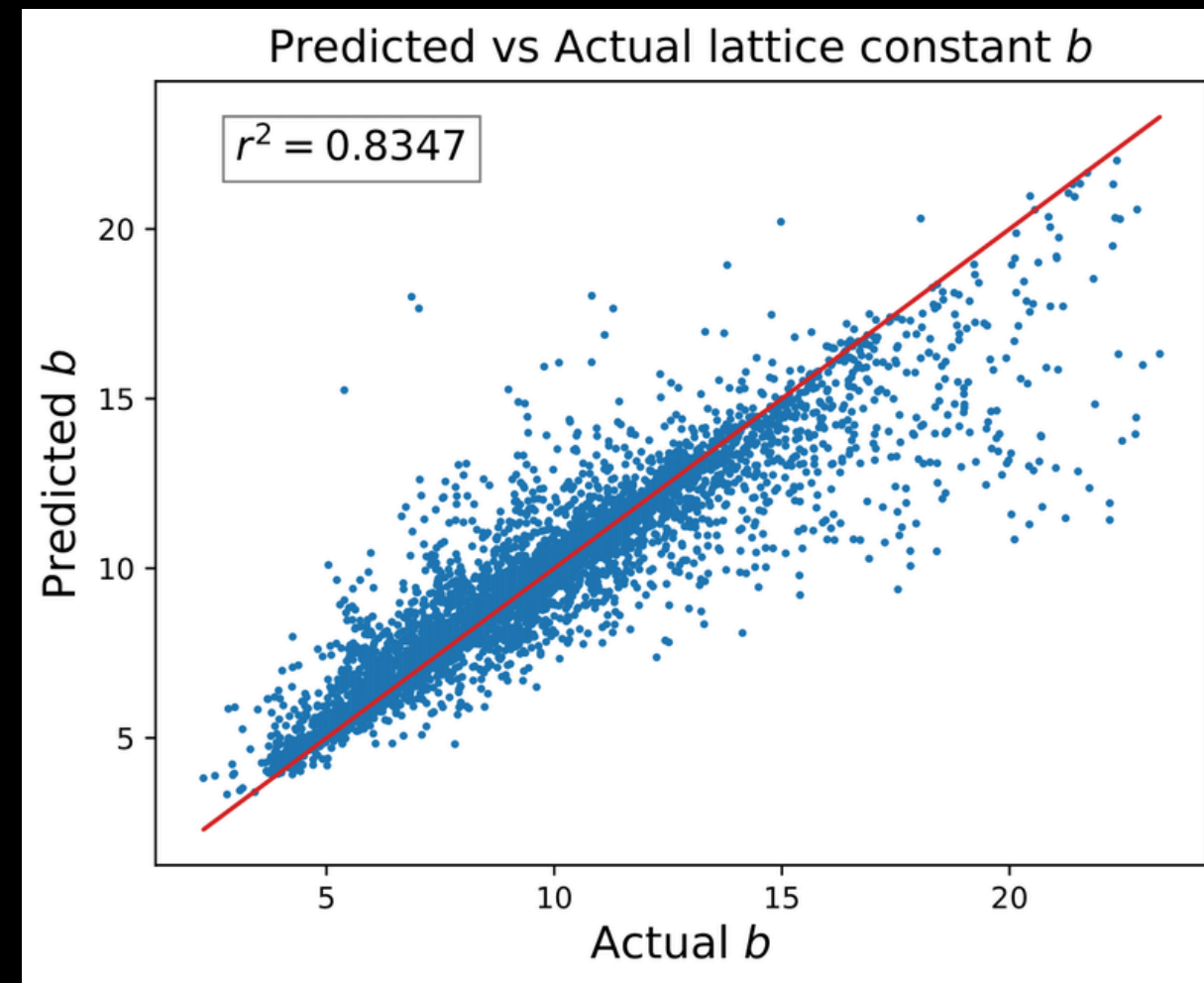
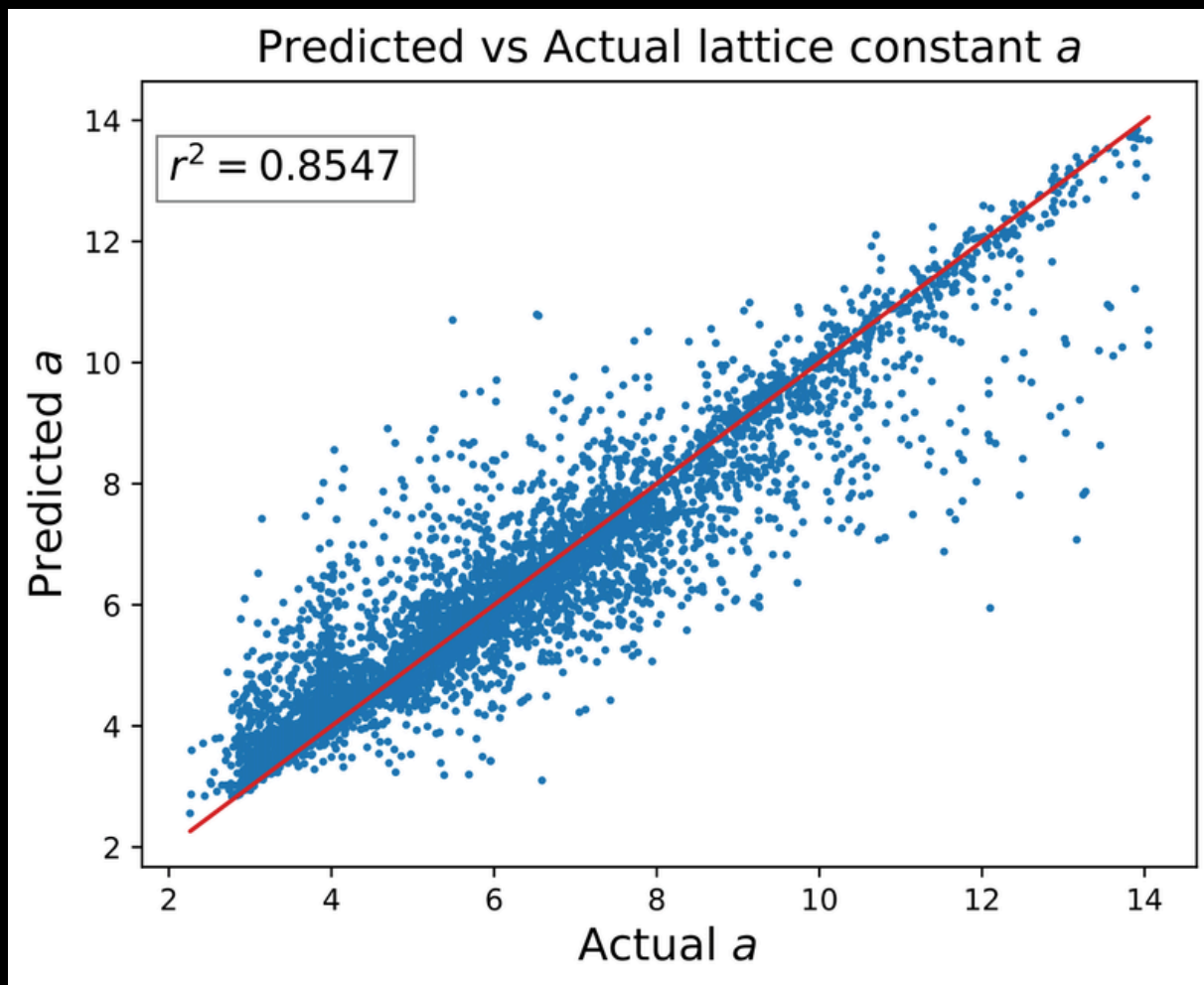
# Method to the Madness

I took three metrics: r-squared score, mean absolute error, and mean squared error to gauge the performance of the model.

# Fruits of Labor



# Fruits of Labor





# What did we learn?

Through random forest regression  
we can predict orthorhombic  
lattice constants to an acceptable  
degree of accuracy.

# What did we learn?

By cleaning the outliers in the training data, we improve the performance of our model.

# What did we learn?

However, this predictive model **cannot** replace DFT ab initio calculations. It can only supplement the investigation methodology through the vast speed-up of the scouring process.

# What did we learn?

By further tailoring our data sets,  
getting more comprehensive data  
and optimizing our model hyperparameters,  
we can improve the performance and  
utility of this method.

# References

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Thank you.