Part B

formal description

The performance of the IRT algorithm in part A is not satisfactory. We analyzed that the main reason is that a single decision tree is a high-variance model. We believe that the decision tree in part A overfits the training data, which makes the model's generalization ability poor.

Therefore, we decided to reduce the variance by averaging the predictions of multiple decision trees using a random forest. Each tree is trained on a different random subset of the training data. The results are aggregated at the end to make the model more robust by smoothing the data.

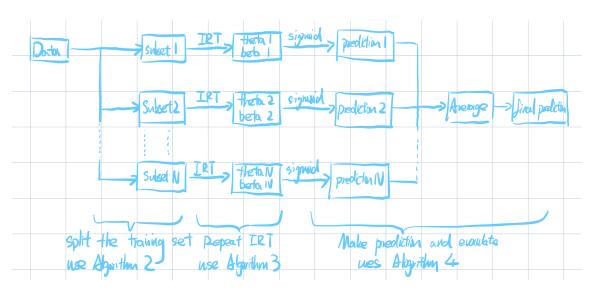
Algorithm Box

```
Result: train a single irt model
Output: thetas, betas
# regular irt algorithm, no changes needed
                            Algorithm 2: split the training data
Result: split the training data into subsets
Output: a subset of the training data
# randomly split the training data into subsets
                            Algorithm 3: irt ensemble algorithm
Result: train an ensemble of irt models
theta_lst = []
beta_lst = []
int repeat
for i in range(repeat):
    # split the training data
   train_data = split_data(data)
    # train a single irt model
    thetas, betas = irt(train_data)
    # store the thetas and betas in a list
    theta_lst.append(thetas)
    beta_lst.append(betas)
return theta_lst, beta_lst
                            Algorithm 4: evaluate the ensemble
Result: evaluate the ensemble of irt models
Output: the accuracy of the ensemble model
total_correct = 0
for i in range(len(data["user_id"])):
    prdict_lst = []
    for j in range(len(theta_lst)):
       # make prediction for each model and store it in a list
    # take the average of the predictions
# then calculate the accuracy
return total_correct / len(data["is_correct"])
```

By training multiple decision trees on different data subsets in the form of random forests,

the overfitting problem of the model can be reduced.

Idea Diagram



Comparison or Demonstration

For comparison, when using the single irt algorithm, we obtained the following statistics. We use this group of data as the baseline models:

```
Final Validation Accuracy: 0.7063223257126728
Final Test Accuracy: 0.707310189105278
```

When using a random forest consisting of two decision trees, we get the following statistics:

```
Ensemble Validation Accuracy: 0.7022297488004516
Ensemble Test Accuracy: 0.7044877222692634
```

When using a random forest consisting of three decision trees, we get the following statistics:

```
Ensemble Validation Accuracy: 0.7054755856618685
Ensemble Test Accuracy: 0.7044877222692634
```

When using a random forest consisting of four decision trees, we get the following statistics:

```
Ensemble Validation Accuracy: 0.705193338978267
Ensemble Test Accuracy: 0.703076488851256
```

After comparison, our model does not significantly improve the accuracy.

experiment to test our hypothesis

We use the accuracy of the training data set and the accuracy of the validation data set to determine whether the model has signs of overfitting. If the training accuracy is significantly higher than the validation accuracy, it means that the model is too sensitive to the training data set and has signs of overfitting. On the contrary, if the training accuracy and validation accuracy are close, it means that the model does not have overfitting.

The test accuracy of the original model is as follows:

Final train Accuracy: 0.7398744002257973

Obviously, the training accuracy of the original model is significantly higher than the validation accuracy, indicating that the original model may be overfitting.

The random forests consisting of 2, 3, and 4 decision trees have the following statistics:

Ensemble Training Accuracy: 0.7379339542760373

Ensemble Training Accuracy: 0.7391158622636184

Ensemble Training Accuracy: 0.7374576629974597

Unfortunately, Random Forest only slightly reduces the gap between training accuracy and validation accuracy, and the overfitting problem still exists.

Limitations

As mentioned above, implementing Random Forest algorithm did not improve our original model significantly. This could be due to the following reasons:

The given dataset is not large enough or diverse enough. Random Forest is known to perform well on large-scale datasets with many features. In our improved model, we split the dataset into many subsets and trained a Random Forest model on each subset. However, the dataset may not be large enough to benefit from this approach.

Hence, changing the dataset may improve the performance of the Random Forest algorithm.

In our bagging implementation, we assumes all predictions from the base models are equally important. This may not be the case in practice. Some models may perform better on certain types of data, and simply averaging their predictions may not be the best approach.

We can use a weighted average of the predictions to give more importance to the better performing models, but this requires tuning the weights, which can be time-consuming.

The Random Forest algorithm is computationally expensive. It requires a lot of time to tune the hyperparameters and train the model. We may not have been able to find the best hyperparameters in the time we had.

IRT model itself may be limited in predicting the student's performance. If we integrate more complex models in the ensemble, we may improve the performance of the model.