Code Explanation

1. Target and Type of regression

The dataset file path and metadata such as the target variable and prediction type are extracted from algoparams_from_ui.json file.

- **af** holds the full dataset.
- target variable defines the column to be predicted.
- prediction_type specifies whether the task is regression or classification.

2. Custom Feature Handling

I built a custom transformer called Custom_Feature_Handling to clean and prepare both numerical and text features — all based on instructions from a JSON configuration.

Handling Numerical Features

For each numerical column in the dataset, we check:

- Is the column selected for modelling?
- Are there missing values? If yes, we fill them using:
 - o The average (mean) of the column, or
 - o A **custom value** defined in the config file.

This makes sure that model doesn't break because of missing values and keeps things consistent.

Handling Text Features

Text columns are handled slightly different:

- If there's any missing text, fill it with "missing text" to avoid errors.
- If the configuration says to use "Tokenize and hash":
 - We apply a hashing trick to turn the text into a fixed number of numerical columns.

This step ensures that text data can be used in the model, without adding unnecessary complexity.

3. Custom Custom Feature Reduction

We created a custom class called Custom_Feature_Reduction that reduces dataset's size by keeping only the most relevant columns.

And the best part?

It adapts its strategy based on what's defined in the configuration JSON.

Modes Strategies:

Depending on what's set in the config, the transformer can use **one of four strategies**:

1. Correlation with Target

- Measures how strongly each feature relates to the target variable.
- Keeps the **top** K features that show the strongest correlation.

2. Tree-Based Importance

- Trains a **Random Forest** to find which features are most important in making predictions.
- Ranks features by their importance scores from the forest.
- Selects the **top K** features that contribute most to accuracy.

3. Principal Component Analysis (PCA)

- PCA compresses many correlated features into a few uncorrelated components.
- It's like blending information from several features into fewer new ones.
- Keeps only the top **K** components that explain the most variance.

4. No Reduction

• If the config says "No Reduction", the transformer keeps all columns (or a fixed number of them).

4. Hyper parameter tuning (GridSearchCV)

Most important step is **choosing the best model** to make predictions. But different models have different strengths — and they need the **right hyperparameters** to perform well.

Selects, tunes and evaluates models using GridSearchCV

Step 1: MODEL for tuning() Mapping of models to hyperparameters

If I want to use model X, then these are the hyperparameters I should try tuning and here's how to extract those tuning ranges from the config data.

MODEL for tuning declares what to tune and how to fetch the range definition from the input config.

The following regression and classification models are tuned:

• Regression:

- o LinearRegression, Ridge, Lasso, ElasticNet
- o RandomForestRegressor, DecisionTreeRegressor
- o GradientBoostingRegressor, ExtraTreesRegressor, XGBRegressor
- o MLPRegressor (neural network)

• Classification:

- o LogisticRegression, RandomForestClassifier, DecisionTreeClassifier
- o GradientBoostingClassifier, ExtraTreesClassifier, XGBClassifier
- o SVM, KNeighborsClassifier, MLPClassifier, SGDClassifier

Step 2: building pmt grid() Creates the actual parameter grid

This function interprets the hyperparameter definitions from MODEL_for_tuning and builds the final parameter grid that will be passed to GridSearchCV.

building_pmt_grid converts abstract definitions into real values to form a concrete pmt_grid

Step 3: Custom ModelSelection With GridSearch

This project automates that process through a custom scikit-learn-compatible class called ModelSelection With GridSearch.

It is designed to **choose, configure, and fine-tune** machine learning models using parameters provided in a JSON configuration file.

What This Component Does

The ModelSelection_With_GridSearch class acts as the final estimator in the pipeline. It performs the following steps programmatically:

- 1. Parses the JSON configuration to determine:
 - o Whether the task is **regression** or **classification**
 - o Which models are selected ("is selected": true)
 - o What parameter ranges should be explored during hyperparameter tuning
- 2. Builds model instances using scikit-learn classes like RandomForestRegressor, LogisticRegression, SVC, MLPClassifier, etc., depending on the type of task.
- 3. Creates parameter grids from the JSON file using keys like range (min, max) or range float (min, max, step) that are parsed into actual hyperparameter value ranges.
- 4. Runs GridSearchCV on each selected model:
 - o Uses **5-fold cross-validation** (K-fold) to evaluate each hyperparameter combination
 - o Identifies the best-performing combination for each model
 - o Stores the model and its best parameters for later use
- 5. **Returns a list of best estimators**, one for each selected model, sorted by their validation performance.

Each of these models can be **turned on/off** and **customized** just by changing values in the JSON file — making the system highly modular and reusable.

5. Execute pipeline

The pipeline is constructed with three core components:

```
Custom_Feature_Handling
Custom_Feature_Reduction
ModelSelection With GridSearch
```

Output

At the end of this step:

- We get the **best version of each selected model**, tuned and validated.
- These models are stored in self.selected_models_ as a list of tuples: (model_name, best_estimator_).

This design makes it **extremely reusable and configurable**. Just change the JSON file, and you can handle a completely new dataset, apply different reduction strategies, or try different model/hyperparameter combinations — **without changing the code**.