# **ML Intern Assignment Report**

#### 1. Introduction

This report summarizes the steps taken to process hyperspectral imaging data, apply dimensionality reduction, train a machine learning model, and evaluate its performance for predicting mycotoxin levels in corn samples.

## 2. Data Preprocessing

- Handling Missing Values: The dataset was checked for missing values, and none were found.
- Feature Selection: Removed the hsi id column as it was not relevant.
- **Normalization:** Applied standardization using StandardScaler to ensure all spectral features were on the same scale.

#### 3. Dimensionality Reduction

- **PCA:** Principal Component Analysis (PCA) was applied to reduce feature dimensions while retaining the maximum variance.
- **Results:** The first two principal components explained a significant portion of the variance, visualized through a scatter plot.

#### 4. Model Training & Evaluation

- Model Used: Random Forest Regressor was chosen for its robustness in handling high-dimensional data.
- Hyperparameters: Used default settings with 100 trees.
- Performance Metrics:
  - Mean Absolute Error (MAE): Computed and reported in script
  - Root Mean Squared Error (RMSE): Computed and reported in script
  - R<sup>2</sup> Score: Computed and reported in script
- Visualization: Scatter plot of actual vs. predicted DON concentration was generated.

## 5. Key Findings & Suggestions for Improvement

- The model performed reasonably well with Random Forest.
- Adding more advanced models like XGBoost or deep learning (CNN/LSTM) could further improve accuracy.
- Hyperparameter tuning (GridSearch or RandomSearch) could optimize model performance.
- Exploring t-SNE for better visualization of data clustering might provide deeper insights.

#### 6. Conclusion

This project successfully demonstrated preprocessing, dimensionality reduction, and ML-based regression for mycotoxin prediction. Future improvements could focus on deep learning approaches and enhanced feature engineering.