

# pyopmspe11: A Python framework using OPM Flow for the SPE11 benchmark project

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## Software

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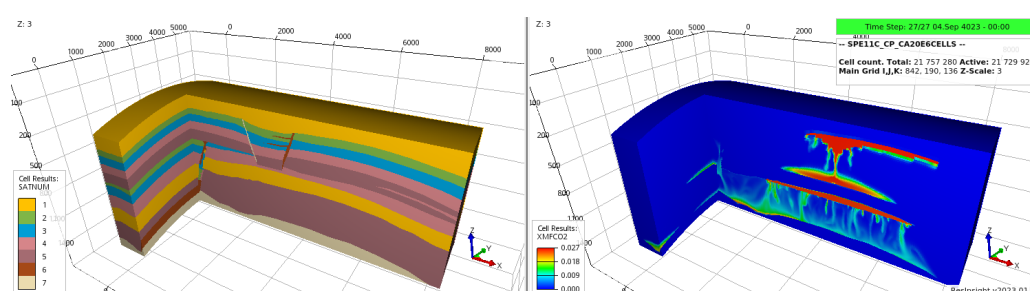
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## Summary

The imperative to achieve climate change goals and the increasing worldwide demand for energy have made geological carbon storage (GCS) technology more relevant today. Since utilizing computational models is essential for planning large-scale GCS projects, it is crucial to benchmark simulation tools to enhance confidence in their results. Inspired by a recent validation study for laboratory-scale CO<sub>2</sub> storage ([Flemisch et al., 2024](#)), a new comparative solution project (CSP) was launched to simulate both lab- and field-scale CO<sub>2</sub> storage ([Nordbotten et al., 2024](#)). This project is called the 11th Society of Petroleum Engineers CSP, and we refer to it as the SPE11 benchmark. The main objective for the SPE11 benchmark is to provide a common platform and reference case for numerical simulation of GCS. A community effort was run by the “Early Access Team” to create utility scripts and input files for popular simulators to make participation more accessible. As part of the “Early Access Team”, we have developed and made open the pyopmspe11 tool which facilitates reproducible solutions to the SPE11 benchmark. This tool serves as a common starting point for developing and testing new GCS simulation technology. Due to its user-friendly functionality (e.g., generation of different types of grids at different grid resolutions, flexibility to choose different rock and fluid properties, flexibility to define well/source locations and schedule for operations), it is expected that its impact will extend far beyond the initial benchmark study (e.g., studies focusing on grid refinement, upscaling/coarsening approaches, numerical solvers, optimization/history matching techniques).



**Figure 1:** Generated model by the configuration file `spe11c_cp_ca20e6cells.txt` in the [examples](#) folder (the model corresponds to the SPE11C Case using a corner-point grid with 21729920 active cells).

## Statement of need

Geological carbon storage (GCS) applications benefit from both commercial and open-source simulators. OPM Flow is an open-source simulator for subsurface applications such as hydrocarbon recovery, CO<sub>2</sub> storage, and H<sub>2</sub> storage. The typical workflow in GCS simulations starts with defining the simulation model (e.g., grid, heterogeneity, physics, fluid properties,

boundary conditions, wells), then setting the simulator parameters (e.g., tolerances, linear solvers, partition algorithms), then running the simulation, and finally visualization/analysis of the simulation results (e.g., CO<sub>2</sub> plume distance to the boundaries, caprock integrity). Here we refer to the first two steps as preprocessing and the final step as postprocessing. Notable works are available in JOSS for pre-/postprocessing of simulation data, e.g., Beucher et al. (2019), Sullivan & Kaszynski (2019), Fraters et al. (2024), Kaus et al. (2024). However, preprocessing and postprocessing can be challenging for everyone even if you know what you are doing. Additionally, setting up and running simulations requires computational expertise. To bridge this gap, developers can simplify the setup of numerical studies by using user-friendly approaches, such as configuration files. This not only ensures reproducibility of results but also facilitates flexible testing of different simulator parameters and allows for easy extension to further studies.

Based on the acquired knowledge by contributing to other open-source projects such as OPM Flow, then we have developed and made open the `pyopmspe11` tool which facilitates reproducible solutions to the SPE11 benchmark, which focus on GCS at different scales (Nordbotten et al., 2024). A previous benchmark study for GCS can be found in Class et al. (2009). One key difference of the SPE11 benchmark from the benchmark in Class et al. (2009) is that no specific size and type of grids were given in the description, i.e., one of the main task for the SPE11 benchmark participants was to create suitable grids (e.g., structured grids such as Cartesian or unstructured grids such as corner-point grids) and to run the cases, i.e., computational grids. To ease the comparison of results between participants, the SPE11 benchmark organizers requested the reporting of the spatial maps to follow a specific format using Cartesian grids with fixed cell sizes, i.e., reporting grids. The participants were encouraged to share data (e.g., input decks, code, submitted results), with the opportunity to store the data for open access. This is where developing tools that made all steps reproducible (i.e., preprocessing and postprocessing) became handy, and for this benchmark study, one available tool is `pyopmspe11`. Examples of pre-/postprocessing simulation tools which also have application in GCS and rely on the OPM Flow simulator include: `pyopmnearewell` (Landa-Marbán & pschultzendorff, 2023) and `expreccs` (Landa-Marbán, 2024). The former focuses on near well dynamics, while the latter on seamless, dynamic, and non-invasive exchange of pressure-related information between local and regional scales.

`pyopmspe11` is a simplified and flexible Python tool to execute the three cases in the SPE Comparative Solution Project using configuration files. `pyopmspe11` relies on the OPM Flow numerical simulator (Rasmussen et al., 2021), where the implementation of the CO<sub>2</sub> model can be found in Sandve et al. (2021). The primary contribution of `pyopmspe11` lies in its data preprocessing and postprocessing capabilities. It offers flexibility in generating various types of grids, including Cartesian, tensor, and corner-point grids. These grids adhere to the standard industry format (i.e., Eclipse grid format), making them compatible not only with OPM Flow but also with other simulators. Here, we mention two existing widely-used visualization/postprocessing software for OPM Flow: `ParaView` and `ResInsight`. While these tools are very useful, to the authors knowledge, there are no available built-in options via the GUIs of these tools to handle all necessary postprocessing to generate all data reporting as required in the SPE11 benchmark study. Instead, this could be achieved in these tools by code development using the Python interface. The benefit of having a specialized tool like `pyopmspe11`, which integrates all stages from preprocessing to simulation and postprocessing, is that it simplifies usage.

`pyopmspe11` supports varying resolutions, having been tested to generate up to 160 million cells. In the context of data postprocessing, `pyopmspe11` not only generates the necessary reporting data as specified by the benchmark, but it also produces .png figures for rapid inspection of individual simulations and for making comparisons between different runs (e.g., to assess sensitivities). The postprocessing methods efficiently interpolate quantities over time and map non-overlapping cell values (both intensive and extensive quantities) between the computational grid and the reporting grid. The Python package Scipy (Virtanen et al., 2020),

specifically the `interp1d` Class, is used for the time interpolation. The Python package `Shapely` (Gillies et al., 2024), specifically the `Polygon` Class, is the base for the developed methods in `pyopmspe11` to handle the mapping from the computational grid to the reporting grid.

## Outlook

`pyopmspe11` is designed for use by researchers, engineers, and students. During the preliminary intercomparison workshops for the benchmark study, the authors received positive feedback about the framework. Some participant groups have utilized `pyopmspe11`, particularly for grid generation. Additionally, the authors have been contacted to provide specific support on how to use the tool for setting up further studies, such as optimizing injection strategies. Looking ahead, the plan for `pyopmspe11`'s future development includes extending its functionality to support the generation of input decks to run simulations of physical models available in OPM Flow in addition to CO<sub>2</sub> storage (e.g., hydrogen storage, salt precipitation, biofilm effects).

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