

Remote Landfill Atmospheric Gas Monitoring Tools - How to Guide

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1. Introduction

Methane (CH₄), a greenhouse gas with a warming potential 28 times greater than carbon dioxide over a 100-year period, has seen its atmospheric concentration increase over 250% since the industrial revolution. Despite CH₄'s shorter atmospheric lifespan, it still contributes to at least a quarter of anthropogenic warming (Pandey et al., 2023; Vigano et al., 2008).

Mismanaged landfills can produce significant CH₄ emissions (Ferronato et al., 2017) and the EU Landfill Directive of 1999 mandates the capture or flaring of CH₄ produced by organic waste decomposition (European Union, 1999; Themelis & Ulloa, 2007). Spain, which sent 11.5 million tonnes of waste to landfills in 2017 (European Environment Agency, 2022), has experienced large emission events despite landfill management on par with other EU countries (Castillo-Giménez et al., 2019; European Space Agency, 2021).

Satellite missions have in recent years improved their CH₄ measurement capabilities, offering advantages over ground-based detectors (Parker et al., 2011). PreZero, a waste management company operating 23 landfills in Spain, currently relies on ground-based detectors in a one detector per hectare grid. The tools outlined in this guide have been created while consulting with them. They expressed interest in other gasses beyond CH₄ so these have been included where possible (Aguasca, 2024; Hidalgo, 2024; Salami, 2024).

Three Python tools were created to display atmospheric gas data from Sentinel-5 and CH₄ data from Sentinel-2 satellites.

- **Sentinel-5 Atmospheric Gas Timeseries (S5-AGT):** A time series for each landfill location showing atmospheric gas concentrations of carbon monoxide (CO), formaldehyde (HCHO), nitro dioxide (NO₂), ozone (O₃), sulphur dioxide (SO₂) and methane (CH₄).
- **Sentinel-5 Atmospheric Gas Mapper (S5-AGM):** Produces a map of daily atmospheric gas concentrations of CO, HCHO, NO₂, O₃, SO₂ and CH₄ and provides zonal statistics for these at landfills.
- **Sentinel-2 Multi-Band-Single-Pass CH₄ Mapper (S2-MBSP):** Produces a high-resolution map that can show CH₄ point sources for selected landfills.

The tools can be downloaded or cloned from the following Github repository. This guide outlines their installation and use. The files for which can be found at the following URL:

[https://github.com/zelcon01/Landfill Atmospheric Gas Monitor Tools](https://github.com/zelcon01/Landfill_Atmospheric_Gas_Monitor_Tools)

2. Setup

Anaconda Navigator, containing the Python programming language and 'Conda', a package manager for creating shareable environments that tools like this can run on will be installed. This includes Jupyter Lab which will be used to run this tool's code.

2.1 Anaconda Navigator

To download Anaconda, navigate to <https://docs.anaconda.com/anaconda/install/> and follow the instructions of your operating system.

2.2. Creating a Conda Environment

In the Anaconda Navigator side bar, click the 'Environments' tab. You will see the installed packages (fig.1).

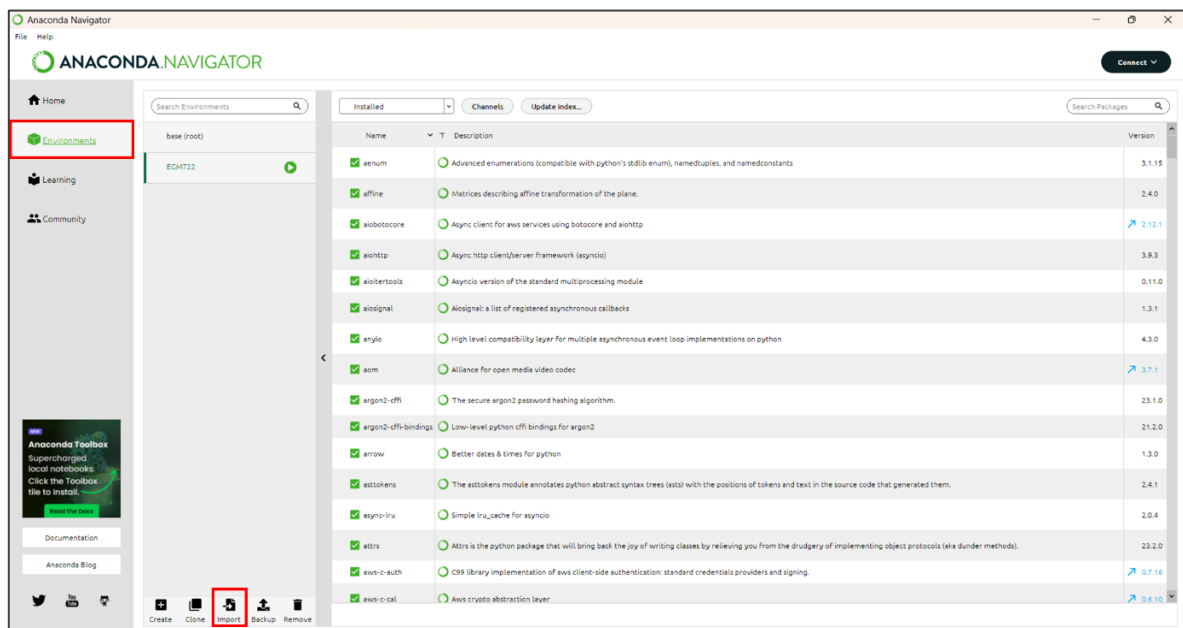


Figure 1: Environments tab of Anaconda Navigator with environments tab and import button highlighted in red.

Next click on the imports tab (fig.1) and select the file 'environment.yml' contained in the .zip file of the tool's download, choosing an appropriate name for the environment (fig.2).

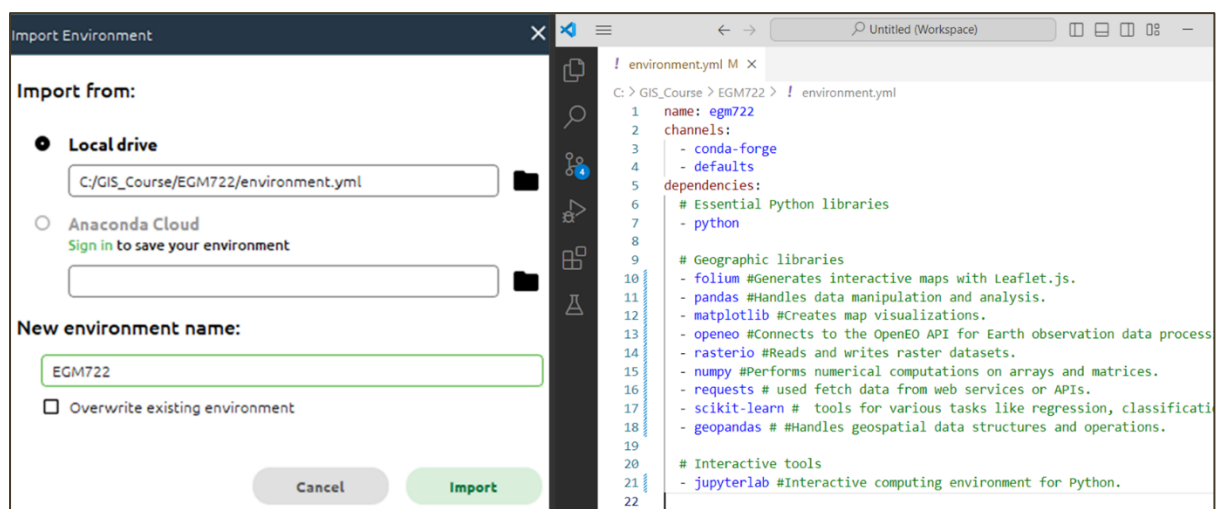


Figure 2: The import config box (left) and the contents of 'environment.yml' (right).

Click Import. The process of installing all the packages may take several minutes. Once finished you will be returned to the environments tab (fig.1)

Next click on the 'Home' tab in Anaconda Navigator's sidebar (fig.3).

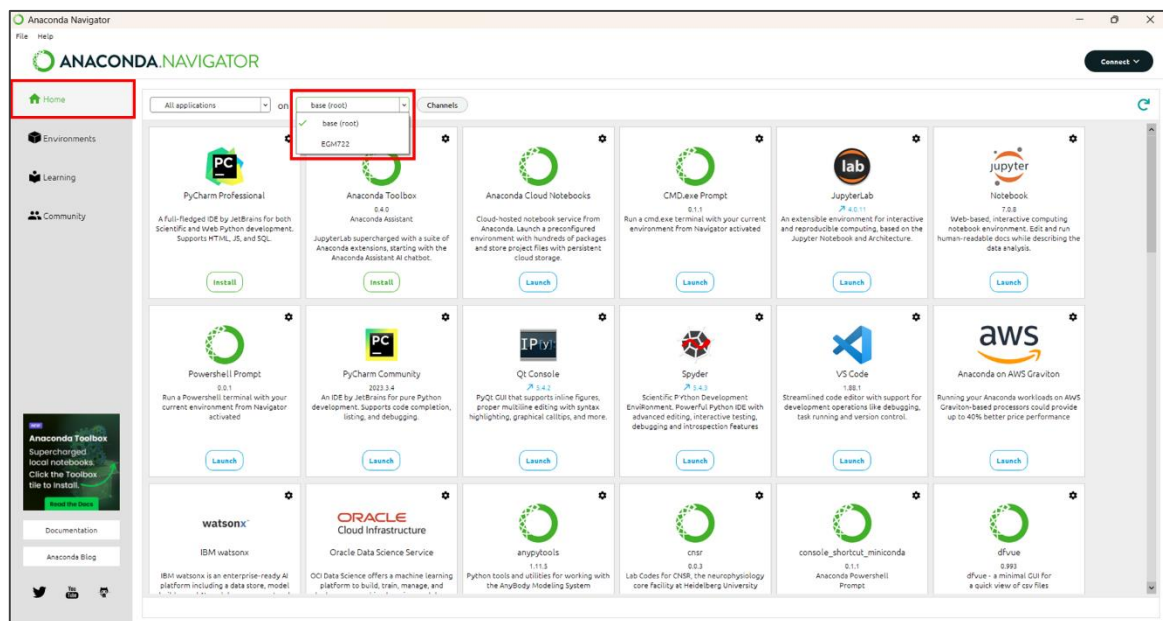


Figure 3: Anaconda Navigator with home tab and environment switching dropdown in red.

The dropdown highlighted in figure 3 should display two options, 'base (root)' and the name of your new environment (in figure 3 this is 'EGM722'). **Ensure your environment name is always selected here or the dependencies installed earlier will not be available to it.**

2.3 Setting up Jupyter Lab

A configuration file ('.config') needs to be created to change the settings used by Jupyter Lab by default. Launch the CMD.exe prompt (fig.4)

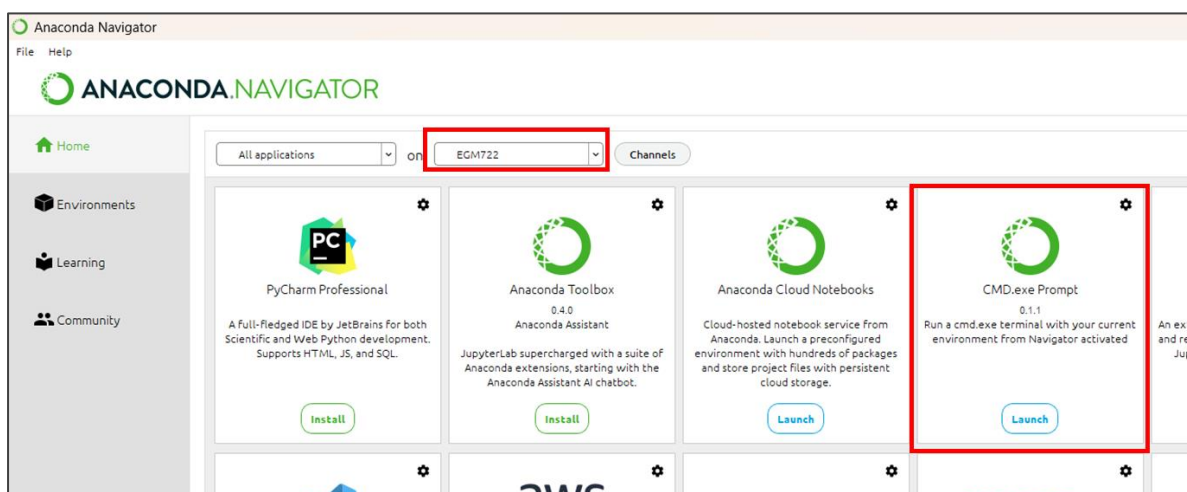


Figure 4: Highlighted locations of selected environment and CMD.exe Prompt

In the command prompt, enter the command:

```
jupyter lab --generate-config
```

This will create a new folder in your user directory called 'jupyter' containing a python script `jupyter_lab_config.py`. On Windows this is usually 'C:\Users\<your_username>'.

Jupyter lab will by default open in your user directory. Due to security restrictions, it is not possible to navigate to the parent directory of the launch location. So if Jupyter launches in 'C:\Users\RockyBalboa', it is not possible to move to 'C:\Users' or 'C:\EGM722'. If the directory you are keeping your data in is outside your user directory, you will need to change the default opening folder to your data directory.

This location is also where you should store the following files and folder:

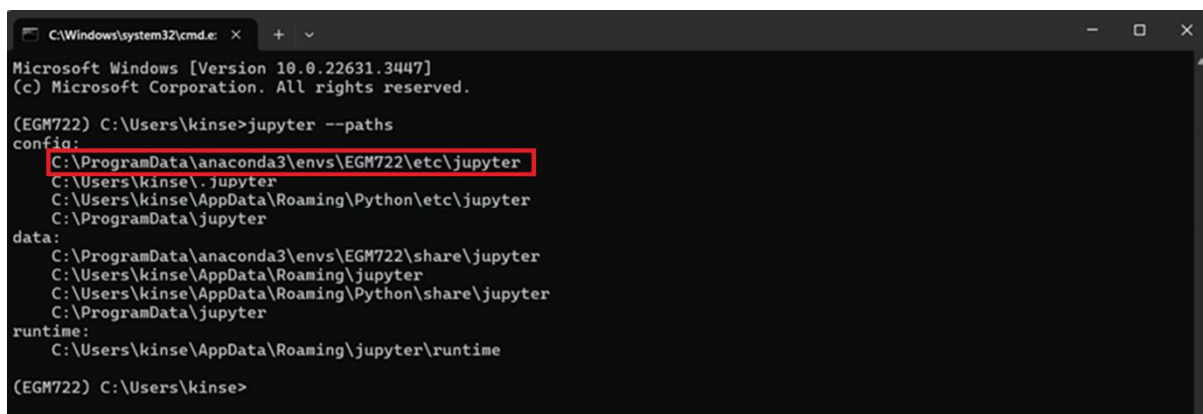
- Sentinel_5P_Atmospheric_Gas_Time_Series.ipynb
- Sentinel_5P_Atmospheric_Gas_Map.ipynb
- Sentinel_2_CH4_Multi-Band-Single-Pass.ipynb
- The folder "Data"

If your data directory is in your user directory, you should be able to click and navigate there using the interface of Jupyter Lab. If that is not the case, you will need to do the following:

Open an Anaconda Navigator CMD.exe prompt and type the following command:

```
jupyter --paths
```

This will show something like figure 5 although your file paths will be unique to you.



```

C:\Windows\system32\cmd.exe
Microsoft Windows [Version 10.0.22631.3447]
(c) Microsoft Corporation. All rights reserved.

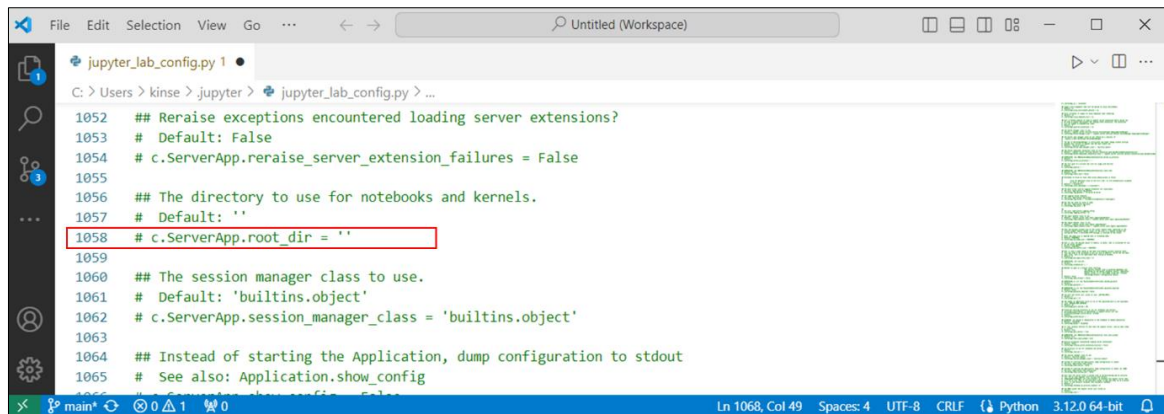
(EGM722) C:\Users\kinse>jupyter --paths
config:
  C:\ProgramData\anaconda3\envs\EGM722\etc\jupyter
  C:\Users\kinse\.jupyter
  C:\Users\kinse\AppData\Roaming\Python\etc\jupyter
  C:\ProgramData\jupyter
data:
  C:\ProgramData\anaconda3\envs\EGM722\share\jupyter
  C:\Users\kinse\AppData\Roaming\jupyter
  C:\Users\kinse\AppData\Roaming\Python\share\jupyter
  C:\ProgramData\jupyter
runtime:
  C:\Users\kinse\AppData\Roaming\jupyter\runtime

(EGM722) C:\Users\kinse>
  
```

Figure 5: results of 'jupyter --paths' command showing path used by environment highlighted in red.

The 'jupyter_lab_config.py' file mentioned earlier needs to be copy pasted into that folder.

Once 'jupyter_lab_config.py' file has been moved, open it in Notepad++, Visual Studio Code or if you don't have those, Notepad. Using the shortcut 'CTRL + F' type in the following line: 'c.ServerApp.root_dir' (without quotes) and you should find the section highlighted in figure 6.



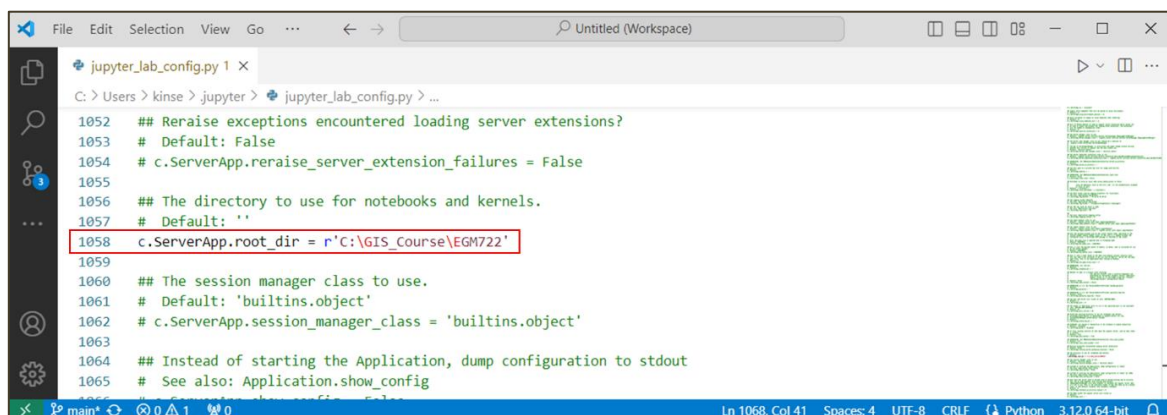
```

1052 ## Reraise exceptions encountered loading server extensions?
1053 # Default: False
1054 # c.ServerApp.reraise_server_extension_failures = False
1055
1056 ## The directory to use for notebooks and kernels.
1057 # Default: ''
1058 # c.ServerApp.root_dir = ''
1059
1060 ## The session manager class to use.
1061 # Default: 'builtins.object'
1062 # c.ServerApp.session_manager_class = 'builtins.object'
1063
1064 ## Instead of starting the Application, dump configuration to stdout
1065 # See also: Application.show_config

```

Figure 6: location of 'c.ServerApp.root_dir' in jupyter_lab_config.py

Remove the '#' and space from the start and add the path used by your environment between the quote marks, adding a 'r' beforehand (fig.7).



```

1052 ## Reraise exceptions encountered loading server extensions?
1053 # Default: False
1054 # c.ServerApp.reraise_server_extension_failures = False
1055
1056 ## The directory to use for notebooks and kernels.
1057 # Default: ''
1058 c.ServerApp.root_dir = r'c:\GIS_Course\EGM722'
1059
1060 ## The session manager class to use.
1061 # Default: 'builtins.object'
1062 # c.ServerApp.session_manager_class = 'builtins.object'
1063
1064 ## Instead of starting the Application, dump configuration to stdout
1065 # See also: Application.show_config

```

Figure 7: path to data directory added to jupyter_lab_config.py

Save and close this file and return to the Anaconda Navigator 'Home' tab. Launch Jupyter Lab and if you have followed the steps correctly, you should see that your data directory is automatically displayed (fig.8).

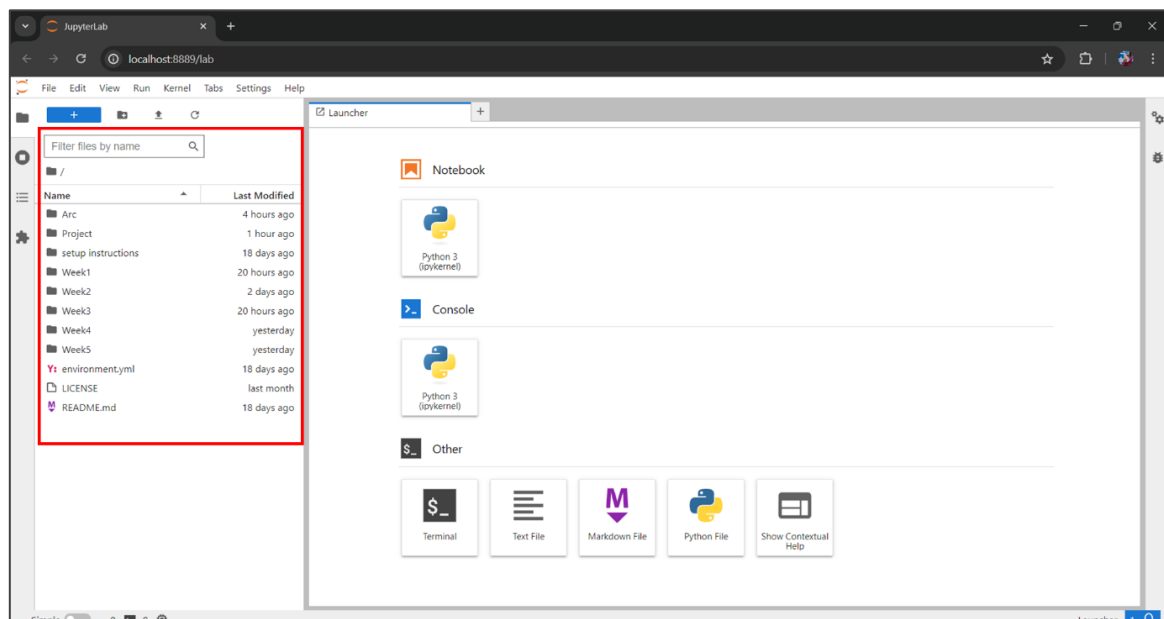


Figure 8: Jupyter Lab showing by default the data directory

2.4 OpenEO setup using Anaconda Navigator

OpenEO is an open-source API that allows access to the earth observation satellite missions run by the Copernicus program. These include the satellites used by this tool.

First search in the Anaconda Navigator environments tab for 'openeo'. Make sure that 'Not installed' is selected (fig.9). If the package appears here, click its tick box and select apply. If you can't see it here, please go to section 2.5.

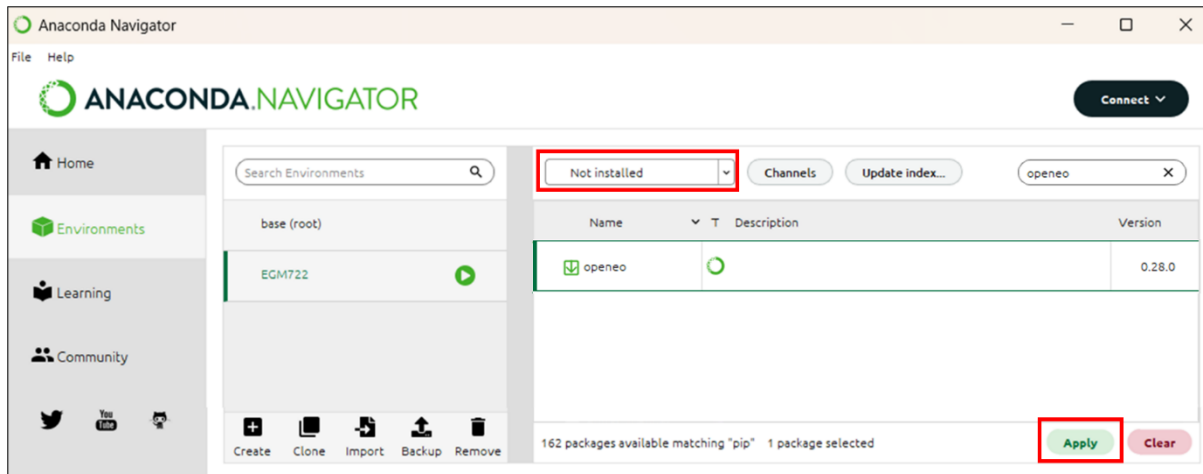


Figure 9: installing OpenEO from Anaconda Navigator.

Next you will be presented with the following screen (fig.10). Once this has finished processing the request. Simply click 'apply' to begin the installation.

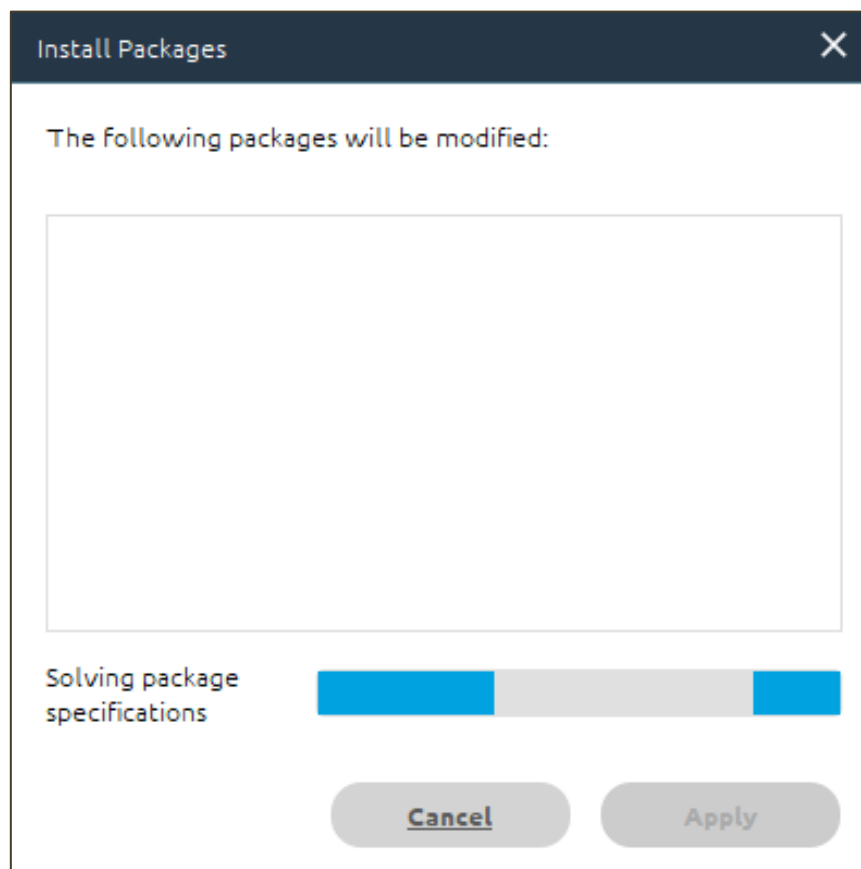


Figure 10: Anaconda Navigator package installer loading screen.

2.5 OpenEO setup using PyPi

If Anaconda Navigator cannot find OpenEO you can use PyPi, the official third-party software library for Python. Search for 'pip', selecting the appropriate tick-box and then clicking apply (fig.11), then clicking apply once the install packages prompt has finished loading (fig.10).

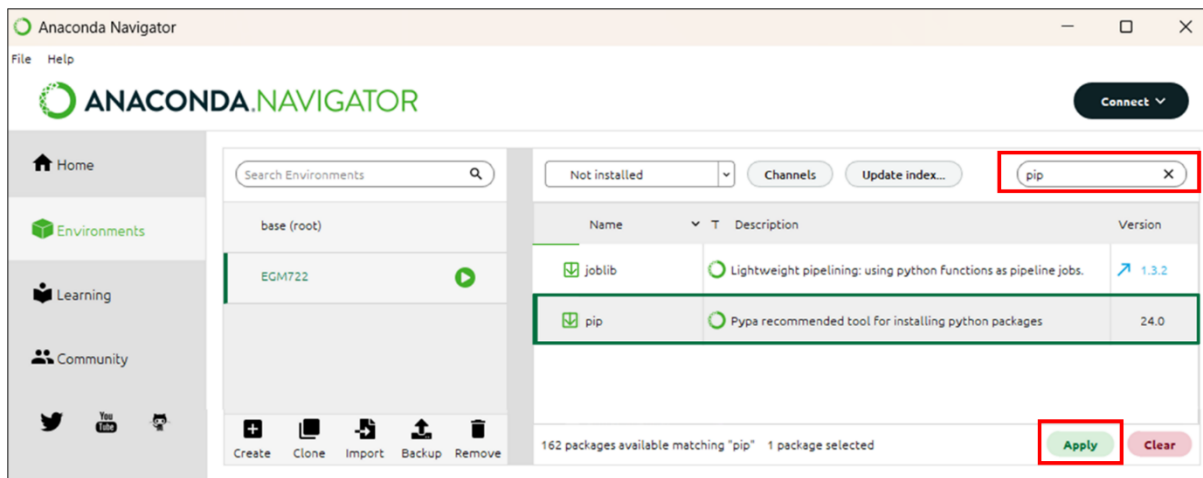


Figure 11: Installing pip via Anaconda Navigator

Open an Anaconda Navigator CMD.exe prompt and type the following command:

```
pip install openeo
```

Once the process has completed, you can close the CMD.exe prompt window.

2.6 Registering with Copernicus Data Space Ecosystem.

Accessing OpenEO requires an authentication. To do this, you need to complete a Copernicus Dataspace Registration. Go to <https://dataspace.copernicus.eu/> and click the green login button (fig.12)

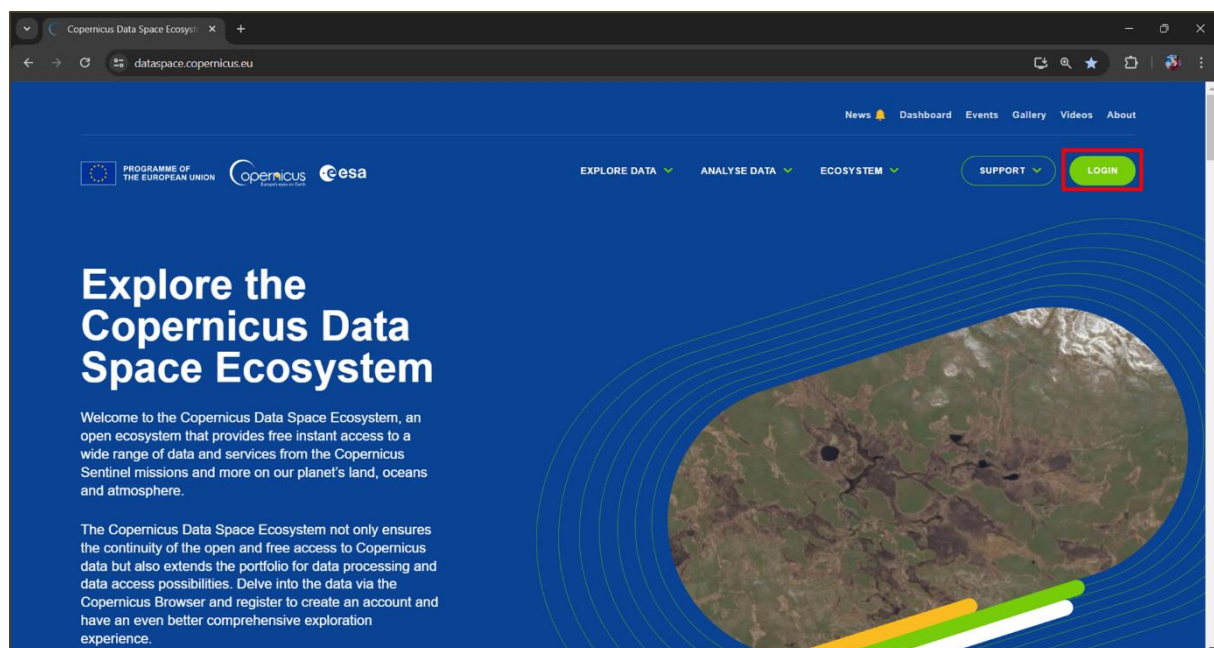


Figure 12: Copernicus Dataspace landing page with login button highlighted in red.

Next click the green 'register' button (fig.13):

Figure 13: Copernicus Dataspace sign in page.

On the following page, fill out the application form and then at the bottom click the green 'register' button (fig.14).

Figure 14: End of Copernicus registration page with register button highlighted in red.

Once registered, you will receive an email asking to verify your address. You can then log-in with your email and chosen password.

For any registration problems, email: help-login@dataspace.copernicus.eu

2.7 Running the tools in Jupyter-lab

Now that (almost) everything has been setup you can launch Jupyter Lab in the Anaconda Navigator (fig.15). Remember as always that your project environment (here 'EGM722') should be selected and not 'base (root)'.

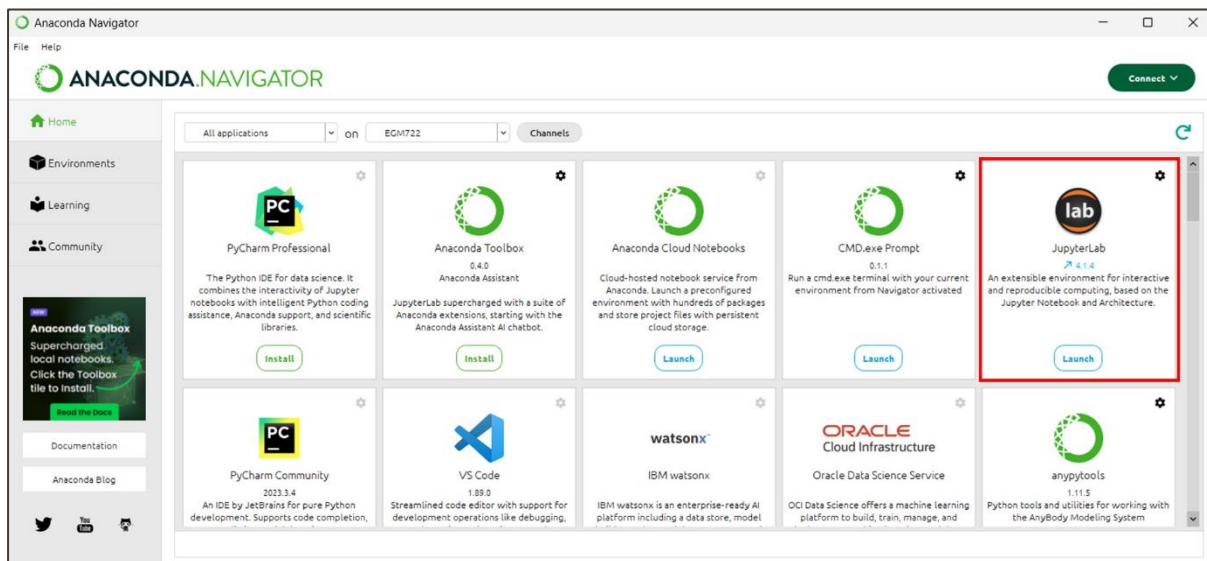


Figure 15: Location of Jupyter Lab in Anaconda Navigator highlighted in red.

Once Jupyter lab opens, you should see the three tools on the left (fig.16).

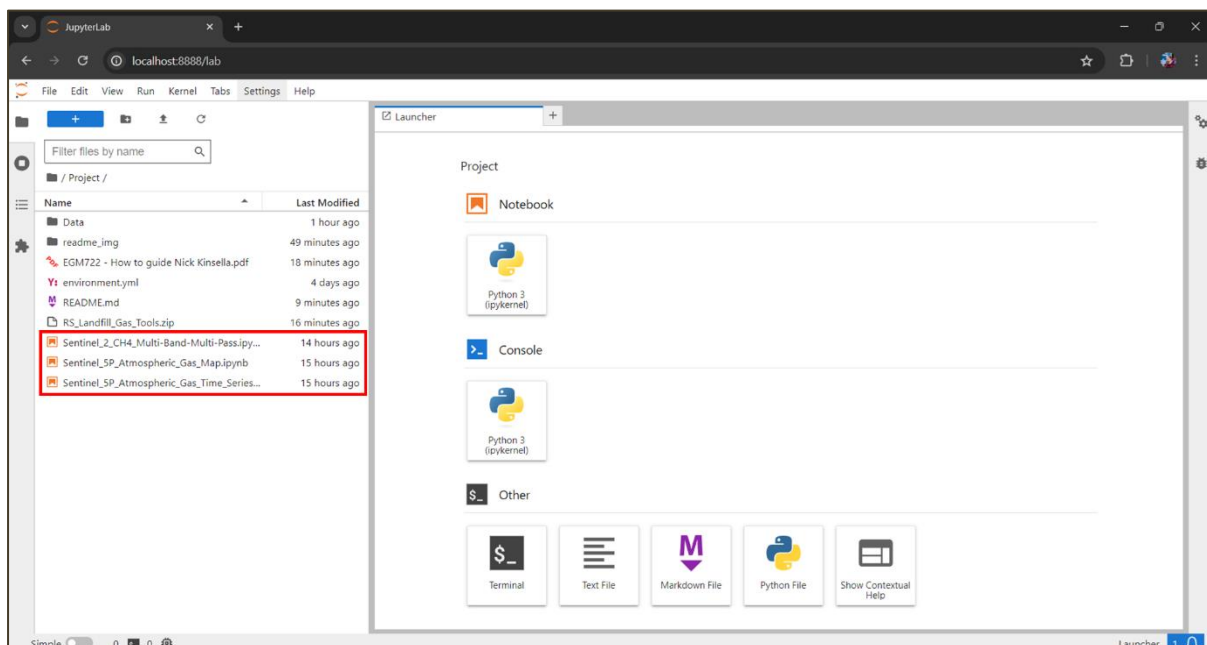


Figure 16: Location of tool scripts in Jupyter lab highlighted in red.

Click on one of the tools to open it. You can then follow the instructions, running the code by clicking the play button (fig.17).

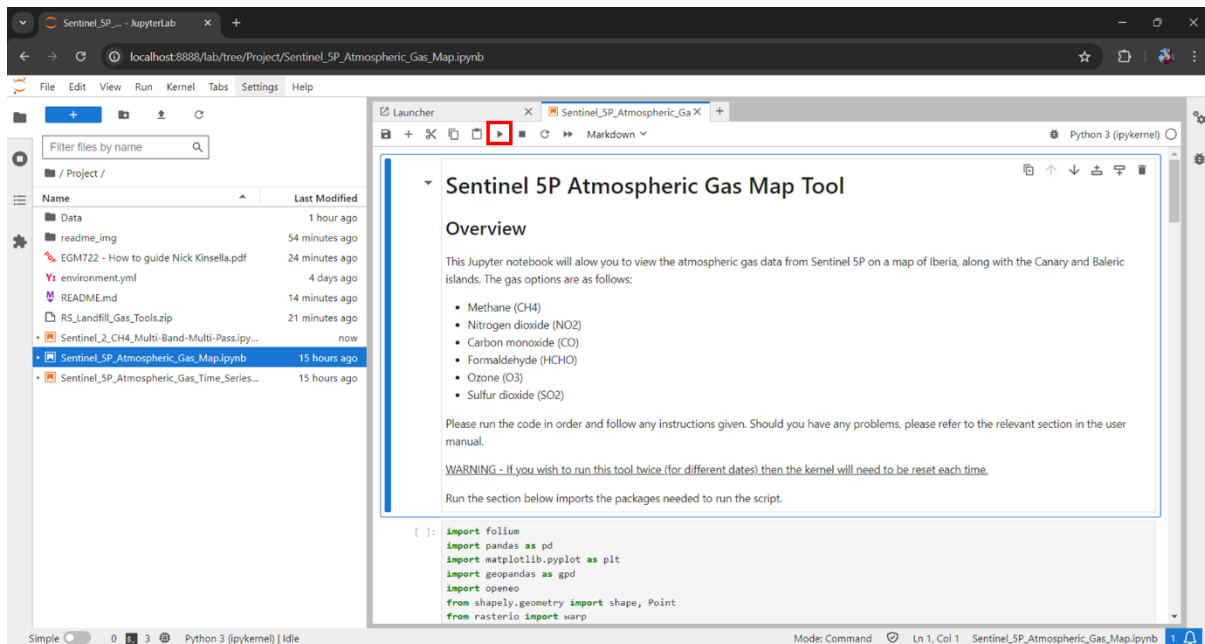


Figure 17: Location of the ‘play’ button which runs each of the code segments of the workbooks.

2.8 Authentication with OpenEO

The very first time one of the tools are run, the following section of code...

```
connection = openeo.connect(url="openeo.dataspace.copernicus.eu")
connection.authenticate_oidc()
```

... will provide you with a URL that will look something like this:

Visit https://auth.example.com/device?user_code=EAXD-RQXV to authenticate.

Copy this into your web browser and login using the Copernicus Data Space. Once this is complete, run tool’s Python script again and it will receive an authentication token, printing the message:

Authorized successfully.

In future you may be prompted with a new URL to create a new authentication token, whereby you should repeat the steps of this section.

3. Data and Dependencies

The datasets used by these tools are outlined in table 1.

Table 1: Datasets used by tools

Description	Data	Source	For use in tool(s)
PreZero Landfill Locations	Vector, point	PreZero International.	S5-AGT, S5-AGM,
PreZero Landfill Bounding	Long/Lat .csv file, point	PreZero Landfill Locations pre-processed.	S2-MBSP
Sentinel 5P (TROPOMI) total columns of CO, HCHO, NO ₂ , O ₃ , SO ₂ and CH ₄	Raster 4,518m x 3,552m	Copernicus Dataspace – OpenEO.	S5-AGT, S5-AGM
Sentinel 2 (MSI) bands 2, 3, 4 and 12	Raster 20m ²	Copernicus Dataspace – OpenEO.	S2-MBSP

The dependencies contained in the environment.yml file are outlined in table 2.

Table 2: dependencies required to use the tools.

Name	Version	Description	For use in tool(s)
Python	3.12.2	Programming language to run the code	S5-AGTS, S5-AGM & S2-MBSP
Jupyterlab	4.1.4	Computing environment for Python	
Folium	0.16.0	Generates interactive maps with Leaflet.js.	S5-AGTS & S5-AGM
Pandas	2.2.1	Data manipulation and analysis.	S5-AGTS, S5-AGM & S2-MBSP
Geopandas	0.14.3	Geospatial data manipulation and analysis.	S5-AGTS & S5-AGM
Matplotlib	3.8.3	Creates map visualizations	S5-AGTS, S5-AGM & S2-MBSP
Rasterio	1.3.9	Reads and edits raster datasets.	S5-AGM & S2-MBMP
Numpy	1.26.4	Performs numerical computations on arrays	S5-AGTS, S5-AGM & S2-MBSP
Requests	2.31.0	Used fetch data from web services or APIs.	S2-MBMP
Rasterstats	0.19.0	Linear regression for brightness correction	S2-MBMP
OpenEO	0.28.0	API for Earth observation data processing	S5-AGT, S5-AGM & S2-MBSP
Shapely	2.0.3	Used to manage points, lines and polygon data	S5-AGTS, S5-AGM

4. Methodology

4.1 Sentinel-5 Atmospheric Gas Timeseries:

This tool creates a timeseries graph of gas concentrations for selected landfills, identifying high concentration days using average concentration in a 10km buffer. It is based on a Python notebook from Copernicus OpenEO (2024), modified to use Sentinel 5 data and PreZero's landfill locations. Despite Sentinel 5's low spatial resolution, it provides near daily measurements of several gases, useful for studying transient atmospheric phenomena like methane emissions (Varon et al., 2021). The processing steps are presented in Figure 18 and the code thereafter.

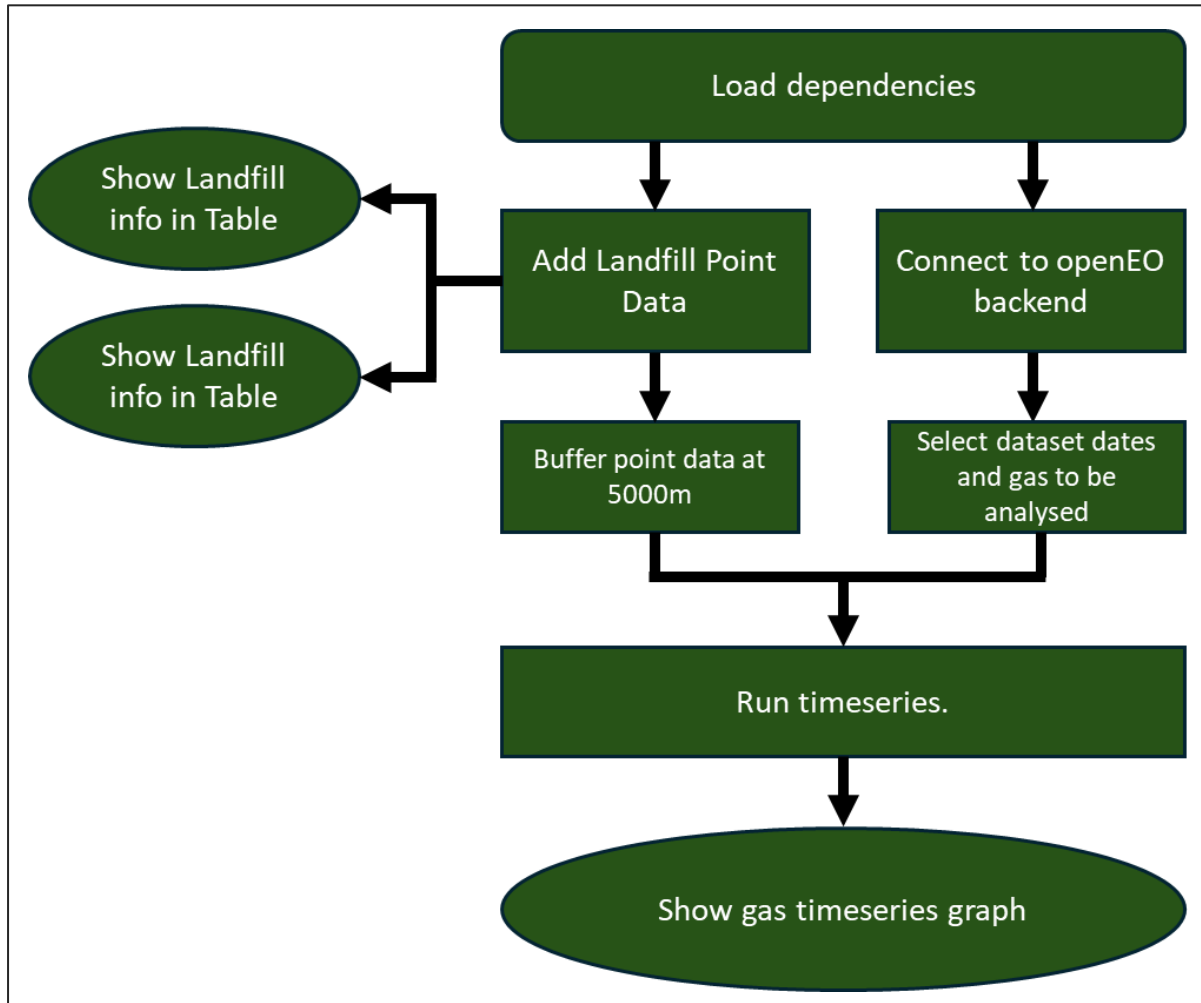


Figure 18: Flowchart of processes used for Gas timeseries analysis.

Code 1 shows the loaded dependencies.

```

import folium
from folium import plugins
import pandas
import matplotlib.pyplot as plt
import geopandas
import openeo
from shapely.geometry import shape
  
```

Code 1: Loading of dependencies for the code to run.

Code 2 connects to the OpenEO backend.

```
connection = openeo.connect(url="openeo.dataspace.copernicus.eu")
connection.authenticate_oidc()
```

Code 2: Connecting to OpenEO

Code 3 reads a file that contains data of the landfill location and then displays it.

```
landfills =
geopandas.read_file(r"C:\GIS_Course\Landfill_Atmospheric_Gas_Monitor_Tools\Data\
PZ_landfill_point4326.geojson")
```

Code 3: Displaying the contents of the landfill file for easy reference.

Code 4 shows the locations of the landfills on a map using folium. The user can click on a map pin to view the landfill's address. The map is centred on the loaded geometries by calculating the average x and y location of all the datapoints.

```
# This creates the map and centres it on the geometries.
centroids = landfills.geometry.centroid
centre = [centroids.y.mean(), centroids.x.mean()]
site_map = folium.Map(location=centre, zoom_start=5, control_scale=True)

# Adding the landfill locations to the map
for feature in landfills.iterfeatures():
    # Extract feature number from properties
    feature_number = feature['properties']['Landfill'] #this is for the
PreZero ones

    # Extract coordinates of the feature
    coordinates = feature['geometry']['coordinates']

    # Create a marker with label for each feature
    folium.Marker(location=[coordinates[1], coordinates[0]],
                  popup=f"Feature {feature_number}").add_to(site_map)

# Display the map
site_map
```

Code 4: Code for displaying the landfill points on a map for easy reference.

Code 5 adds a buffer of 10,000 metres to the landfill point data, and then formats it as a .geojson, as the OpenEO API requires it in that format.

```
# loading dataframe
landfill_10000m = landfills

# The dataset is projected in EPSG:4326 with its units in degrees. This
needs to be converted to CRS to EPSG:2062, which is in metres.
landfill_10000m = landfill_10000m.to_crs(epsg=2062)

# Now the dataframe is converted, a buffer of 5000m is added to each point
landfill_10000m['geometry'] = landfill_10000m.buffer(10000)
```

```
# Now the buffered data needs to be converted back to EPSG:4326 because the
Sentinel data is projected in EPSG:4326.
landfill_10000m = landfill_10000m.to_crs(epsg=4326)

# The time series analysis requires that A GeoJSON format file is used for
the analysed areas, so this bit produces a file suitable for that.
landfill_10000m_geojson = landfill_10000m.__geo_interface__
```

Code 5: Code for adding buffers to landfill point data

Code 6 selects the specific date range and gas type.

```
s5cube_timeseries = connection.load_collection(
    "SENTINEL_5P_L2",
    temporal_extent=["2021-08-01", "2021-10-31"], # format YYYY-MM-DD
    bands=["CH4"], # Gas options 'CO', 'HCHO', 'NO2', 'O3', 'SO2', 'CH4'
)
```

Code 6: Selecting date for time series and gas to be monitored.

Code 7 runs the time series analysis and saves the results as a .csv file. This process takes around 5 minutes per month selected.

```
timeseries =
s5cube_timeseries.aggregate_spatial(geometries=landfill_5000m_geojson,
reducer="mean")

#This saves the results as a .CSV file which can be viewed in Microsoft
Excel or a similar package. It will be saved in the indicated location.
job = timeseries.execute_batch(out_format="CSV", title="Gas timeseries")

job.get_results().download_file("Gas_Timeseries_results/Gas_timeseries.csv")

pd.read_csv("Gas_Timeseries_results/Gas_timeseries.csv", index_col=0)
```

Code 7: Running the data collection for the time series.

Code 8 plots the data on a timeseries graph using much of the script taken from the NDVI OpenEO notebook (OpenEO, 2024). Modifications include scaling the value of the Y axis to whatever data is loaded and allowing the user to select which landfills they want to view on the graph.


```

def plot_timeseries(filename, selected_landfill_ids=None, figsize=(15, 7)):
    df = pandas.read_csv(filename, index_col=0)
    df.index = pandas.to_datetime(df.index)
    df = df.sort_index()

    fig, ax = plt.subplots(figsize=figsize, dpi=90)

    if selected_landfill_ids:
        df_selected = df[df['feature_index'].isin(selected_landfill_ids)]
        for landfill_id, group in df_selected.groupby("feature_index"):
            group["avg(band_0)"].plot(marker="o", label=f"Landfill
{landfill_id}", ax=ax)
    else:
        df.groupby("feature_index")["avg(band_0)"].plot(marker="o", ax=ax)

    ax.set_title(filename.split("/")[-1])
    ax.set_ylabel("Parts per billion for CH4 or mol/m2 for all other gases")

    ymin = df["avg(band_0)"].min()
    ymax = df["avg(band_0)"].max()
    ymin_with_margin = ymin - 0.1 * (ymax - ymin)
    ymax_with_margin = ymax + 0.1 * (ymax - ymin)
    ax.set_ylim(ymin_with_margin, ymax_with_margin)

    ax.legend(title="Landfill id", loc='upper left', bbox_to_anchor=(1.02,
1), ncol=2)
    ax.xaxis.set_major_locator(plt.MaxNLocator(30))
    ax.grid(True)

# Change the selected landfill ids for the ones you want to view.
plot_timeseries("Gas_Timeseries_results/Gas_timeseries.csv",

                selected_landfill_ids=[6, 14, 23])

```

Code 8: For plotting the time series data on a graph

4.2 Sentinel-5 Atmospheric Gas Concentration Mapper:

The tool makes a map of atmospheric gas concentrations over the Iberian peninsula, the Balearic and Canary islands, with marked locations for where PreZero has its landfills, allowing the user to continue their analysis. Figure 19 provides an overview of the processing steps.

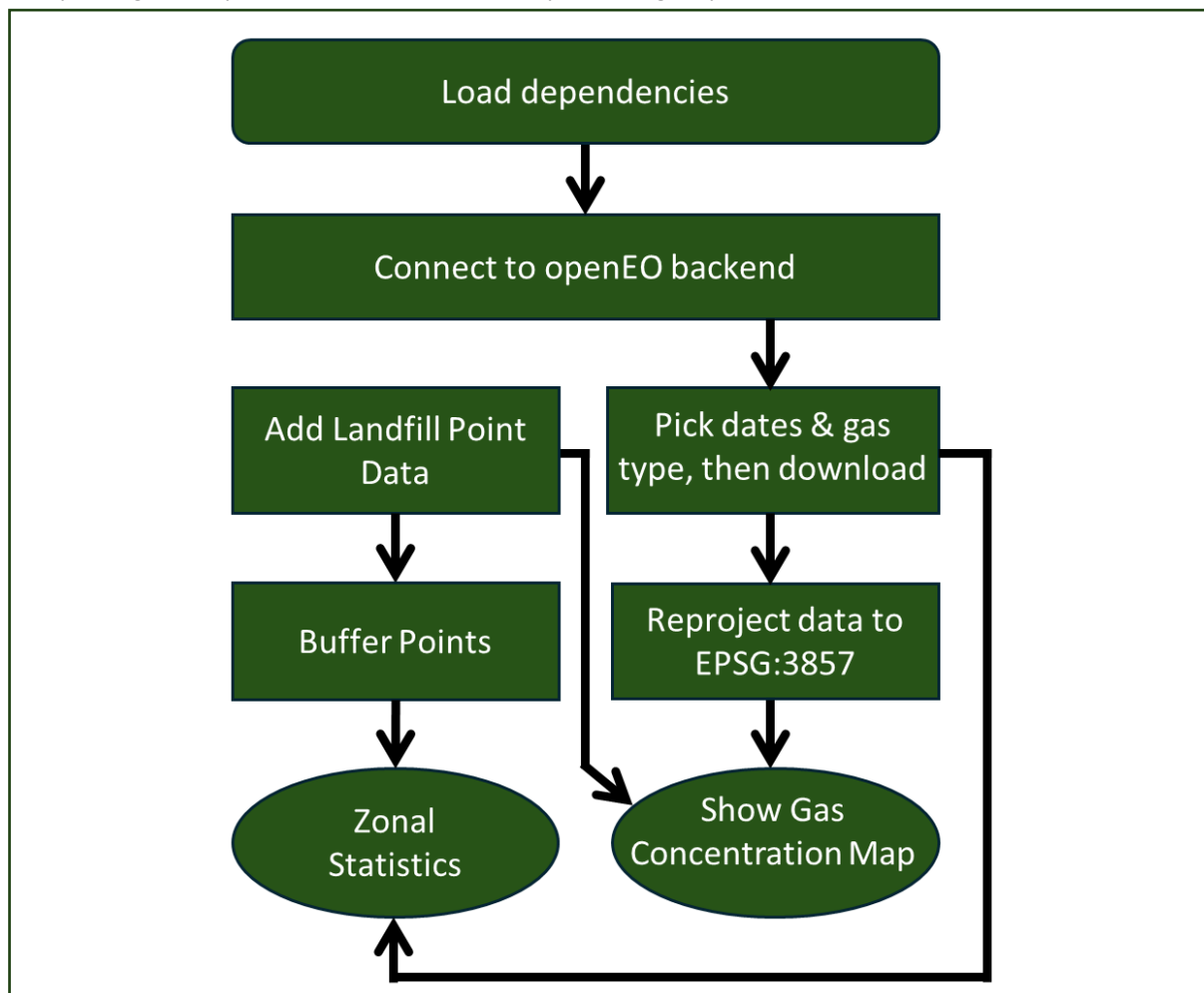


Figure 19: Flowchart of processes used for Gas Concentration Map

Code 9 shows the loaded dependencies.

```

import folium
import pandas
import matplotlib.pyplot as plt
import geopandas
import openeo
from shapely.geometry import shape, Point
from rasterio import warp
from matplotlib import cm
from matplotlib.colors import Normalize
import folium.raster_layers
import rasterio
from rasterio import warp
from rasterstats import zonal_stats
import numpy
  
```

Code 9: Loading of dependencies for the code to run.

Code 10 connects to the OpenEO backend.

```
connection = openeo.connect(url="openeo.dataspace.copernicus.eu")
connection.authenticate_oidc()
```

Code 10: Connecting to OpenEO

Code 11 shows the list of landfills as in S5-AGT

```
landfills =
geopandas.read_file(r"C:\GIS_Course\Landfill_Atmospheric_Gas_Monitor_Tools\D
ata\PZ_landfill_point4326.geojson")
```

Code 11: Displaying the contents of the landfill file for easy reference.

Code 12 selects the date to be viewed and what gas is to be monitored. It then downloads it as a .GTiff file.

```
cube = connection.load_collection(
    collection_id="SENTINEL_5P_L2",
    temporal_extent=["2023-06-01", "2023-06-01"], # format YYYY-MM-DD. Only
one date should be selected so the to and from fields should be identical.
    spatial_extent={"west": -19.5, "south": 27.0, "east": 5.0, "north":
44.5},
    bands=["CH4"], Gas monitoring options: 'CO', 'HCHO', 'NO2', 'O3', 'SO2',
'CH4'
)
cube.download("Sentinel-5P_Spain.GTiff")
```

Code 12: Code for selecting dates and gas, then downloading the data.

Code 13 takes the Sentinel-5P_Spain.GTiff file and reprojects its CRS from EPSG:4326 to EPSG:3857, the projection of the folium map. This code was provided in the course material (McNabb, 2024).

```
# The target coordinate reference system (CRS)
dst_crs = 'EPSG:3857'

# Open the gas data file that is in EPSG:4326 and calculate its bounds
with rasterio.open('Sentinel-5P_Spain.GTiff') as src:
    transform, width, height = rasterio.warp.calculate_default_transform(
        src.crs, dst_crs, src.width, src.height, *src.bounds)

    # Copy and update the metadata from the source dataset
    kwargs = src.meta.copy()
    kwargs.update({
        'crs': dst_crs,
        'transform': transform,
        'width': width,
        'height': height
    })

    # Create a new gas data file in EPSG:3857
    # Loop through each band in the source dataset
    for ind in range(1, src.count + 1):
        # Reproject each band and write it to the destination dataset
```

```

rasterio.warp.reproject(
    source=rasterio.band(src, ind),
    destination=rasterio.band(dst, ind),
    src_transform=src.transform,
    src_crs=src.crs,
    dst_transform=transform,
    dst_crs=dst_crs,
    resampling=rasterio.warp.Resampling.nearest)

```

Code 13: Code for reprojecting raster

Code 14 loads the folium map. The pins are clickable for the landfill information and the map is centred on those features. Because the CH₄ dataset is not a continuous raster, a piece of code ignoring those values has been included to aid basemap visibility. The colourmap is set using matplotlib. The raster bounds are set using a west, south, east, north longitude and latitude. Finally the legend bar scales to whatever atmospheric gas dataset is shown on the map.

```

# This section loads the map

#This is the reprojected gas data
gas_data = r"C:\GIS_Course\Landfill_Atmospheric_Gas_Monitor_Tools\Sentinel-
5P_Spain3857.GTiff"

# Open raster file, load values and prepare them to be displayed.
dataset = rasterio.open(gas_data, 'r')
rasdata = dataset.read()[0]
rasdata_normed = rasdata / rasdata.max() * 10

# When the data is displayed, this says to ignore values of zero.
non_zero_values = rasdata[rasdata != 0]
min_value = non_zero_values.min()
max_value = non_zero_values.max()
normalized_data = (non_zero_values - min_value) / (max_value - min_value)

# Create a colourmap for the non-zero values
colourmap = cm.turbo
colourmap_index = numpy.zeros_like(rasdata, dtype=numpy.float64)
colourmap_index[rasdata != 0] = normalized_data

#Loading in the landfill locations
PZ_landfill_Locations =
geopandas.read_file(r"C:\GIS_Course\Landfill_Atmospheric_Gas_Monitor_Tools\D
ata\PZ_landfill_point4326.geojson")

# This creates the map and centres it on the geometries.
centroids = PZ_landfill_Locations.geometry.centroid
center = [centroids.y.mean(), centroids.x.mean()]
gas_concentration_map = folium.Map(location=center, zoom_start=5,
tiles='CartoDB Positron', control_scale=True)

```



```

# Adding the PreZero landfill locations to the map and making them clickable
for info
for feature in PZ_landfill_Locations.iterfeatures():
    # Extract feature number from properties
    feature_number = feature['properties']['Landfill']
    # Extract coordinates of the feature
    coordinates = feature['geometry']['coordinates']
    # Create a marker with label for each feature
    folium.Marker(location=[coordinates[1], coordinates[0]],
                  popup=f"Feature
{feature_number}").add_to(gas_concentration_map)

# Adding the gas concentration dataset to the map
folium.raster_layers.ImageOverlay(
    image=colourmap(colourmap_index),
    name='gas concentration in atmosphere',
    opacity=0.6,
    bounds=[[20.0, -19.5], [44.4, 9.0]], # this should be the same as the
spatial extent of cube
    interactive=False,
    cross_origin=False,
    zindex=1
).add_to(gas_concentration_map)

# Creating the legend for gas concentration
fig, ax = plt.subplots(figsize=(8, 0.2))
cbar = plt.colorbar(cm.ScalarMappable(norm=Normalize(vmin=min_value,
vmax=max_value),
                        cmap=colourmap),
                    cax=ax, orientation='horizontal')
cbar.set_label('Parts per billion for CH4 or mol/m2 for all other gasses')

# Display the map
gas_concentration_map

```

Code 14: Code for configuring and loading Gas Concentration Map

Code 15 adds buffers to the point data and generates zonal statistics for each of the landfills.

```

# Same buffering code as S5-AGT without the geojson conversion
landfill_10000m = landfills
landfill_10000m = landfill_10000m.to_crs(epsg=2062)
landfill_10000m['geometry'] = landfill_10000m.buffer(10000)
landfill_10000m = landfill_10000m.to_crs(epsg=4326)

# Generate the zonal statistics
stats = zonal_stats(landfill_10000m, "Sentinel-5P_Spain.GTiff",
                    stats="sum min mean max median")

# Load zonal stats into geodataframe
stats_df = pandas.DataFrame(stats)

```

```
# Sort the data into columns
stats_df = stats_df[['sum', 'min', 'mean', 'max', 'median']]

# Display the statistics
print(stats_df)
```

Code 15: Code for configuring and loading Gas Concentration Map

4.3 Sentinel-2 Multi-Band-Single-Pass CH₄ Mapper

Sentinel 5's spatial resolution is very coarse, so to help users locate emission sources on a landfill, Sentinel 2's MSI instrument with a 20m² spatial resolution and return frequency of 3-5 days was employed.

Varon et al. (2021) demonstrated that methane columns from point sources can be measured by exploiting the SWIR-1 and SWIR-2 bands of Sentinel-2. Their recommended approach utilizes a multi-band-multi-pass (MBMP) retrieval method. However a multi band single pass (MBSP) method was employed to simplify the tool for users. A Jupyter workbook containing the MBMP method has been included in this project's repository for those interested.

The multi-band-single-pass equation is as follows:

$$MBSP = B11 - cB12$$

Where:

- B11 is the Sentinel-2 SWIR-1 band.
- B12 is the Sentinel-2 SWIR-2 band.
- c is calculated using the difference in the median of B12 to B11 and then adding the difference to B12.

Figure 20 provides an overview of the main steps in processing.

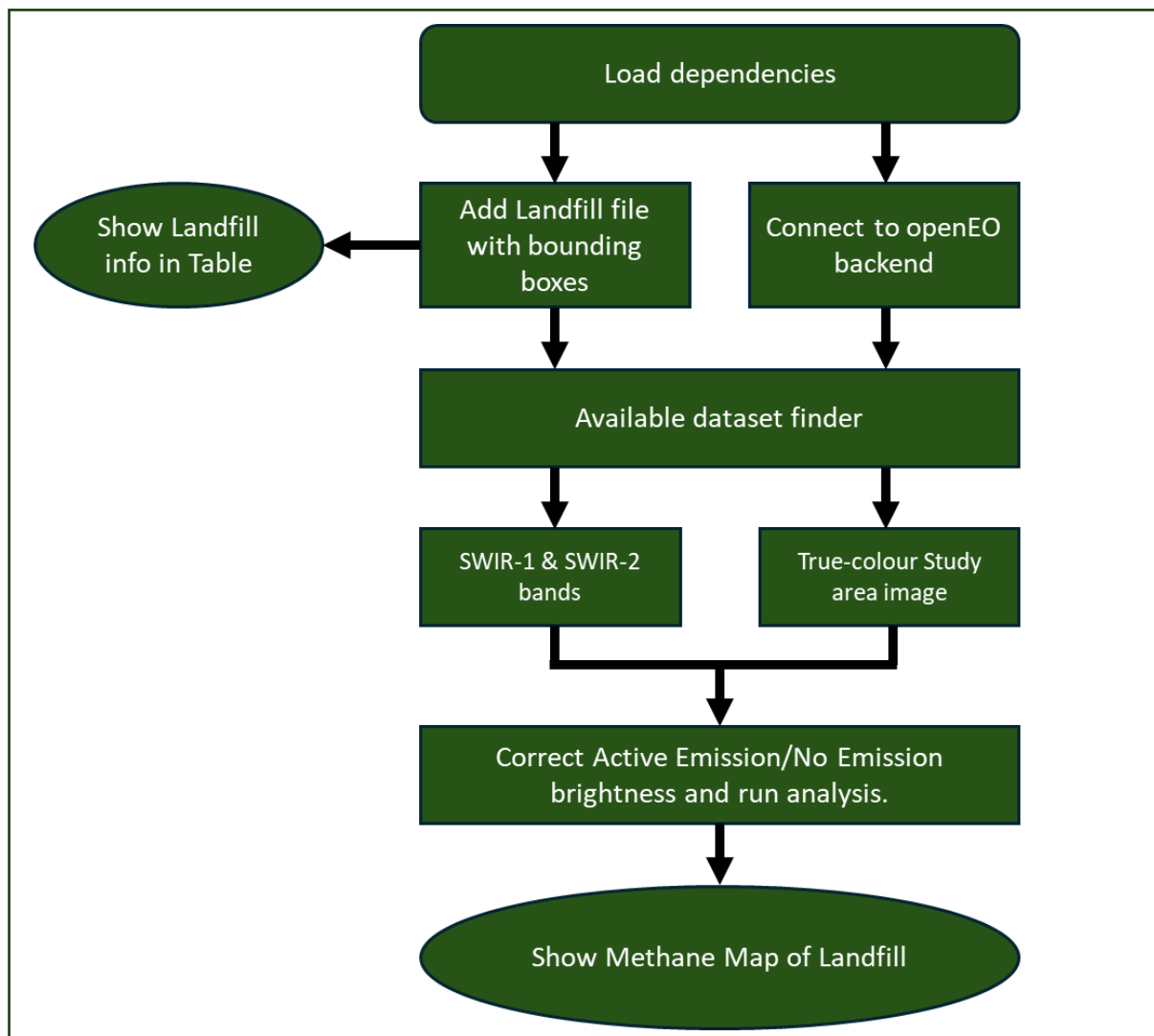


Figure 20: Flowchart of processes used for Sentinel 2 Landfill CH₄ Map

Code 16 shows the loaded dependencies.

```
import folium
import pandas
import matplotlib.pyplot as plt
import openeo
import rasterio
import numpy
import requests
```

Code 16: Loading of dependencies for the code to run.

Code 17 connects to the OpenEO backend.

```
connection = openeo.connect(url="openeo.dataspace.copernicus.eu")
connection.authenticate_oidc()
```

Code 17: Connecting to OpenEO

Code 18 reads a pre-processed file that contains the landfill location bounding boxes, and then displays it for the end user to see.

```
landfill_csv =
pandas.read_csv(r'C:\GIS_Course\Landfill_Atmospheric_Gas_Monitor_Tools\Data\
PreZero_Landfill_Bounding.csv')
landfill_csv
```

Code 18: Displaying the contents of the landfill file for easy reference.

Code 19 provides the available datasets for a date range and landfill selected by the user. It also has a cloud filter set at 15% as any cloud in the scene interferes with the tool. This section of code was provided by Sonneveld, E. (2024)

```
def get_spatial_extent(landfill_id):
    landfill = landfill_csv[landfill_csv['id'] == landfill_id].iloc[0]
    return {
        "west": landfill['west'],
        "south": landfill['south'],
        "east": landfill['east'],
        "north": landfill['north']
    }

def fetch_available_dates(landfill_id, temporal_extent):
    spatial_extent = get_spatial_extent(landfill_id)
    catalog_url =
f"https://catalogue.dataspace.copernicus.eu/resto/api/collections/Sentinel2/
search.json?box={spatial_extent['west']}%2C{spatial_extent['south']}%2C{spat
ial_extent['east']}%2C{spatial_extent['north']}&sortParam=startDate&sortOrde
r=ascending&page=1&maxRecords=1000&status=ONLINE&dataset=ESA-
DATASET&productType=L2A&startDate={temporal_extent[0]}T00%3A00%3A00Z&complet
ionDate={temporal_extent[1]}T00%3A00%3A00Z&cloudCover=%5B0%2C{cloud_cover}%5
D"

    response = requests.get(catalog_url)
    response.raise_for_status()
    catalog = response.json()
```



```

    dates = [date.split('T')[0] for date in map(lambda x:
x['properties']['startDate'], catalog['features'])]
    return dates

# Please enter your parameters here.
landfill_id = 23 # Specify the landfill ID.
temporal_extent = ["2023-01-01", "2023-12-30"] # Specify the the date range
you want to check for available data.
cloud_cover = 15

available_dates = fetch_available_dates(landfill_id, temporal_extent)
print("Available dates:", available_dates)

```

Code 19: Code for determining available dates for analysis.

Next the user will select the day that the emission was seen with the S5-AGT tool, along with the appropriate landfill id (code 20).

```

def active_emission(landfill_id, temporal_extent):
    landfill = landfill_csv[landfill_csv['id'] == landfill_id].iloc[0]

    active_emission = connection.load_collection(
        "SENTINEL2_L2A",
        temporal_extent=temporal_extent,
        spatial_extent={
            "west": landfill['west'],
            "south": landfill['south'],
            "east": landfill['east'],
            "north": landfill['north']
        },
        bands=["B11", "B12"],
    )
    active_emission.download("Sentinel-2_active_emissionMBSP.GTiff")

# Enter parameters for the active emission day
landfill_id = 23 # Specify the landfill ID
temporal_extent = ["2023-02-25", "2023-02-25"]
active_emission(landfill_id, temporal_extent)

```

Code 20: Code for downloading active emission dataset.

Code 21 downloads a true colour satellite image to aid in the visualisation of the data.

```

def truecolour_image(landfill_id, temporal_extent):
    landfill = landfill_csv[landfill_csv['id'] == landfill_id].iloc[0]

    truecolour_image_collection = connection.load_collection(
        "SENTINEL2_L2A",
        temporal_extent=temporal_extent,
        spatial_extent={
            "west": landfill['west'],
            "south": landfill['south'],
            "east": landfill['east'],

```

```

        "north": landfill['north']
    },
    bands=["B02", "B03", "B04"],
)
truecolour_image_collection.download("Sentinel-2_truecolourMBSP.GTiff")

# Enter parameters for the no emission day
landfill_id = 23 # Specify the landfill ID
temporal_extent = ["2023-02-25", "2023-02-25"]
truecolour_image(landfill_id, temporal_extent)

```

Code 21: Code for downloading true colour satellite image for data visualisation

Code 22 runs the analysis as outlined at the beginning of this section (4.3).

```

# Define file path
Active_Multiband = "Sentinel-2_active_emissionMBSP.GTiff"

# Define each band from the active dataset
with rasterio.open(Active_Multiband) as Active_img:
    Active_B11 = Active_img.read(1)
    Active_B12 = Active_img.read(2)

# Calculate the median difference for Active_B11 and Active_B12
median_diff_active = np.median(Active_B11) - np.median(Active_B12)

# Adjust Active_B12
Corrected_Active_B12 = Active_B12 + median_diff_active

# Calculate the fractional change
SWIR_diff = Active_B11 - Corrected_Active_B12

```

Code 22: Code running analysis

Code 23 displays the map. It firstly takes the true colour satellite image bands, stacks them and applies a brightness factor to make the image clearer. Then it adds the SWIR_diff data calculated in the previous code and sets it to only display data between 1.5 and 3 standard deviations above the mean. It adds scaled grid lines to help understand distances on the image. Finally it downloads the map as a .jpg file.

```

# This section adds the true colour satellite image.
# Open the and load the true colour image file and define which band is
which.
truecolour_sat = 'Sentinel-2_truecolourMBSP.GTiff'
img = rasterio.open(truecolour_sat)
blue = img.read(1)
green = img.read(2)
red = img.read(3)

# Change this number up or down if the satellite background image is too
dark or bright.
brightness_factor = 0.05
blue = numpy.clip(blue * brightness_factor, 0, 255)
green = numpy.clip(green * brightness_factor, 0, 255)

```

```
red = numpy.clip(red * brightness_factor, 0, 255)

# Stack the blue, green and red bands to make a colour image.
rgb = numpy.dstack((red, green, blue))
rgb = rgb / rgb.max()

# Create a new map
plt.figure(figsize=(10, 8))

# Display the RGB image on that map.
plt.imshow(rgb)

# This section adds the methane column data.
# Calculate mean and standard deviation of SWIR_diff.
mean = numpy.nanmean(SWIR_diff)
std = numpy.nanstd(SWIR_diff)

# Create a mask to only show values 1.5 SD above the mean.
mask = SWIR_diff < (mean + 1.5 * std)

# Apply the mask
masked_SWIR_diff = numpy.ma.masked_array(SWIR_diff, mask=mask)

# Set the minimum and maximum displayed data values to be between 1.5 and 3
standard deviations above the mean.
mean = numpy.nanmean(SWIR_diff)
std = numpy.nanstd(SWIR_diff)
vmin = mean + 1.5 * std
vmax = mean + 3 * std

# Display the masked_SWIR_diff data on top of the RGB image.
plt.imshow(masked_SWIR_diff, cmap='plasma', alpha=1, vmin=vmin, vmax=vmax)

# This next section adds features to the map to aid in interpretation.
# Add a colorbar and labels
cbar = plt.colorbar(label='CH4 Standard Deviation Above Background Levels',
shrink=0.7)
cbar.set_ticks([vmin, vmax])
cbar.set_ticklabels(['1.5 SD', '3 SD'])
plt.title('CH4 Concentration')

# Dataset resolution in metres.
resolution = 20

# Get the dimensions of the image.
height, width, _ = rgb.shape

# Create axis showing scale in metres with grid lines every 500m.
```

```
x = numpy.arange(0, width * resolution, 500)
y = numpy.arange(0, height * resolution, 500)
plt.xticks(x / resolution, x)
plt.yticks(y / resolution, y)
plt.xlabel('X (meters)')
plt.ylabel('Y (meters)')
plt.grid(color='gray', linestyle='--', linewidth=0.5)

# Downloading image for use in report
# Please enter the emission date where it says '[emission date here]' if you
# want to download more than one map.
plt.savefig(f'S2MBSP[emission_date_here].jpg', format='jpg',
bbox_inches='tight')

# Show the map
plt.show()
```

Code 23: Code displaying the map

5. Expected Results and Demo

This section will illustrate the functionality of the tools by showcasing documented methane plumes and their corresponding depiction in the results.

5.1 Sentinel-5 Atmospheric Gas Timeseries:

The S5-AGT tool generates line graphs similar to those depicted in figures 21 and 22. In these graphs, dots connected by lines represent consecutive days of methane emissions.

On April 23, 2023, a methane plume was observed near the PreZero-managed Cañada Hermosa (Jet Propulsion Laboratory, 2024). The data suggests that the CH₄ emission may have occurred over a span of 9 days, from February 15 to February 24, 2023 (see fig. 21). However, it's possible that this was a short-duration event within a broader pattern of methane presence in the atmosphere.

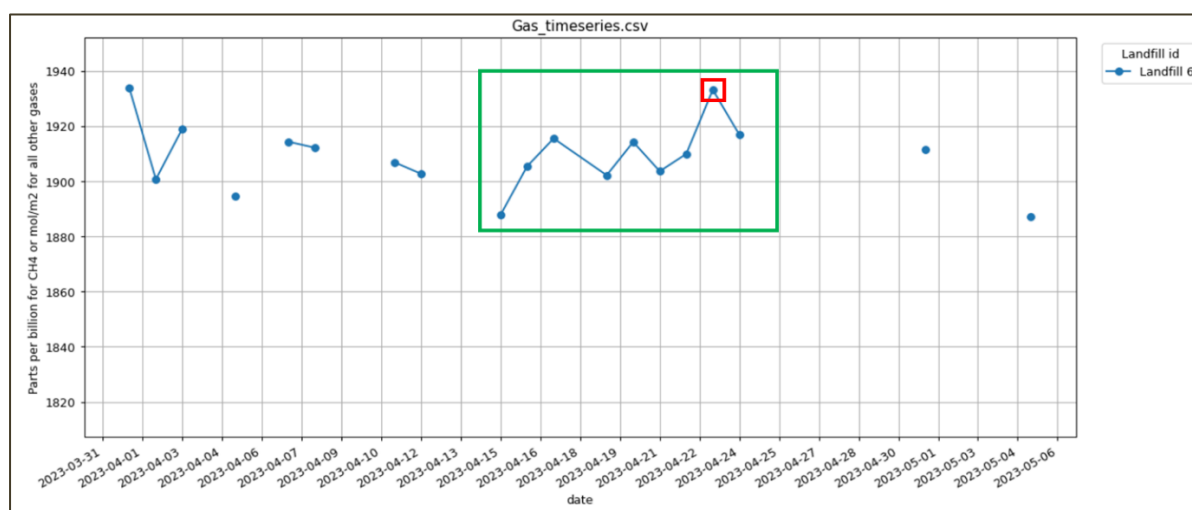


Figure 21: timeseries of landfill 6 'Cañada Hermosa', with the signal relating to the plume on the 23rd of April 2023 highlighted in red and sequential days of emissions highlighted in green.

Less ambiguous candidates for plumes are singular high emissions. Figure 22 shows a known emission from the Pinto biomethanization landfill on the 24th of February 2023 near Madrid (The Guardian, 2024) and 2 other isolated high values, which may warrant further investigation.

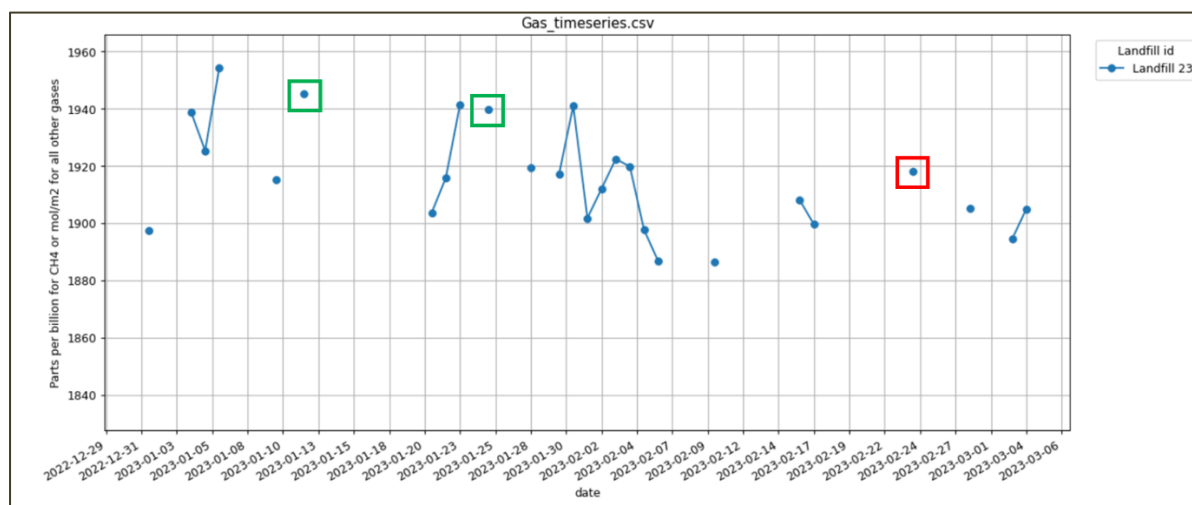


Figure 22: timeseries of landfill 23 'Pinto biomethanization plant', with the signal relating to the 24th of February 2023 plume highlighted in red and two isolated high values highlighted in green.

5.2 Sentinel-5 Atmospheric Gas Concentration Mapper:

Following on from the S5-AGT tool, the S5-AGM can show a specific Sentinel 5 dataset for further investigation. Figure 23 and 24 below show the data for the 24th of February 2023.

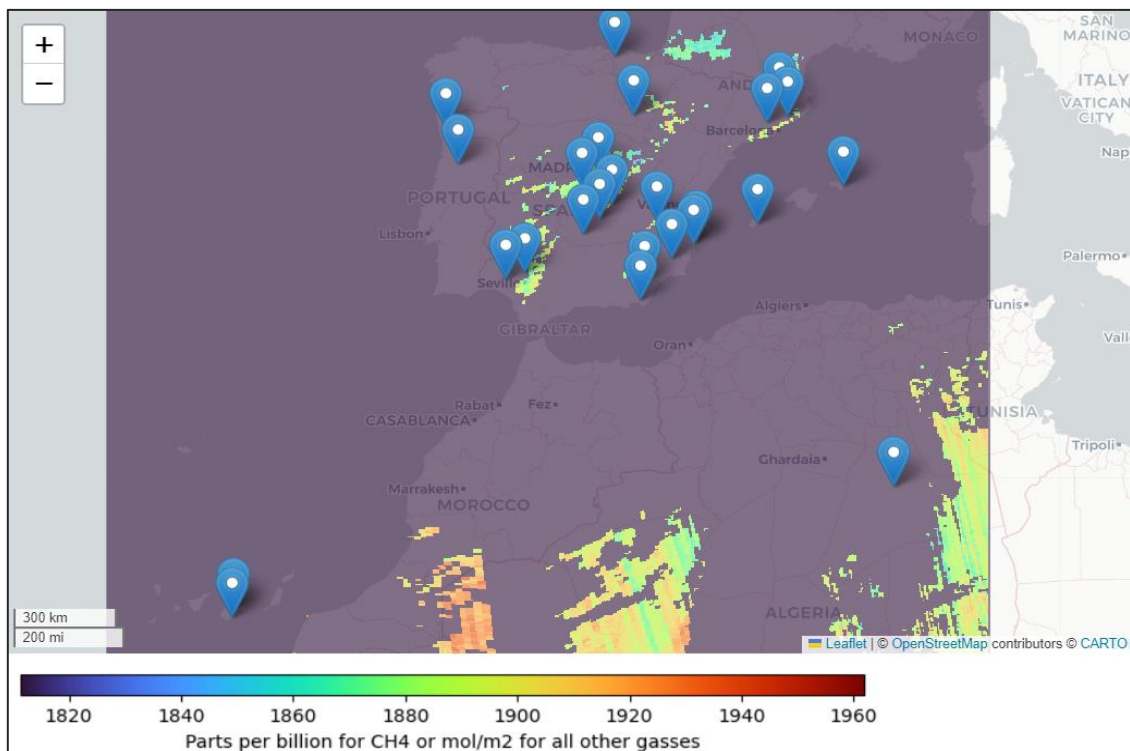


Figure 23: Screenshot of CH₄ map, over Spain with landfill locations shown as blue pins.

The user can then zoom in to an area of interest. In figure 24 the Pinto Waste Site is shown along with a distinct hotspot of dark orange, characteristic of a CH₄ plume (European Space Agency, 2021).

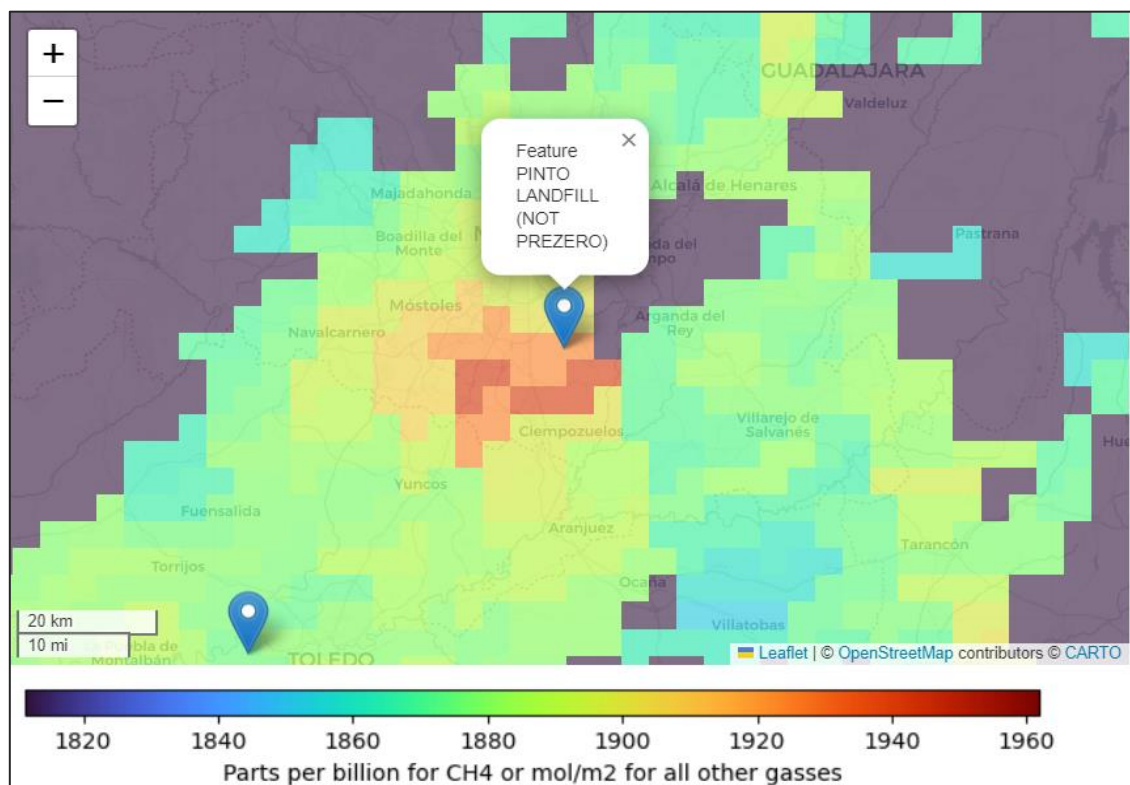


Figure 24: Sentinel 5P data showing methane plume from the Pinto Biomethanization Landfill

Examples of gas datasets for CO, HCHO, NO₂, O₃ and SO₂ are shown in figure 25.

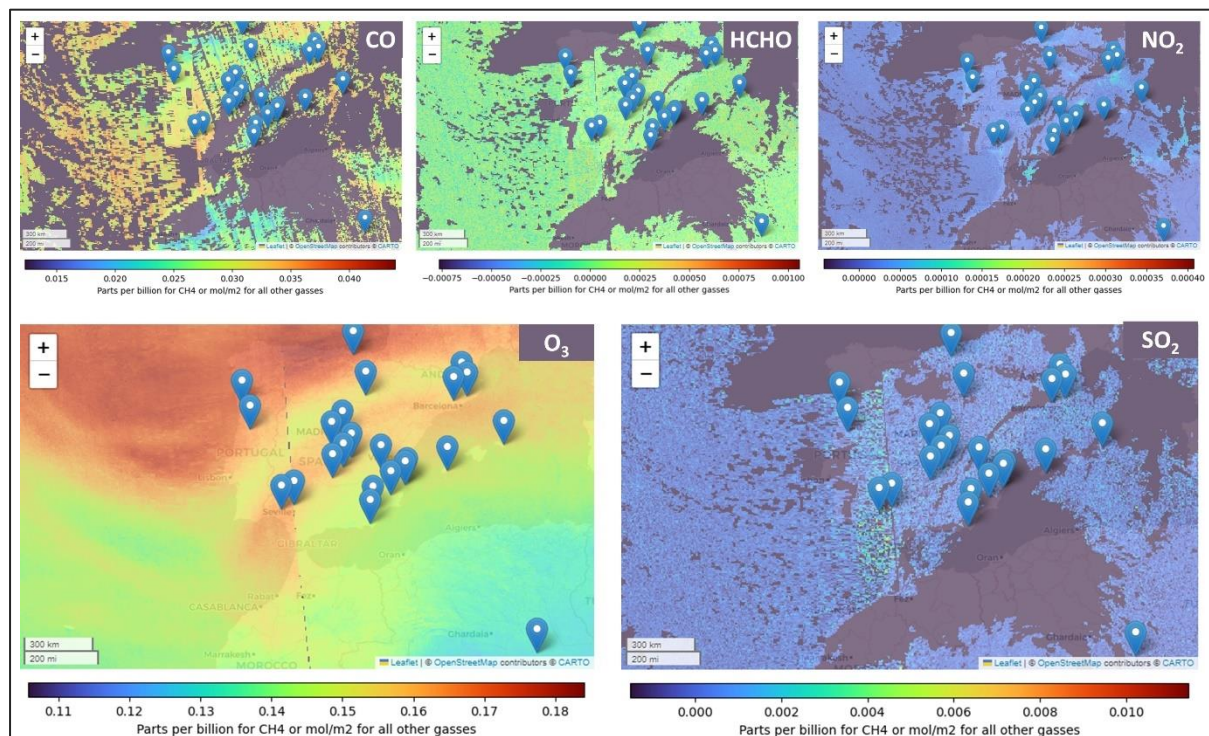


Figure 25: Gas datasets for CO, HCHO, NO₂, O₃ and SO₂ for the 24th of February 2023

5.3 Sentinel-2 Multi-Band-Multi-Pass CH₄ Mapper:

To ensure that this tool works as intended, a known emission in Algeria used by Varon et al. (2021) was imaged (fig.26).

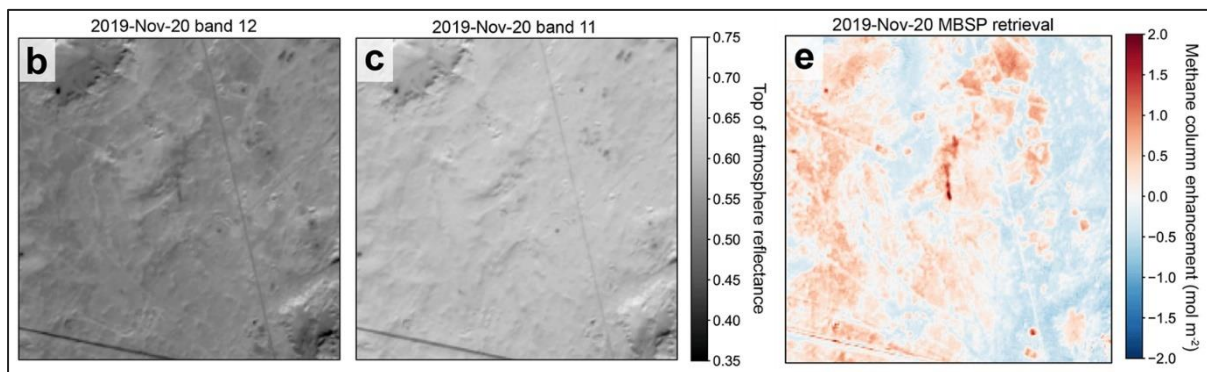


Figure 26: Sentinel 2 SWIR-2 (b), SWIR-1 (c) and Methane column retrieval for Multi Band Single Pass method (e) from Varon et al. (2021)

In figure 27, you can see that the methane column has been retrieved successfully. This plume can be found in the PreZero_Landfill_Bounding.csv file if you want to run this test.

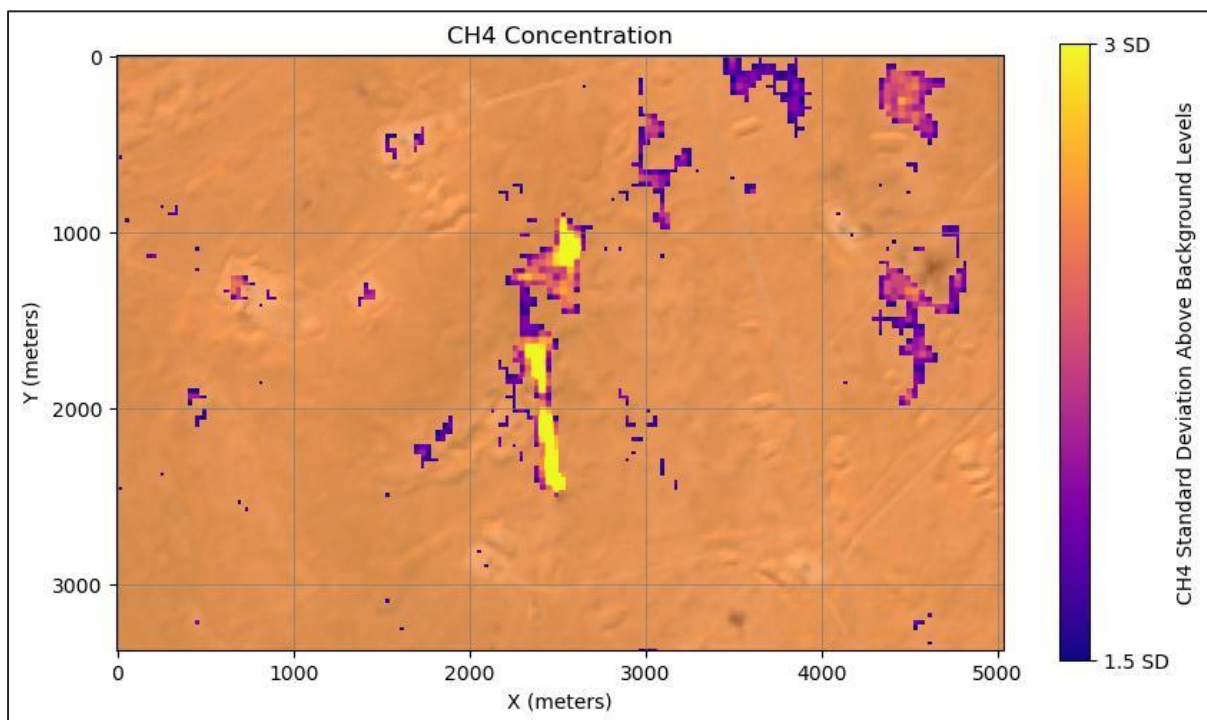


Figure 27: Methane column retrieval using S2-MBSP tool.

The emission date of 24th Feb 2023 for the Pinto Biomethanization Landfill, wasn't available due to Sentinel 2's return period (Sentinel Hub, 2024). However, the day after was available (fig.28)

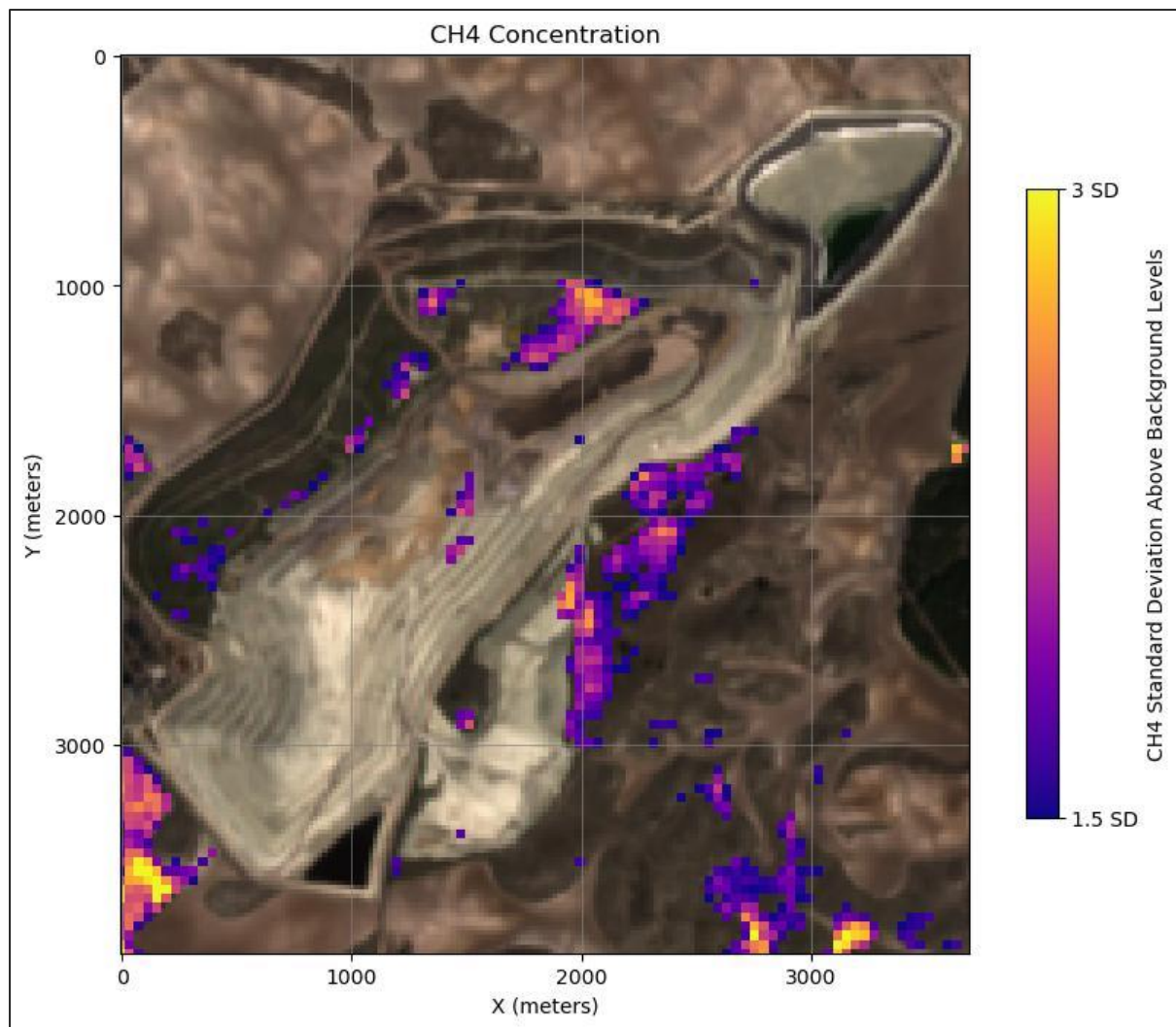


Figure 28: Pinto Madrid on the 25th of February 2023 with possible methane emission sources.

6. Troubleshooting

Input errors have been covered in the tool notebooks, however there are errors unrelated to a user input which can cause problems. These are detailed below.

6.1 Remote disconnected error (All tools)

Error: Remote disconnected

Or

OpenEoApiError: [500] Internal: Server error: KazooTimeoutError('Connection time-out') (ref: r-2405108830e742b59ef8ac2f28647fb5)

This can occur when there are issues with the Copernicus network. In the event that you see an error like this you can check page <https://dataspace.copernicus.eu/news> for any downtime messages and you can also contact the Copernicus dataspace team via the form at <https://helpcenter.dataspace.copernicus.eu/hc/en-gb/requests/new>

6.2 Concurrent job error (S5-AGM)

OpenEoApiError: [400] ConcurrentJobLimit: Job was not started because concurrent job limit (2) is reached. (ref: r-240413b5d1b240118da9f9ed90807c58)

This can happen when S5-AGT tool is run, cancelled and then run again. If this happens the process is still running in the background and needs to be cancelled.

To do this go to the following URL: <https://openeo.dataspace.copernicus.eu/>

You will be presented with the following screen (fig.29). Please click the highlighted login button.

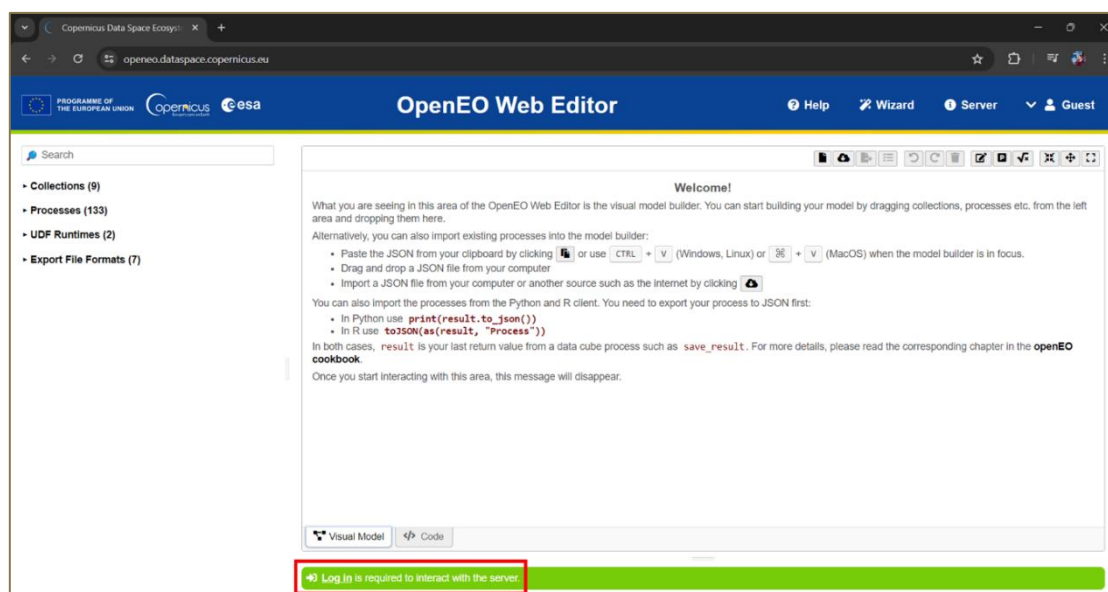


Figure 29: OpenEO Web Editor with login button highlighted in red

You will then see the following screen (fig.30).

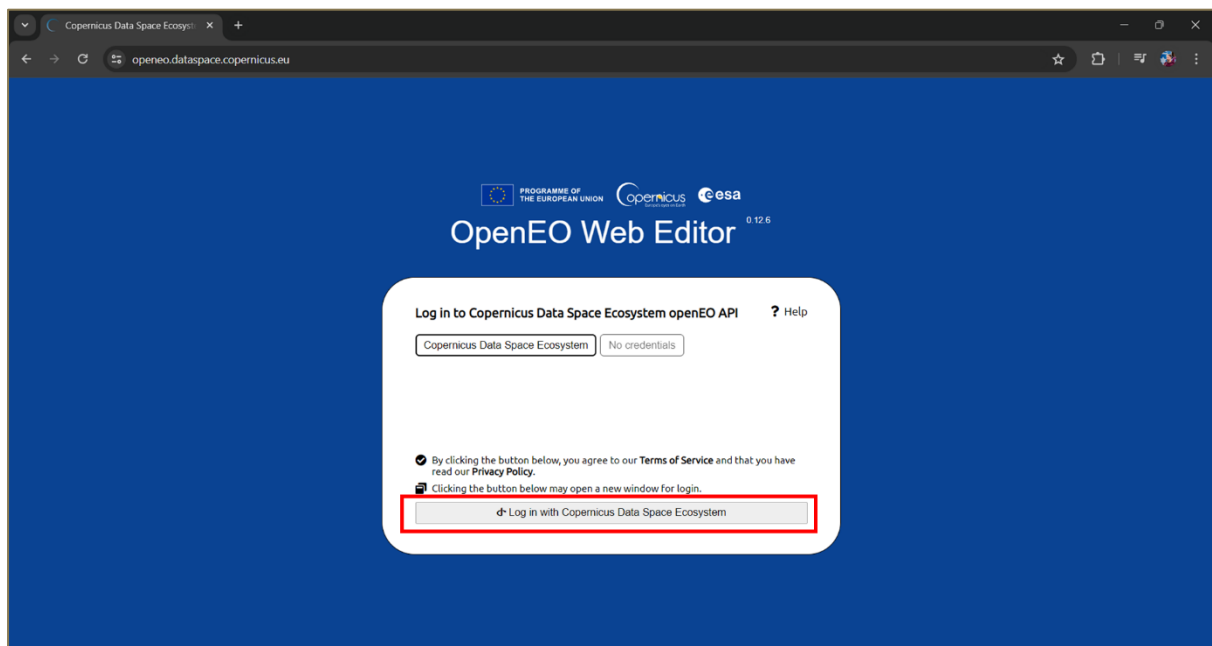


Figure 30: OpenEO Web editor login prompt with login button highlighted in red.

Simply click the highlighted button (fig.31) and follow the process. You should then be returned to the OpenEO web editor but now you will see a list of processes including the active ones.

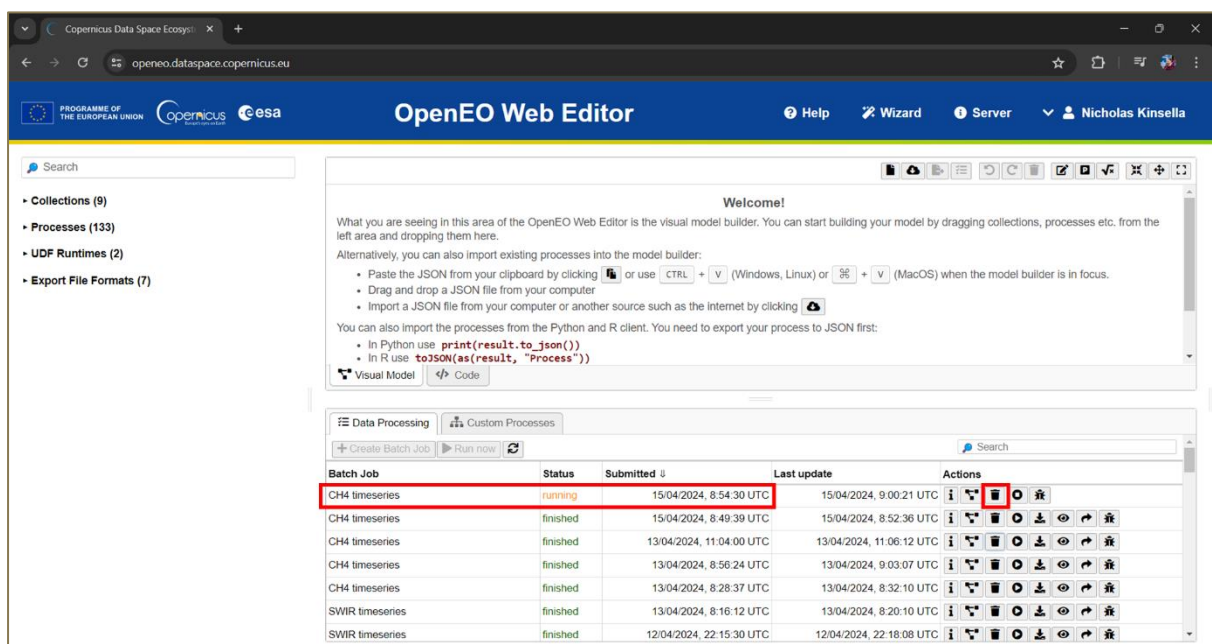


Figure 31: OpenEO Web Editor showing batch job screen, with running job and delete button highlighted.

To stop the process, simply click the highlighted bin button (delete). This should allow the tool to work normally again.

7. References

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