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# Background

Concerns regarding tight sandstone reservoirs have become crucial issues in unconventional oil and gas resources. Because of the heterogeneity of tight sandstone reservoirs, the pore structures are complex and the lithology and electrical properties vary greatly. Achieving ideal results from the evaluation of tight sandstone reservoirs is difficult from the perspective of regional adaptability using the routine analysis methods.

# Problem Statement

The microscopic pore structure of the reservoir directly affects the reservoir's seepage capacity, controls oil and gas migration and reservoir distribution, and ultimately determines the distribution of reservoir productivity. This project focuses on prediction of reservoir quality of tight sandstone reservoirs using high-pressure mercury-injection data. We used two popular machine learning methods for the reservoir quality prediction, including K-means and random forest models.

#### Datasets

This project collected the high-pressure mercury intrusion data of eight wells, divided into two groups for analysis and prediction. The distributions of the data sets are as follows:



# Machine Learning Technique for The Quantitative Evaluation of **Tight Sandstone Reservoirs using High-Pressure Mercury-Injection Data**

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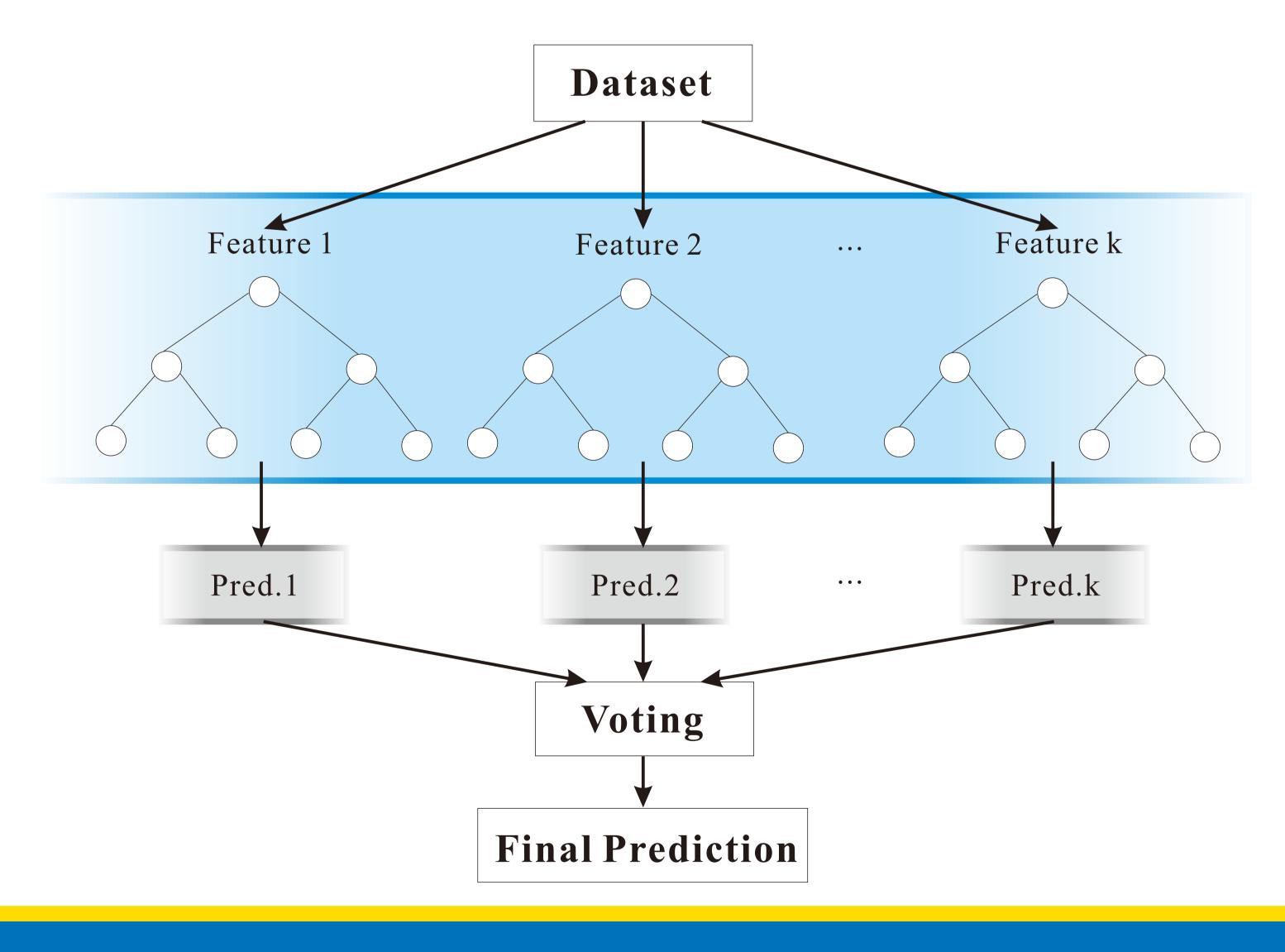
#### Models-K-means

The K-means clustering method is a non-genetic clustering method based on an iterative algorithm that clusters samples into k sets. The number k of classes can be predetermined or determined during the clustering process. The method has a small calculation cost and fast processing speed, which can be applied to a large data group.

# **Models-Random Forest**

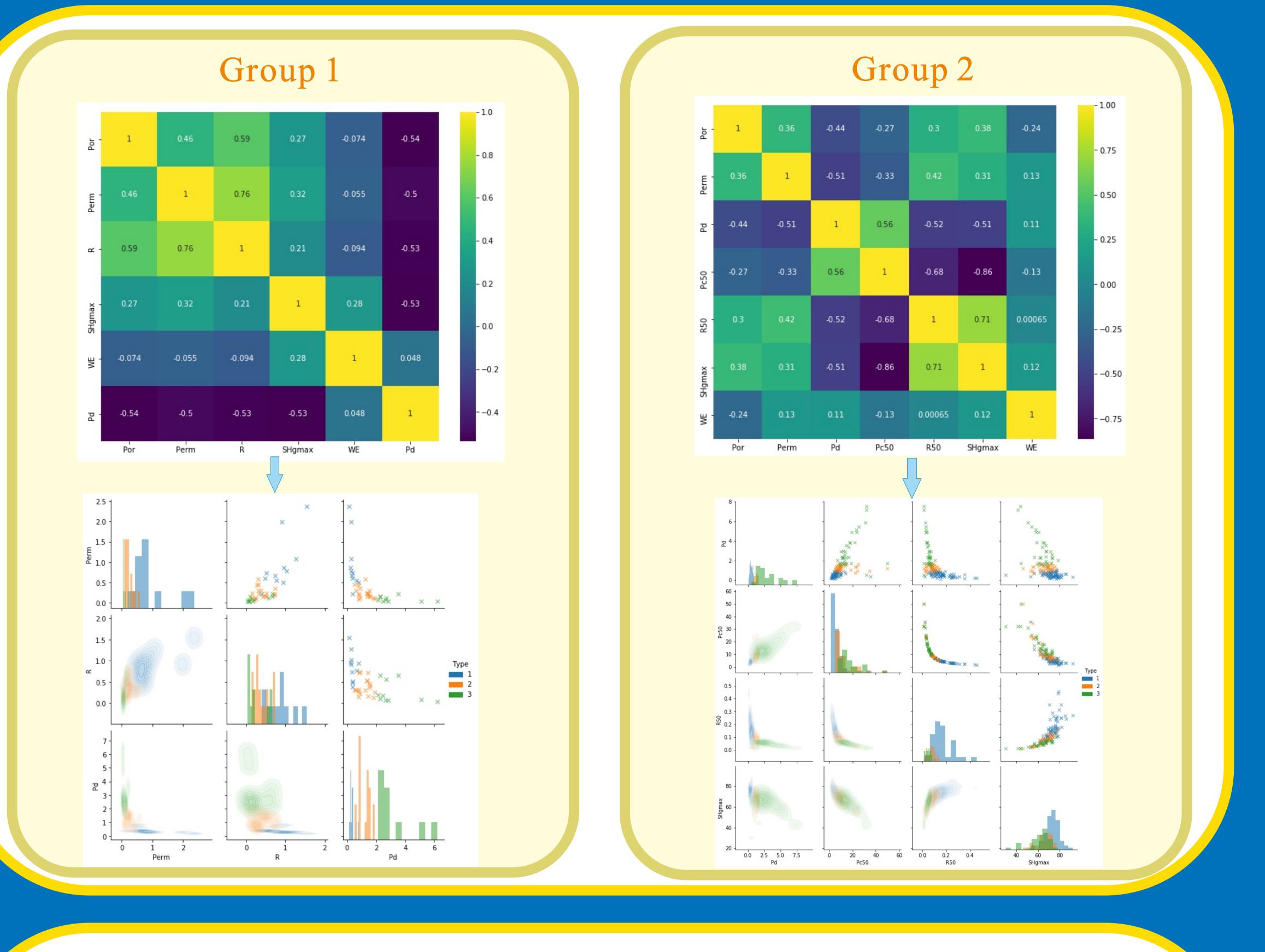
00000 39.000000 39.000000
02051 1.482821 1.974359
81262 1.310086 0.742938
70000 0.170000 1.000000
50000 0.565000 1.000000
20000 1.260000 2.000000
85000 2.015000 2.500000
90000 6.200000 3.000000
R50 SHgmax WE Type
140.000000 140.000000 140.000000 140.000000
0.116121 68.514571 35.407429 1.735714
0.082982 9.024134 6.146080 0.818937
0.015000 29.860000 15.560000 1.000000
0.059250 64.227500 33.260000 1.000000
0.098000 69.525000 35.805000 1.500000
0.151250 73.945000 39.285000 2.000000
0.455000 91.820000 46.280000 3.000000

Random forest is an ensemble learning method composed of several decision trees. It contains multiple decision trees trained by Bagging, that is, Bootstrap Aggregating learning technology. When the samples to be classified are input, the final result is output by many decision trees by voting. This method solves the problem of performance bottlenecks of decision tree, and has good parallelism and scalability for high-dimensional data classification problems, and has good tolerance for noise and outliers.



### Feature selection

Heatmap is a good way to express the relationship between features, reflecting the importance and correlation of features. When the dimension of the data set is large, it will increase the calculation amount of the classifier and the running time. You can remove some of the insignificant features through the correlation analysis of the features to increase the effectiveness of the analysis task.



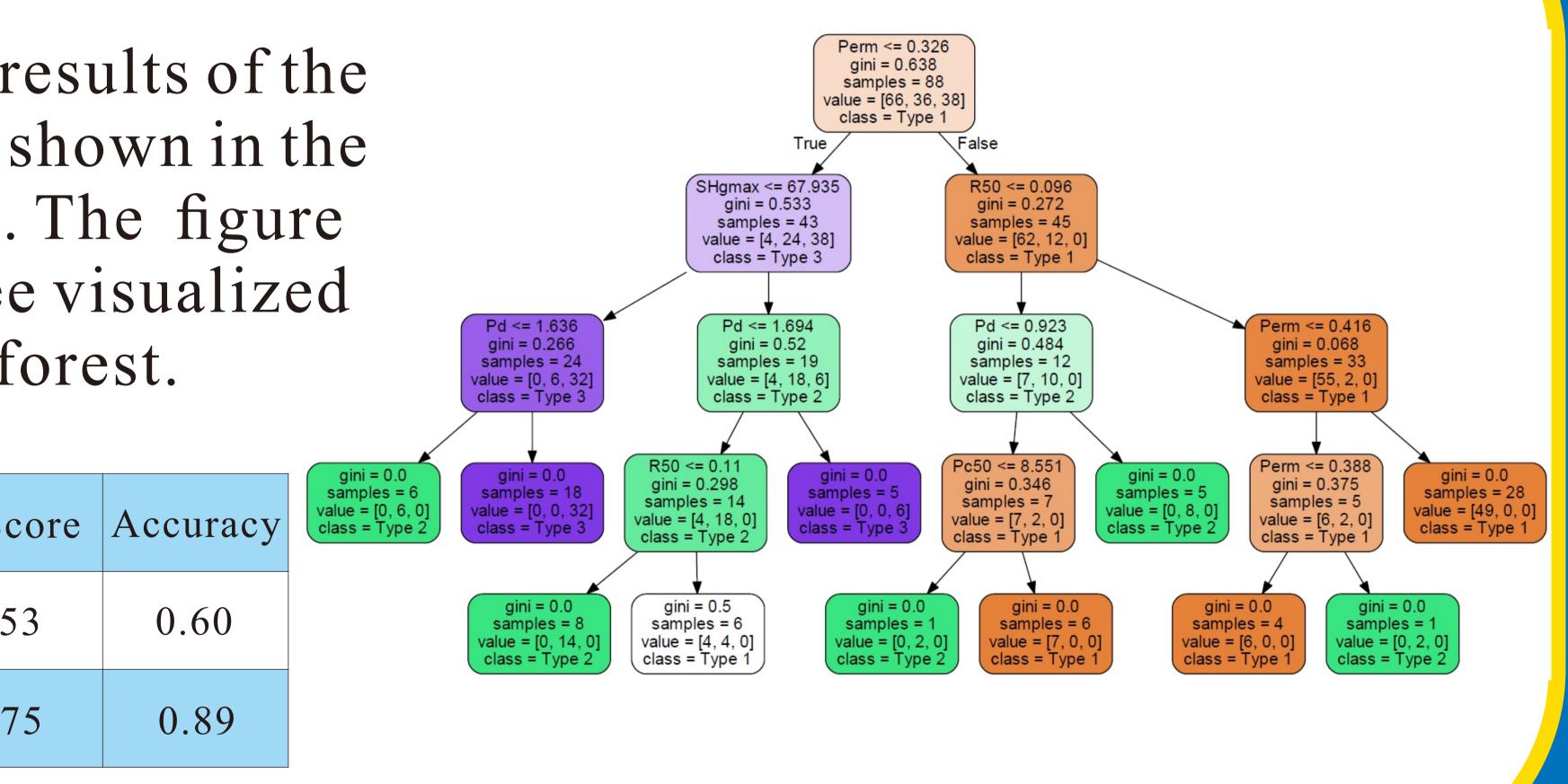
The prediction results of the two models are shown in the following table. The figure is a decision tree visualized from a random forest.

Model	F1 Sc
K-means	0.5
Random Forest	0.7

1.Based on K-means clustering and random forest analysis, the tight sandstone reservoirs in the study area were divided into three types. The prediction model achieved relatively good accuracy. 2. The low computational complexity, low complexity, and low subjective factor interference ensure the quantification and accuracy of the evaluation results. It may apply to the prediction of other parameters in the petroleum field.



#### Results



#### Conclusions