



# Kedro

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## Basics

Create and activate conda env (this is optional)

```
conda create --name my_kedro_env python==3.8.1
conda activate my_kedro_env
```

Install kedro (the latest version in pip)...

```
pip install kedro
```

... or install kedro (specific version as needed)

```
pip install kedro==0.17.5
```

Verify kedro installation

```
kedro info
```

Create new kedro project in cwd

```
kedro new
```

Create new kedro project in cwd with example project code

```
kedro new --starter=pandas-iris
```

Create new kedro project in cwd with pyspark configuration

```
kedro new --starter=pyspark
```

Initialise git repo in your new kedro project

```
cd path_to_your_project && git init
```

Install all project dependencies

```
kedro install
```

Run the kedro pipeline(s)

```
kedro run
```

Package your project, build .whl and .egg files

```
kedro package
```

Visualize your pipelines with kedro viz

```
pip install kedro-viz && kedro viz
```

## Constructs

Kedro is broken up into a few major building blocks:

### Node

Simply a wrapper for a Python function that defines optional inputs/outputs (data from the Data Catalog).

A Kedro Node must have at least 1 input or output.

### Pipeline

Organises the dependencies and execution order of nodes.

### Data Catalog

A registry of all the data sources that a project can utilise. It handles loading and saving of data for you. Define once here and reuse everywhere.

### Runner

An object that runs the pipeline. You can choose from:

[Sequential](#) – executes pipeline nodes one-by-one

[Parallel](#) – enables concurrency via multiprocessing

[Threaded](#) – uses threading for concurrent execution

### Parameters

One or many yml files which are used to store re-usable project configuration such as model tuning params or splits. Remove all hardcoding from your code by defining once here.

## Folder structure

```
my_project      # Kedro Project Parent dir
├── conf         # Project configuration files
├── data         # Local project data*
├── docs         # Documentation
├── logs         # Output logs*
├── notebooks   # Jupyter notebooks
├── README.md    # Project README
├── setup.cfg    # Config for `pytest` and `isort`
└── src         # Project source code

* = Not committed to version control.
```

[conf/](#)

[base](#) – project-specific settings to share with others.

[local](#) – settings that should not be shared, e.g., sensitive credentials

[data/](#)

[raw](#) – raw unprocessed data should live here, unchanged

[intermediate](#) – optional data transformations applied to raw

[primary](#) – cleaned, transformed and wrangled data

[feature](#) – primary data grouped, and with features added

[model\\_input](#) – specific feature data for a given analysis

[models](#) – stored, serialised pre-trained model(s)

[model\\_output](#) – results generated by the model(s)

[reporting](#) – used to drive dashboards/views

## Handling dependencies

There are two places dependencies are defined (both inside src/), as per [pip compile](#) best practices:

[requirements.in](#) – define your project dependencies

[requirements.txt](#) – generated file of pinned dependencies

For a new project, without a *requirements.in* file:

Add your dependencies to *requirements.txt* and generate *requirements.in*. After this, update only *requirements.in*.

```
kedro build-reqs
```

For a project, with a *requirements.in* file:

Add your dependencies to *requirements.in* only, then:

```
kedro build-reqs
```

To install all project dependencies from *requirements.txt*:

```
kedro install
```

## Executing nodes & pipelines

Execute the default pipeline, sequentially

```
kedro run
```

Execute the default pipeline, with multiprocessing. Note: doesn't work with Spark data sets - use ThreadRunner.

```
kedro run --runner=ParallelRunner
```

Execute the default pipeline, with multithreading

```
kedro run --runner=ThreadRunner
```

Execute the pipeline asynchronously

```
kedro run --async
```

View all pipelines available

```
kedro pipeline list
```

Run a pipeline by name

```
kedro run --pipeline=my_pipeline
```

Run tagged nodes/pipeline

```
kedro run --tag=my_tag
```

Run a node by name

```
kedro run --node=my_node_name
```



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## Jupyter and ipython

Run all of these commands from Kedro project root.

Use Kedro within ipython

```
kedro ipython
```

Use Kedro within jupyter notebook

```
kedro jupyter notebook
```

Use Kedro within jupyter lab

```
kedro jupyter lab
```

From here, you have access to Kedro's 'context' and 'session' objects.

Load or save item from the Data Catalog

```
df = context.catalog.load("my_data")
context.catalog.save("catalog_item_name", my_df)
```

Access project parameters from parameters.yml

```
parameters = context.params # type: Dict
parameters["test_train_split"]
```

Execute the Pipeline in a notebook cell - see [run options](#)

```
session.run()
session.run(pipeline_name='my_pipeline')
```

Reload kedro variables (e.g. if you modify the catalog)

```
%reload_kedro
```

Load kedro variables into a standard ipython session

```
%load_ext kedro.extras.extensions.ipython
%reload_kedro <path_to_project_root>
```

...or

```
%load_ext kedro.extras.extensions.ipython
%init_kedro <path_to_project_root>
%reload_kedro
```

## Data catalog code

Data Catalog code lives inside one or many catalog.yml files.

```
companies:
  type: pandas.ExcelDataSet
  filepath: data/01_raw/companies.xlsx

reviews:
  type: pandas.CSVDataSet
  filepath: data/01_raw/reviews.csv
```

● **Unique name** of the catalog entry

● **Location on disk**, inside the data folder. Note the layer (01\_raw)

● **Type of data.** Here we're using a Pandas CSV dataset. There are many [out of the box](#) data formats supported including Excel, S3, Azure, Spark, SQL Databases.

## Node & pipeline code

Python node code typically lives within: src/<your\_project>/pipelines/<your\_pipeline>/nodes.py

Pipeline definition code, as below, lives within: src/<your\_project>/pipelines/<your\_pipeline>/pipeline.py

```
from kedro.pipeline import Pipeline, node
from .nodes import preprocess_companies, preprocess_shuttles

def create_pipeline(**kwargs):
    return Pipeline([
        node(
            func=preprocess_companies,
            inputs=["companies"],
            outputs="preprocessed_companies",
            name="preprocess_companies_node",
        ),
        node(
            func=preprocess_shuttles,
            inputs=["shuttles"],
            outputs="preprocessed_shuttles",
            name="preprocess_shuttles_node",
        ),
    ])
```

● **Import your node functions.**

● **Define a pipeline.** A pipeline consists of 1 or more nodes. You can add tags if you wish

● **Define a node.** A node consists of a function to call, inputs, outputs, a name and optional tags. Must have 1 input or output

● **Define which function to call.** A function is simply a python function

● **Define optional inputs.** Inputs link to data catalog items, parameters or memory datasets. Can be a single item, a list or a dict.

● **Define optional outputs.** As per inputs above. If you simply enter a string that doesn't exist in params or catalog, it will be saved to an in-memory dataset instead.

Be sure to update the project's pipeline in src/<your\_project>/pipeline\_registry.py so kedro knows to run your pipeline

```
from typing import Dict
from kedro.pipeline import Pipeline
from kedro_tutorial.pipelines import data_processing as dp

def register_pipelines() -> Dict[str, Pipeline]:
    data_processing_pipeline = dp.create_pipeline()

    return {
        "__default__": data_processing_pipeline,
        "dp": data_processing_pipeline,
    }
```

● **Set a default pipeline.** This is the pipeline which will execute when you type: kedro run

● **Define a pipeline.** The key of the dict here is used when you call a pipeline, e.g., kedro run --pipeline=dp