




Integrated Project - 2

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# Module	
# Time spent, hours	
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Task statement

Prepare a prototype of a machine learning model for Zyfra. The company develops efficiency solutions for heavy industry.

The model should predict the amount of gold recovered from gold ore. You have the data on extraction and purification.

The model will help to optimize the production and eliminate unprofitable parameters.

You need to:

1. Prepare the data;
2. Perform data analysis;
3. Develop and train a model.

To complete the project, you may want to use documentation from *pandas*, *matplotlib*, and *sklearn*.

The next lesson is about the ore purification process. You will pick the information that is important for the model development.

Technological Process

How is gold extracted from ore? Let's look into the process stages.

Mined ore undergoes primary processing to get the ore mixture or rougher feed, which is the raw material for flotation (also known as the rougher process). After flotation, the material is sent to two-stage purification.

Let's break down the process:

1. Flotation

Gold ore mixture is fed into the float banks to obtain rougher Au concentrate and rougher tails (product residues with a low concentration of valuable metals).

The stability of this process is affected by the volatile and non-optimal physicochemical state of the flotation pulp (a mixture of solid particles and liquid).

2. Purification

The rougher concentrate undergoes two stages of purification. After purification, we have the final concentrate and new tails.

Data description

Technological process

- *Rougher feed* — raw material
- *Rougher additions* (or *reagent additions*) — flotation reagents: *Xanthate*, *Sulphate*, *Depressant*
 - *Xanthate* — promoter or flotation activator;
 - *Sulphate* — sodium sulphide for this particular process;
 - *Depressant* — sodium silicate.
- *Rougher process* — flotation
- *Rougher tails* — product residues
- *Float banks* — flotation unit
- *Cleaner process* — purification
- *Rougher Au* — rougher gold concentrate

- *Final Au* — final gold concentrate

Parameters of stages

- *air amount* — volume of air
- *fluid levels*
- *feed size* — feed particle size
- *feed rate*

Feature naming

Here's how you name the features:

`[stage].[parameter_type].[parameter_name]`

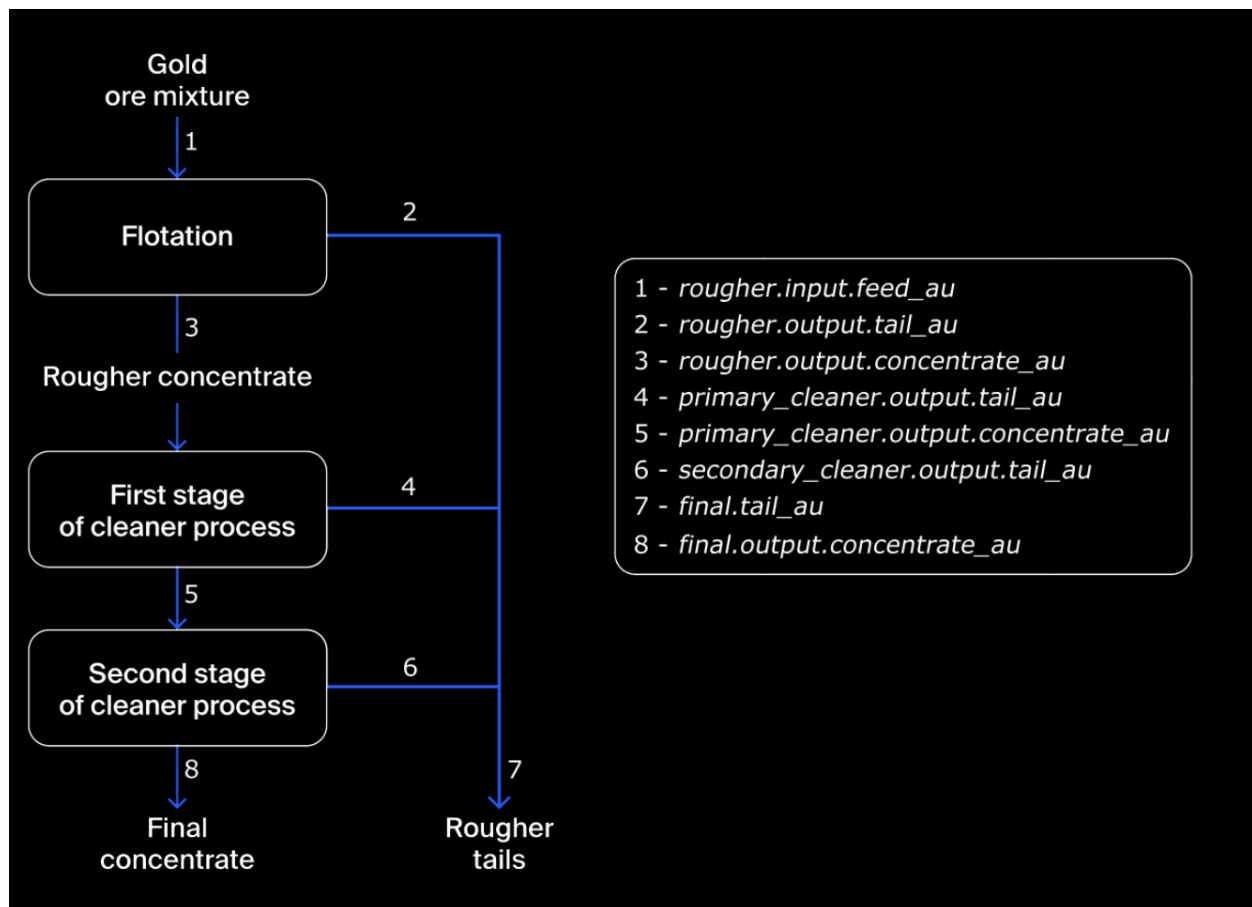
Example: `rougher.input.feed_ag`

Possible values for `[stage]`:

- *rougher* — flotation
- *primary_cleaner* — primary purification
- *secondary_cleaner* — secondary purification
- *final* — final characteristics

Possible values for `[parameter_type]`:

- *input* — raw material parameters
- *output* — product parameters
- *state* — parameters characterizing the current state of the stage
- *calculation* — calculation characteristics



Recovery calculation

You need to simulate the process of recovering gold from gold ore.

Use the following formula to simulate the recovery process:

$$\text{Recovery} = \frac{C \times (F - T)}{F \times (C - T)} \times 100\%$$

where:

- *C* — share of gold in the concentrate right after flotation (for finding the rougher concentrate recovery)/after purification (for finding the final concentrate recovery)

- F — share of gold in the feed before flotation (for finding the rougher concentrate recovery)/in the concentrate right after flotation (for finding the final concentrate recovery)
- T — share of gold in the rougher tails right after flotation (for finding the rougher concentrate recovery)/after purification (for finding the final concentrate recovery)

To predict the coefficient, you need to find the share of gold in the concentrate and the tails. Note that both final and rougher concentrates matter.

Evaluation metric

To solve the problem, we will need a new metric. It is called **sMAPE**, symmetric Mean Absolute Percentage Error.

It is similar to MAE, but is expressed in relative values instead of absolute ones. Why is it symmetrical? It equally takes into account the scale of both the target and the prediction.

Here's how *sMAPE* is calculated:

$$\text{sMAPE} = \frac{1}{N} \sum_{i=1}^N \frac{|y_i - \hat{y}_i|}{(|y_i| + |\hat{y}_i|) / 2} \times 100\%$$

Denotation:

$$y_i$$

- Value of target for the observation with the i index in the sample used to measure quality.

$$\hat{y}_i$$

- Value of prediction for the observation with the i index, for example, in the test sample.

$$N$$

- Number of observations in the sample.

$$\sum_{i=1}^N$$

- Summation over all observations of the sample (i takes values from 1 to N).

We need to predict two values:

1. rougher concentrate recovery `rougher.output.recovery`
2. final concentrate recovery `final.output.recovery`

The final metric includes the two values:

$$\text{Final sMAPE} = 25\% \times \text{sMAPE(rougher)} + 75\% \times \text{sMAPE(final)}$$

Project description

The data is stored in three files:

- `gold_recovery_train.csv` — training dataset [download](#)
- `gold_recovery_test.csv` — test dataset [download](#)
- `gold_recovery_full.csv` — source dataset [download](#)

Data is indexed with the date and time of acquisition (`date` feature). Parameters that are next to each other in terms of time are often similar.

Some parameters are not available because they were measured and/or calculated much later. That's why, some of the features that are present in the training set may be absent from the test set. The test set also doesn't contain targets.

The source dataset contains the training and test sets with all the features.

You have the raw data that was only downloaded from the warehouse. Before building the model, check the correctness of the data. For that, use our instructions.

Project instructions

1. Prepare the data

1.1. Open the files and look into the data.

Path to files:

- `/datasets/gold_recovery_train.csv`
- `/datasets/gold_recovery_test.csv`
- `/datasets/gold_recovery_full.csv`

1.2. Check that recovery is calculated correctly. Using the training set, calculate recovery for the ``rougher.output.recovery``. Find the *MAE* between your calculations and the feature values. Provide findings.

1.3. Analyze the features not available in the test set. What are these parameters? What is their type?

1.4. Perform data preprocessing.

2. Analyze the data

2.1. Take note of how the concentrations of metals (*Au*, *Ag*, *Pb*) change depending on the purification stage.

2.2. Compare the feed particle size distributions in the training set and in the test set. If the distributions vary significantly, the model evaluation will be incorrect.

2.3. Consider the total concentrations of all substances at different stages: raw feed, rougher concentrate, and final concentrate. Do you notice any abnormal values in the total distribution? If you do, is it worth removing such values from both samples? Describe the findings and eliminate anomalies.

3. Build the model

3.1. Write a function to calculate the final *sMAPE* value.

3.2. Train different models. Evaluate them using cross-validation. Pick the best model and test it using the test sample. Provide findings.

Use these formulas for evaluation metrics:

$$sMAPE = \frac{1}{N} \sum_{i=1}^N \frac{|y_i - \hat{y}_i|}{(|y_i| + |\hat{y}_i|) / 2} \times 100\%$$

$$\text{Final } sMAPE = 25\% \times sMAPE(\text{rougher}) + 75\% \times sMAPE(\text{final})$$

Project evaluation

We've put together the evaluation criteria for the project. Read this carefully before moving on to the case.

Here's what the reviewers will look at when reviewing your project:

- Have you prepared and analyzed the data properly?

- What models have you developed?
- How did you check the model's quality?
- Have you followed all the steps of the instructions?
- Did you keep to the project structure and explain the steps performed?
- What are your findings?
- Have you kept the code neat and avoided code duplication?

You have your takeaway sheets and summaries for previous chapters, so you are good to go.

Good luck!