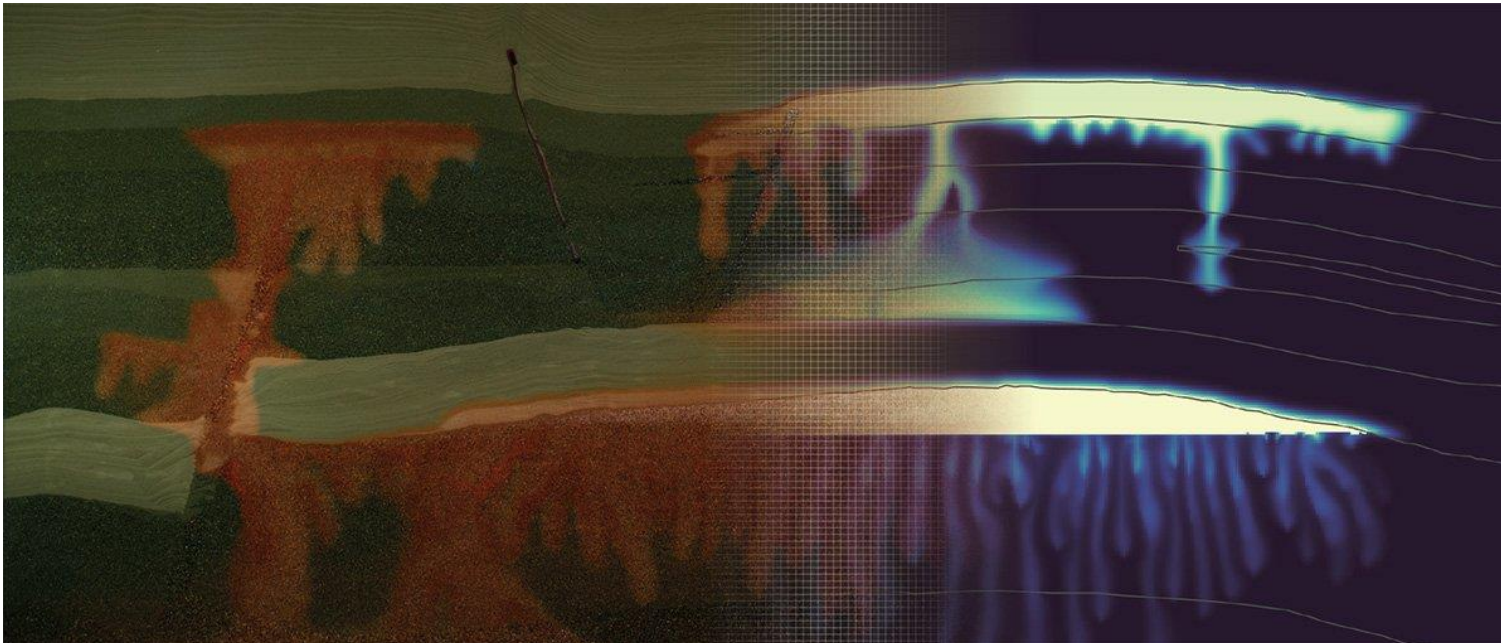


SPE11 CSP using OPM Flow

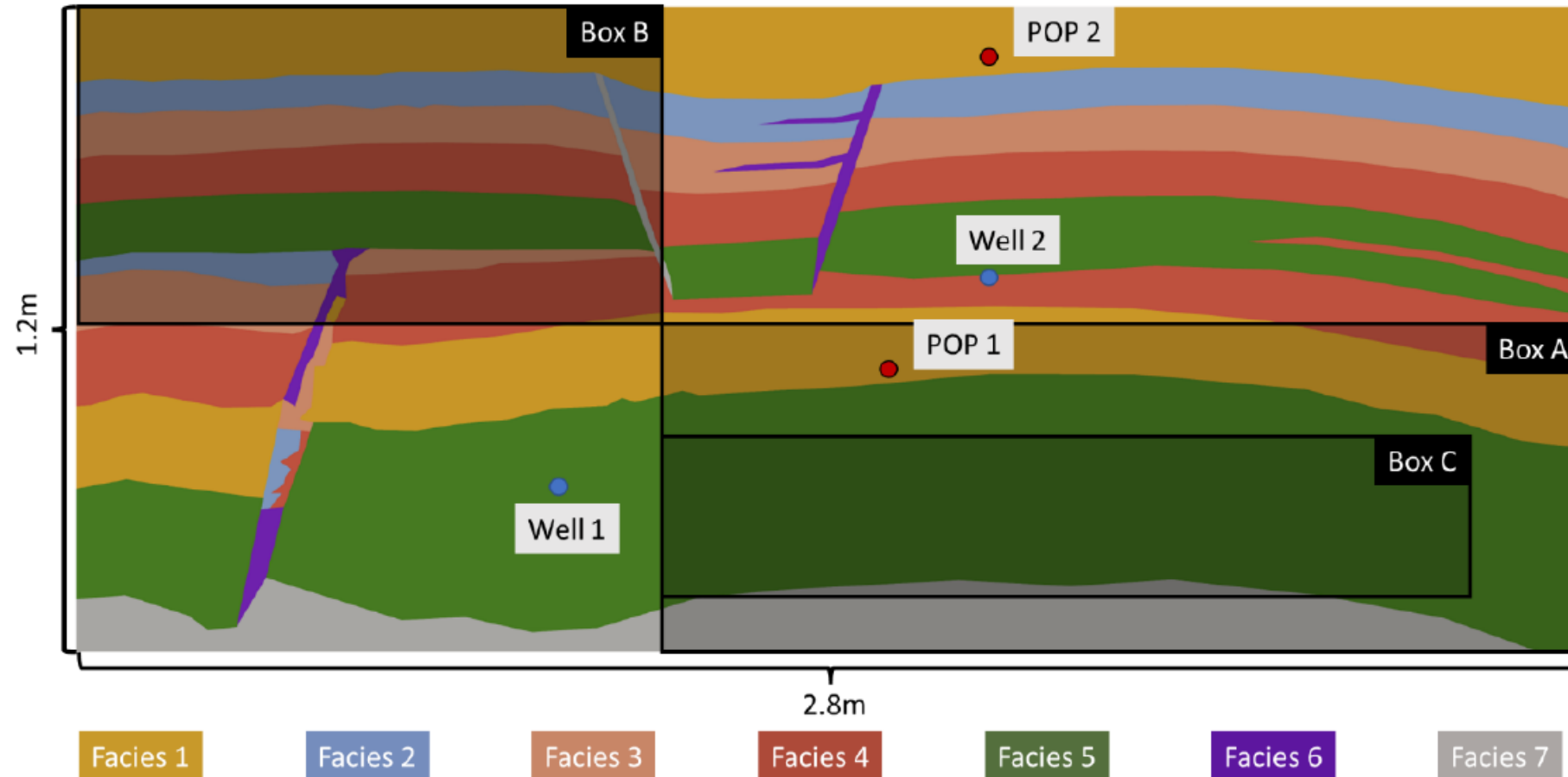
Tor Harald Sandve and David Landa-Marbán

SPE11 CSP using OPM Flow

- The 11th Society of Petroleum Engineers Comparative Solution Project
- 11 participants that submitted results
- <https://www.spe.org/en/csp/>

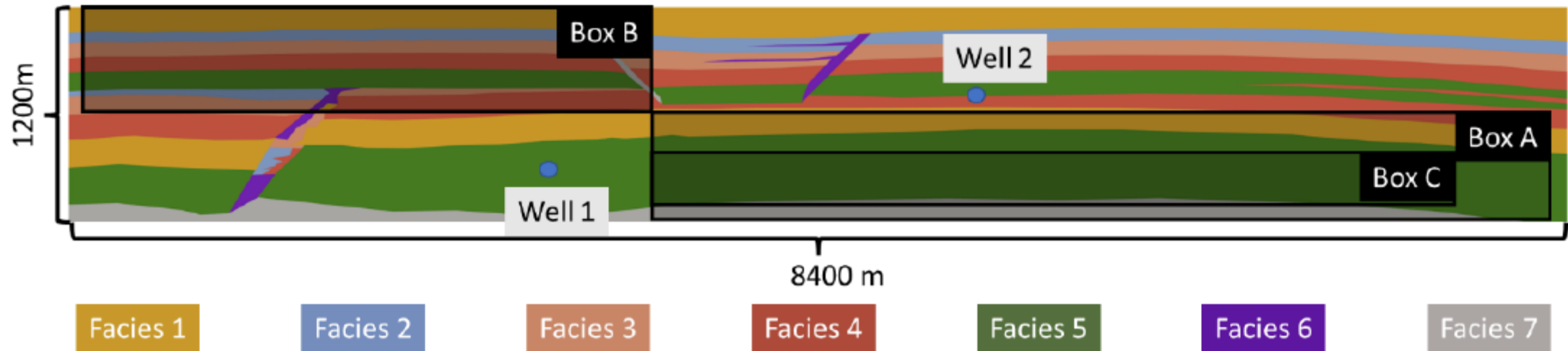


SPE 11 A



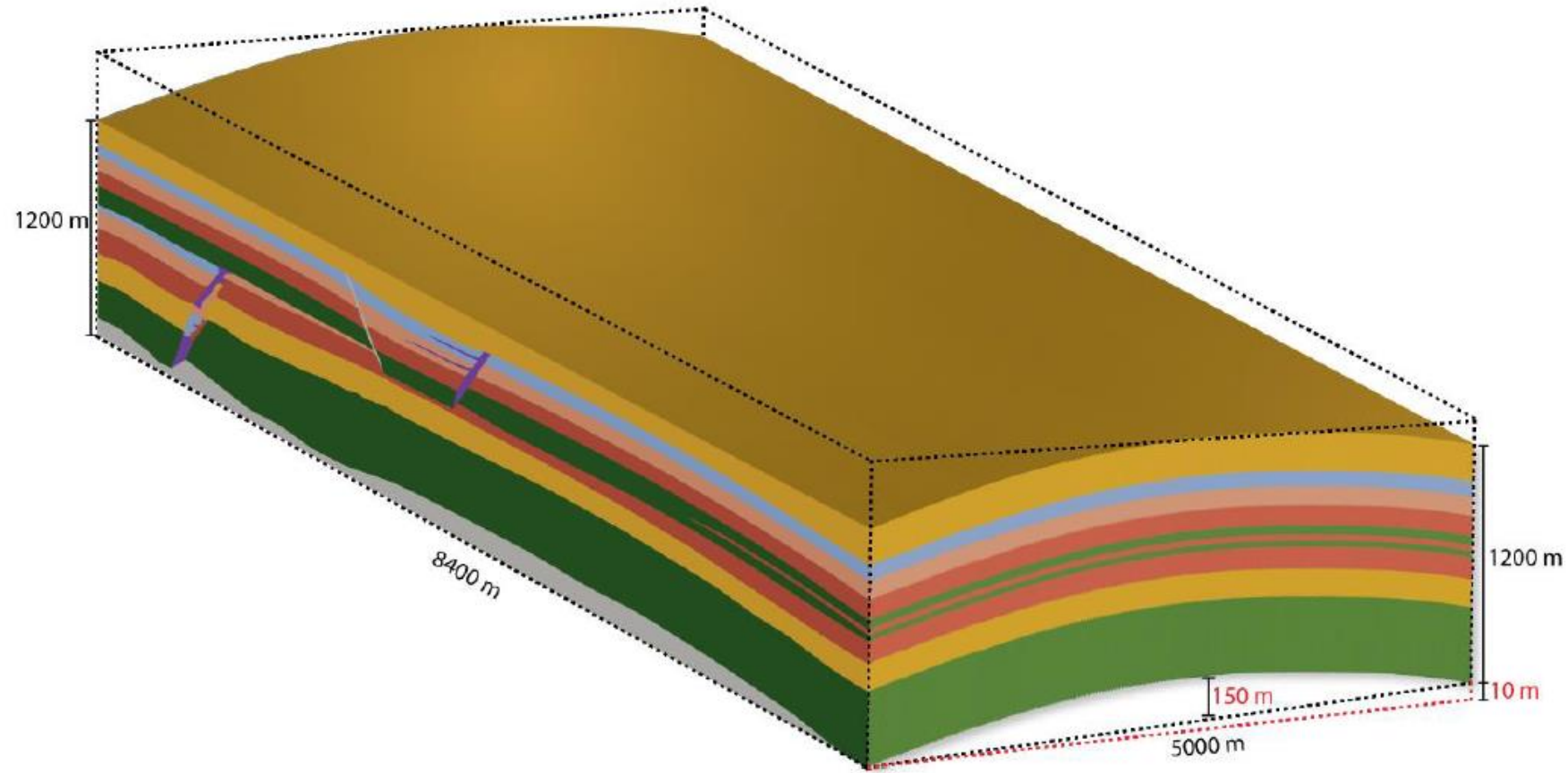
SPE 11A is a 2D geometry at the laboratory scale, inspired by a recent CO₂ storage forecasting and validation study (The FluidFlower).

SPE 11 B



For SPE 11B, the 2D geometry and operational conditions from 11A are rescaled to field conditions characteristic of the Norwegian Continental Shelf.

SPE 11 C



SPE 11C, the geometry of version 11B is extruded to a full 3D field model.

A Python framework using OPM Flow for the SPE11 benchmark project



- Source code: <https://github.com/OPM/pyopmspe11>
- Documentation: <https://opm.github.io/pyopmspe11/index.html>
- Landa-Marbán, D. and Sandve, T. H., pyopmspe11: A Python framework using OPM Flow for the SPE11 benchmark project. Under review JOSS.

```
pyopmspe11 -i configuration_file
```

Configuration file

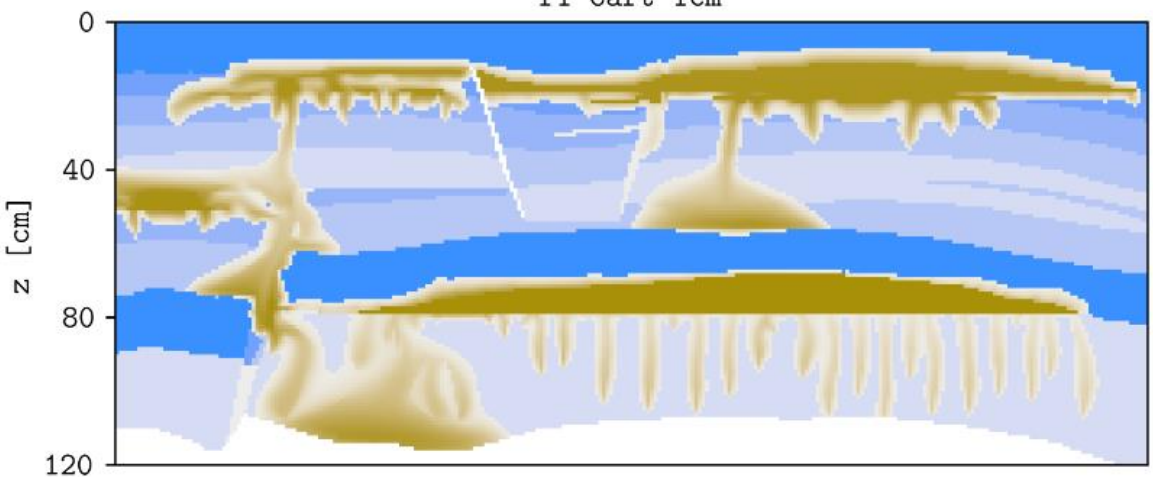


```
1 # Set the full path to the flow executable and flags
2 flow = "flow --enable-tuning=true --enable-opm-rst-file=true --output-extra=
3
4 # Set the model parameters
5 spe11 = "spe11c" # Name of the spe case (spe11a, spe11b, or spe11c)
6 version = "release" # OPM Flow version (release or master)
7 model = "complete" # Name of the co2 model (immiscible, convective, or con
8 co2store = "gaswater" # co2store implementation (gaswater or gasoil [oil p
9 grid = "corner-point" # Type of grid (cartesian, tensor, or corner-point)
10 dims = [8400.0, 5000.0, 1200.0] # Length, width, and depth [m]
11 x_n = [420] # If cartesian, number of x cells [-]; otherwise, variable arr
12 y_n = [30, 40, 50, 40, 30] # If cartesian, number of y cells [-]; otherwis
13 z_n = [5, 3, 1, 2, 3, 2, 4, 4, 10, 4, 6, 6, 4, 8, 4, 15, 30, 9] # If carte
14 temperature = [70.0, 36.12] # Temperature bottom and top rig [C]
15 datum = 300 # Datum [m]
16 pressure = 3e7 # Pressure at the datum [Pa]
17 kzMult = 0.1 # Multiplier for the permeability in the z direction [-]
18 diffusion = [1e-9, 2e-8] # Diffusion (in liquid and gas) [m^2/s]
19 rockExtra = [8.5e-1, 2500.0] # Rock specific heat capacity [kJ/(kg K)] and
20 pvAdded = 5e4 # Extra pore volume per area on lateral boundaries [m] (for
21 widthBuffer = 1 # Width of buffer cells [m] (for spe11b/c)
22 elevation = 150 # Maximum elevation difference (relative to the baseline g
23 backElevation = 10 # Back boundary elevation w.r.t the front boundary [m]
24 dispersion = [10, 10, 10, 10, 10, 10, 0] # Dispersion rock [m], facie 1 to
25 rockCond = [1.9, 1.25, 1.25, 1.25, 0.92, 0.26, 2.0] # Thermal conductivity
26 radius = [0.15, 0.15] # Wells radius [m] (0 to use the SOURCE keyword inst
27 wellCoord = [[2700.0, 1000.0, 300.0], [5100.0, 1000.0, 700.0]] # Well posi
28 wellCoordF = [[2700.0, 4000.0, 300.0], [5100.0, 4000.0, 700.0]] # Well fir
```

```
30 # Set the saturation functions
31 krw = "(max(0, (s_w - swi) / (1 - swi))) ** 1.5"
32 krn = "(max(0, (1 - s_w - sni) / (1 - sni))) ** 1.5"
33 pcap = "penmax * math.erf(pen * ((s_w-sw_i) / (1.-sw_i)) ** -(1.0 / 1.5)) *
34 s_w = "(np.exp(np.flip(np.linspace(0, 5.0, npoints))) - 1) / (np.exp(5.0)
35
36 # Properties sat functions: 1) swi [-], 2) sni [-], 3) pen [Pa], 4) penmax
37 safu = [[0.32, 0.1, 193531.39, 3e7, 1000],
38         [0.14, 0.1, 8654.99, 3e7, 1000],
39         [0.12, 0.1, 6120.00, 3e7, 1000],
40         [0.12, 0.1, 3870.63, 3e7, 1000],
41         [0.12, 0.1, 3060.00, 3e7, 1000],
42         [0.10, 0.1, 2560.18, 3e7, 1000],
43         [0, 0, 0, 3e7, 2]]
44
45 # Properties rock: 1) K [mD] and 2) phi [-], facie 1 to 7
46 rock = [[0.10132, 0.10],
47         [101.324, 0.20],
48         [202.650, 0.20],
49         [506.625, 0.20],
50         [1013.25, 0.25],
51         [2026.50, 0.35],
52         [1e-5, 1e-6]]
53
54 # Define the injection values ([hours] for spe11a; [years] for spe11b/c):
55 inj = [[999.9, 999.9, 100, 1, 0, 10, 1, 0, 10],
56        [ 0.1, 0.1, 0.1, 1, 0, 10, 1, 0, 10],
57        [ 25, 5, 5, 1, 50, 10, 1, 0, 10],
58        [ 25, 5, 5, 1, 50, 10, 1, 50, 10],
59        [ 50, 25, 25, 1, 0, 10, 1, 0, 10],
60        [ 400, 50, 50, 1, 0, 10, 1, 0, 10],
61        [ 500, 100, 100, 1, 0, 10, 1, 0, 10]]
```

SPE11 A

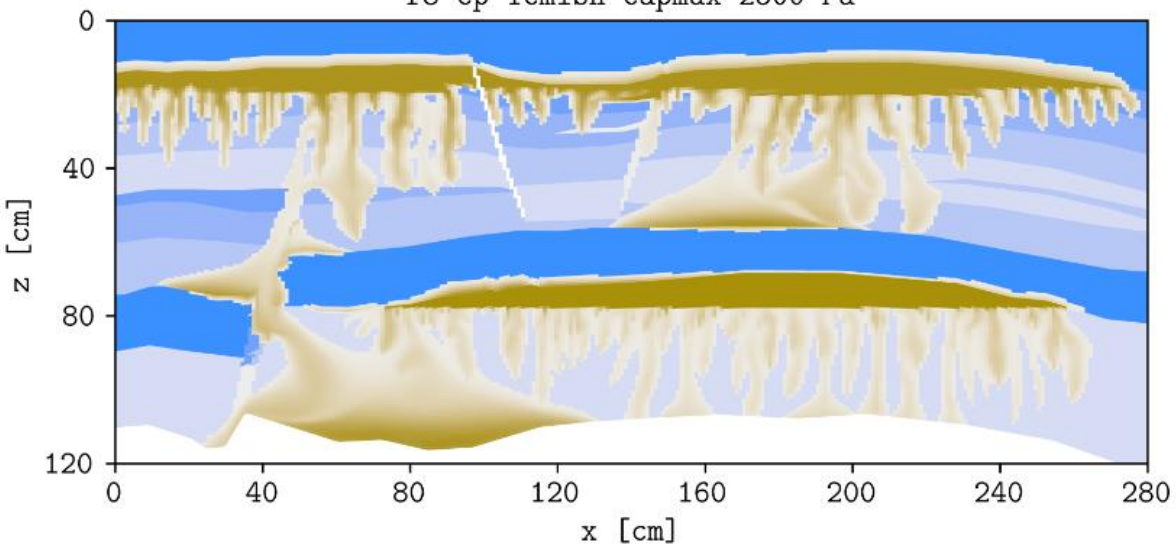
r1 Cart 1cm



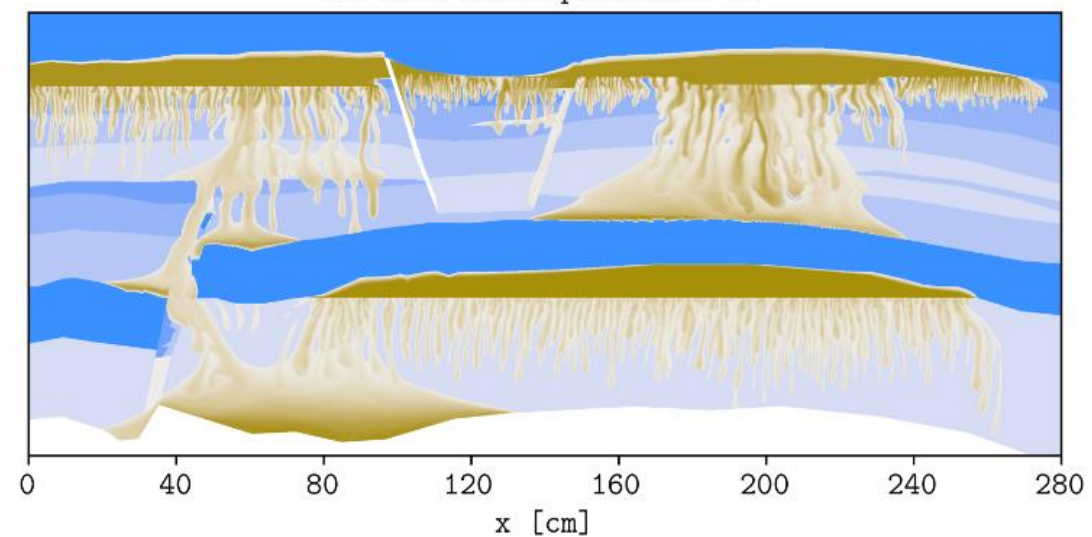
r2 Cart 1cm capmax 2500 Pa



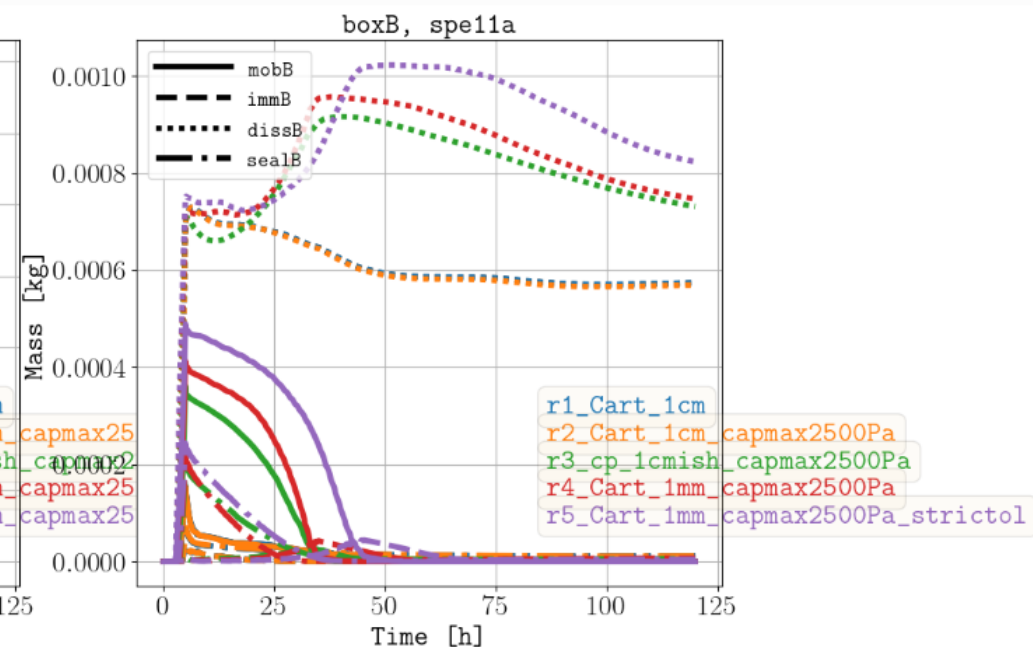
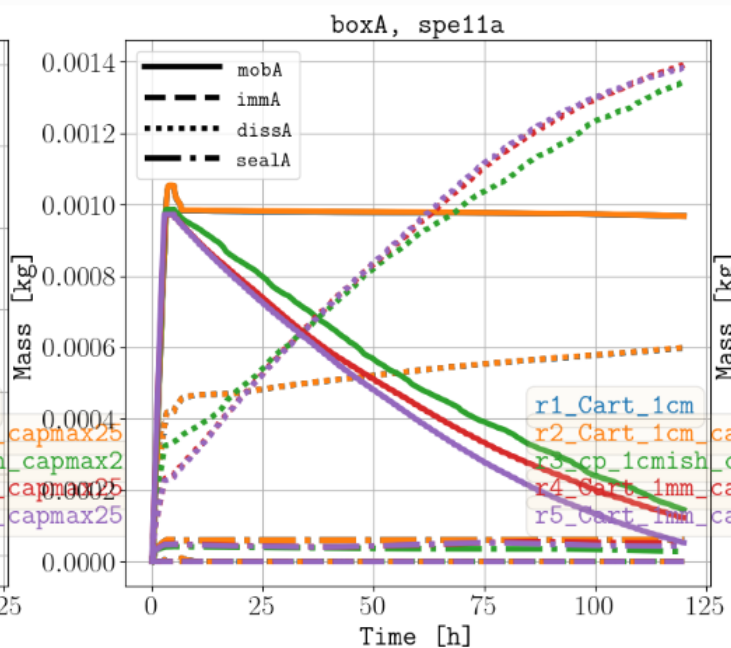
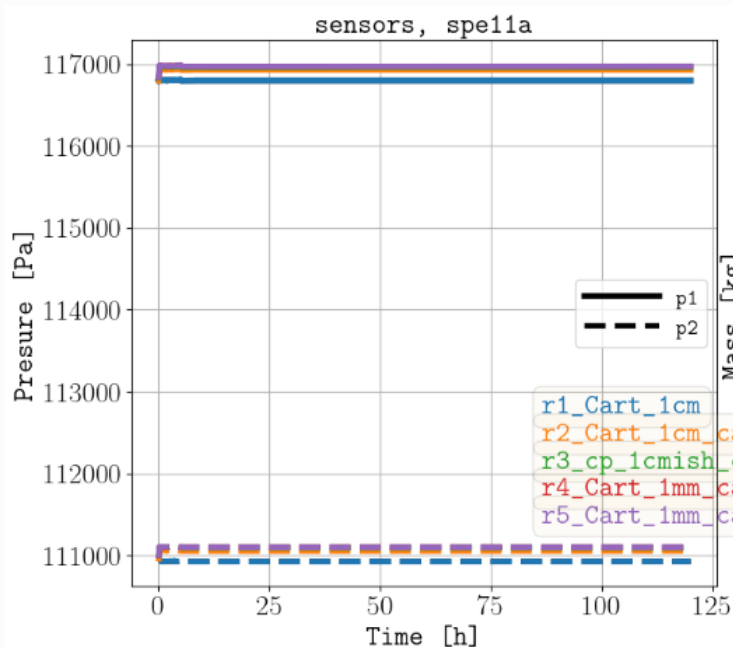
r3 cp 1cmish capmax 2500 Pa



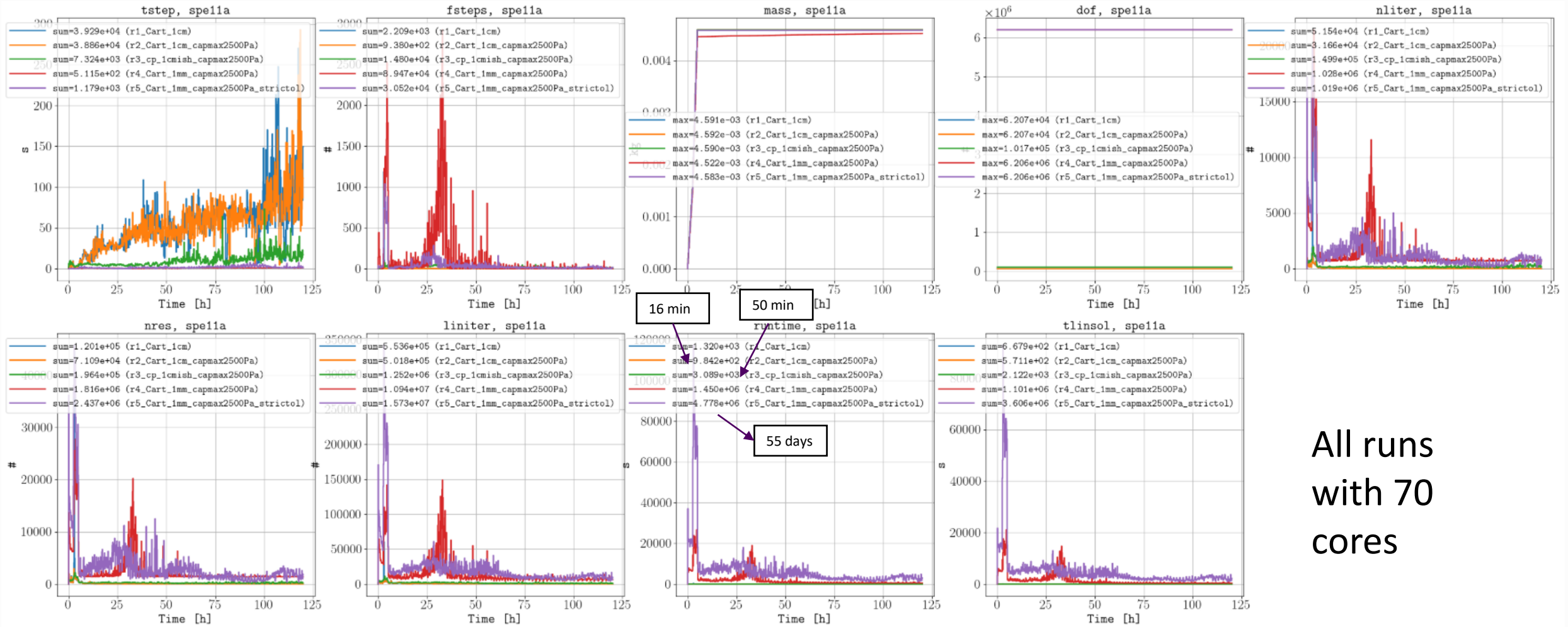
r4 Cart 1mm capmax 2500 Pa



SPE11 A



SPE11 A



All runs
with 70
cores

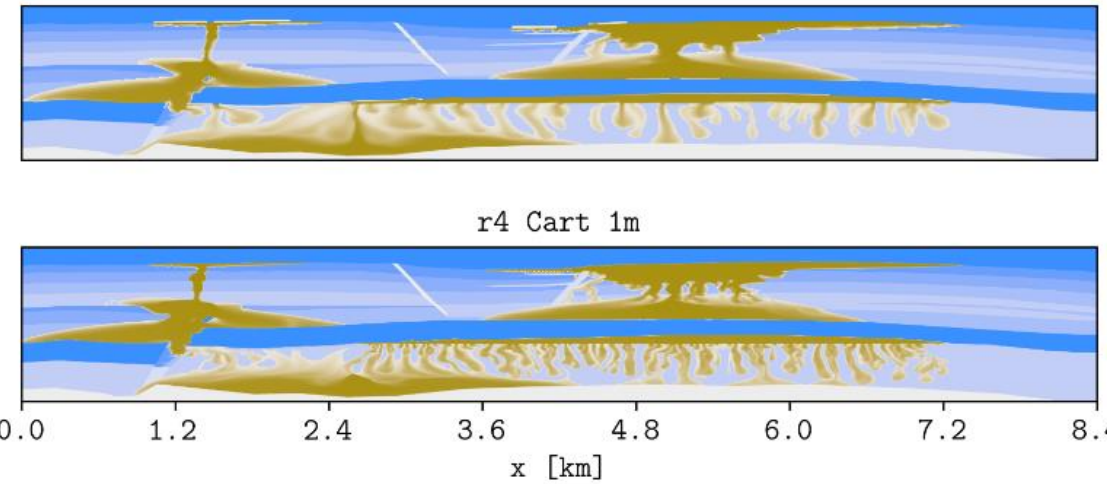
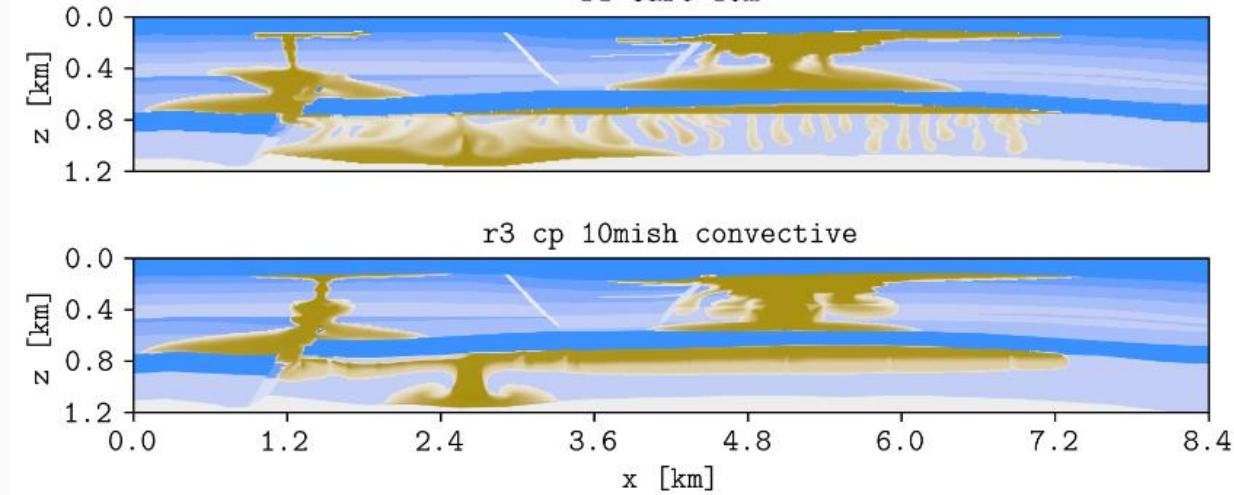
SPE 11 B

SPE11B: CO₂ mass fraction (liquid phase) after 500 years
0.00e+00 3.16e-02 6.31e-02



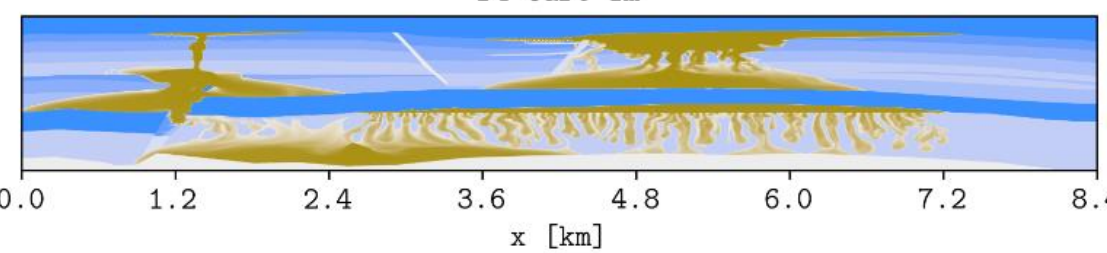
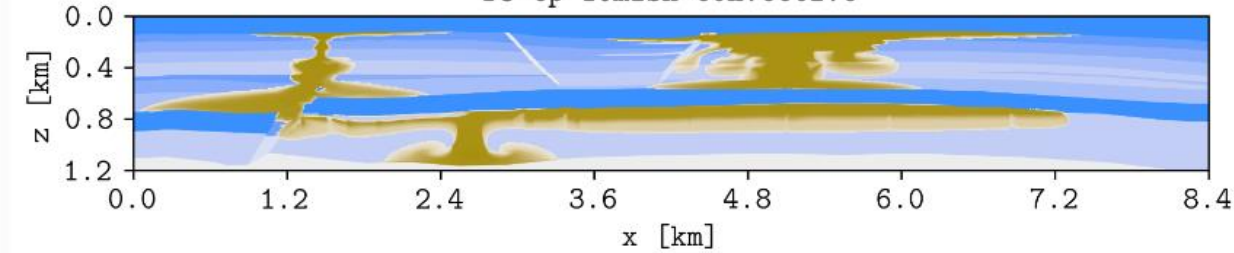
r1 Cart 10m

r2 cp 10mish



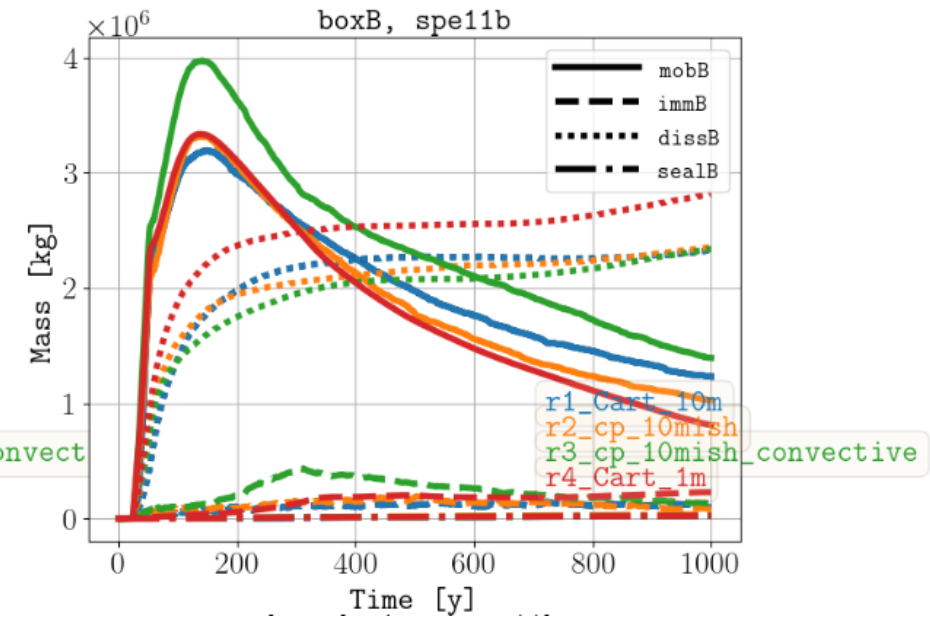
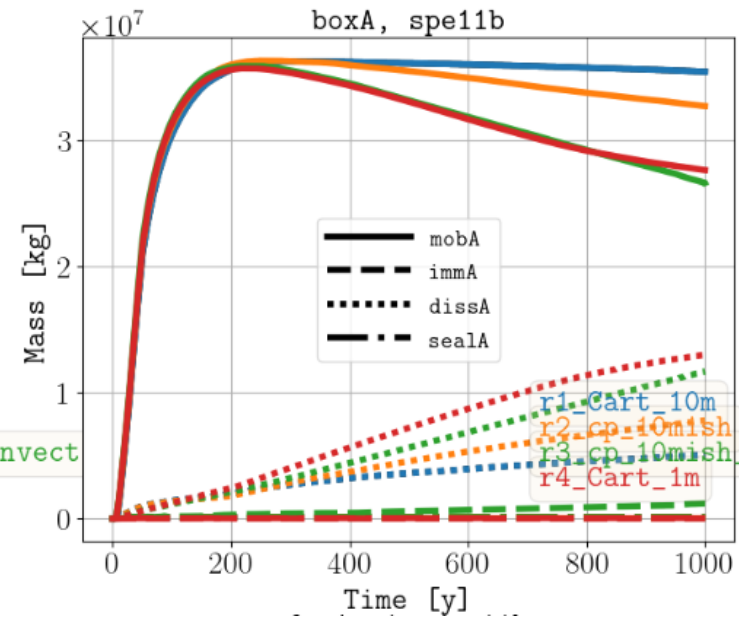
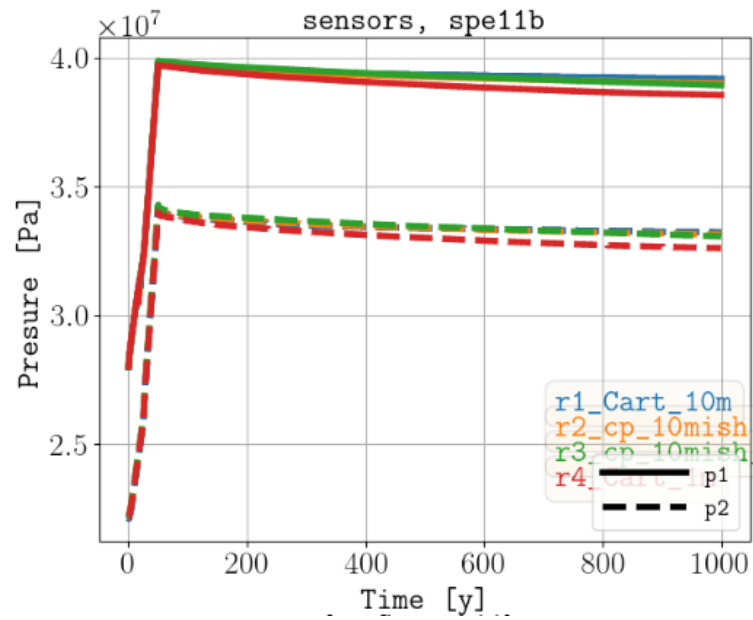
r3 cp 10mish convective

r4 Cart 1m

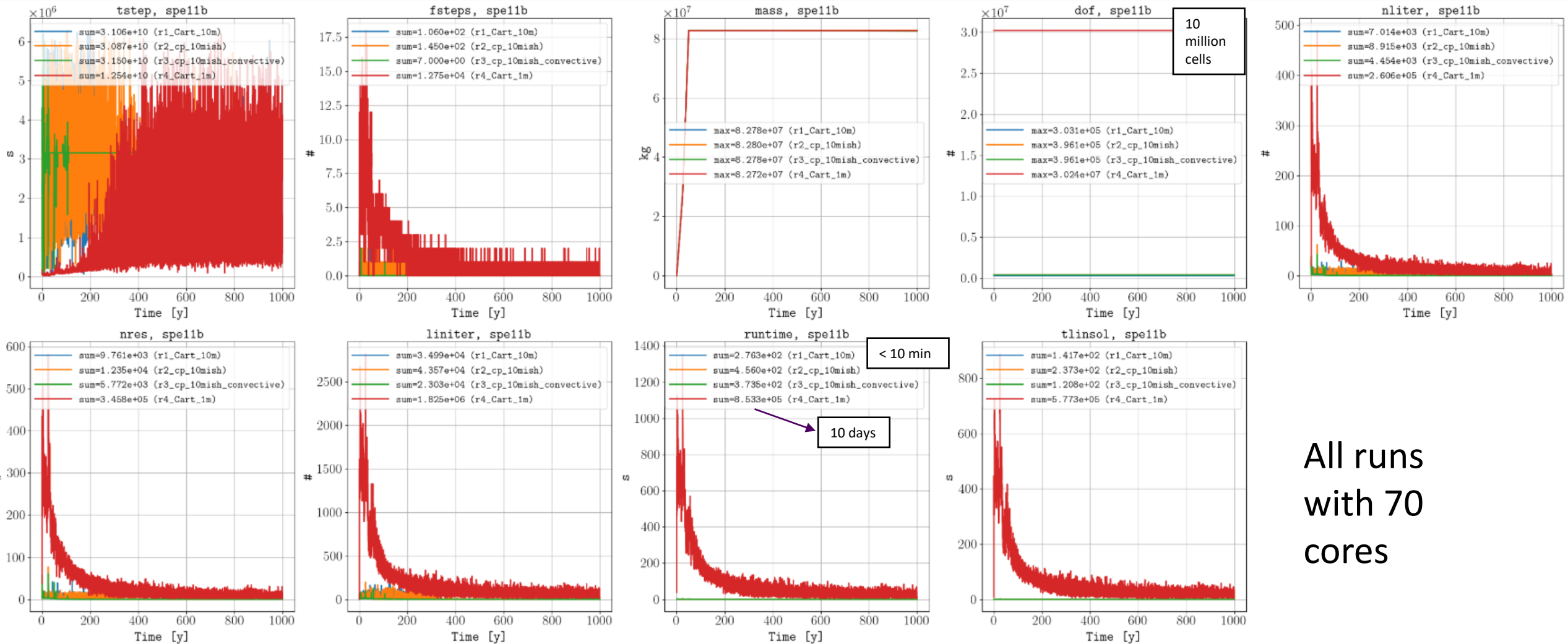


R3: "New sub-grid model for convective mixing in field-scale CO₂ storage simulation". Mykkeltvedt et al. In press TIMP 2024

SPE 11 B



SPE 11 B

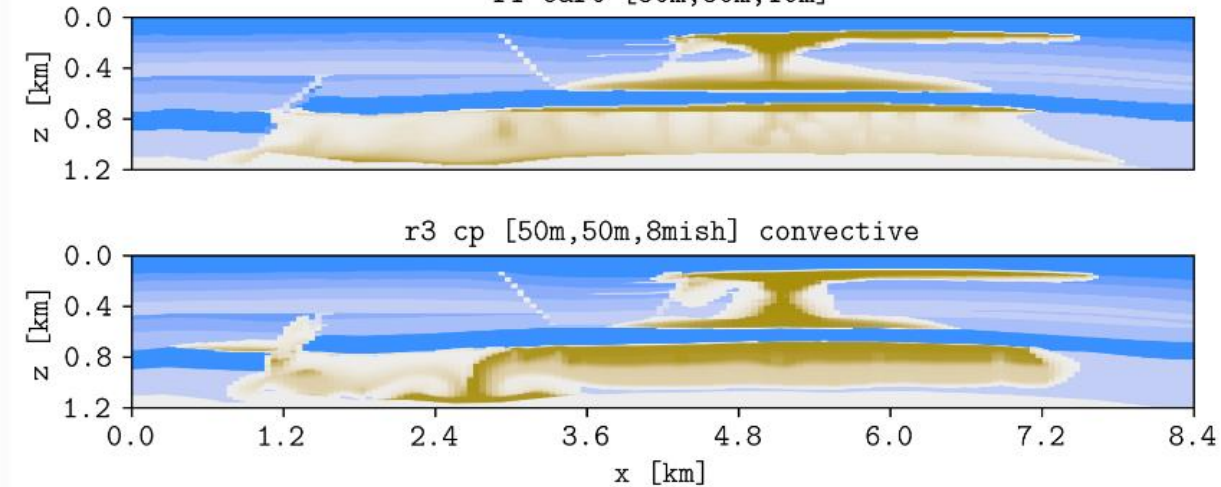


All runs with 70 cores

SPE11 C

SPE11C: CO₂ mass fraction (liquid phase) after 1000 years (y=2.5 [km])
0.00e+00 3.08e-02 6.16e-02

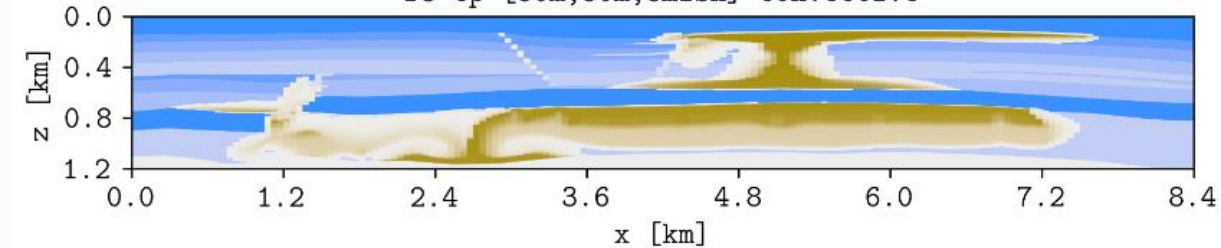
r1 Cart [50m,50m,10m]



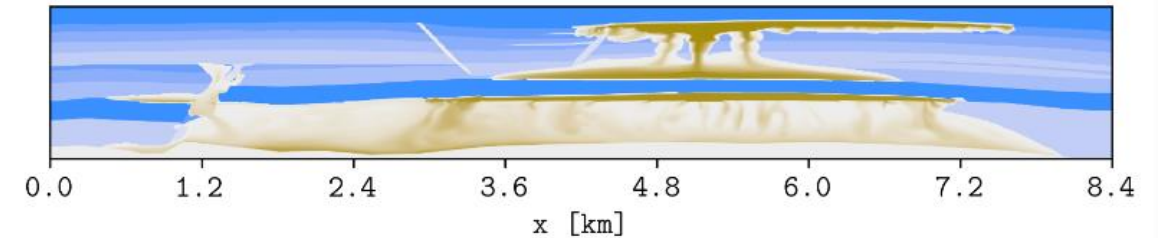
r2 cp [50m,50m,8mish]



r3 cp [50m,50m,8mish] convective

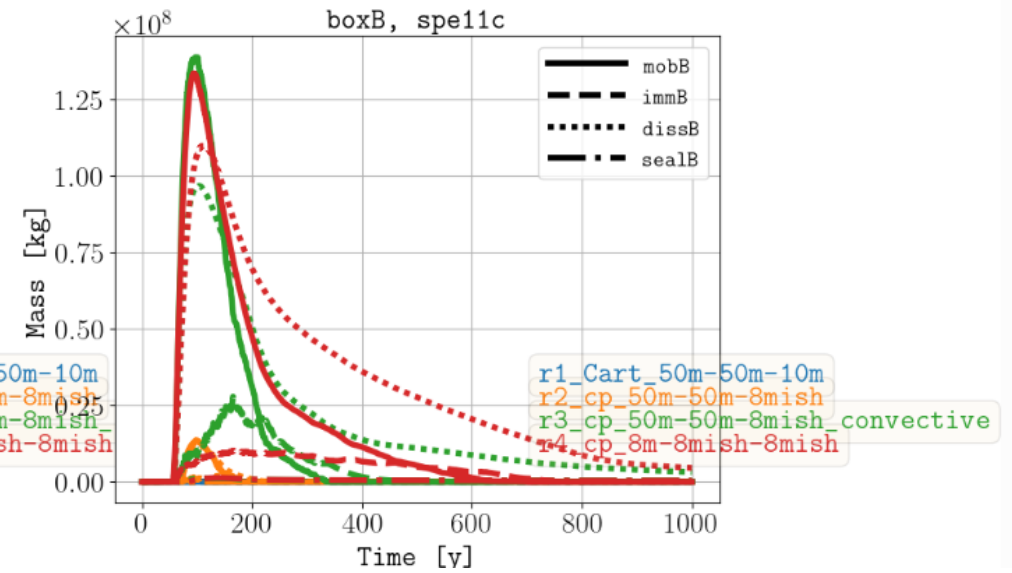
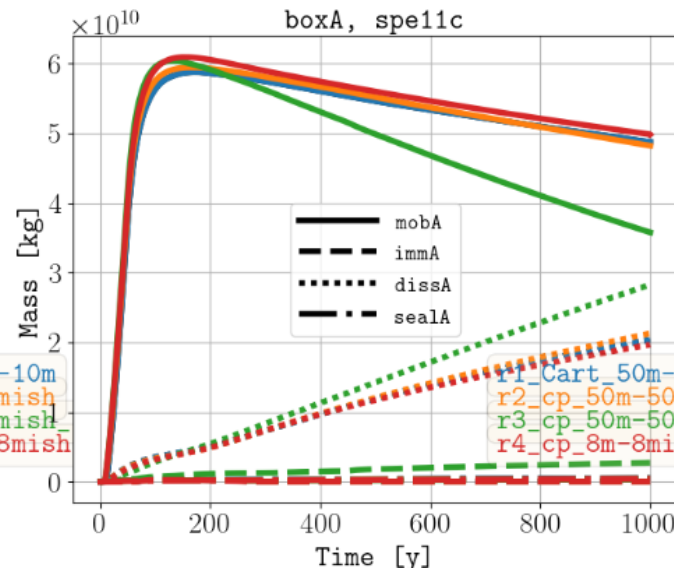
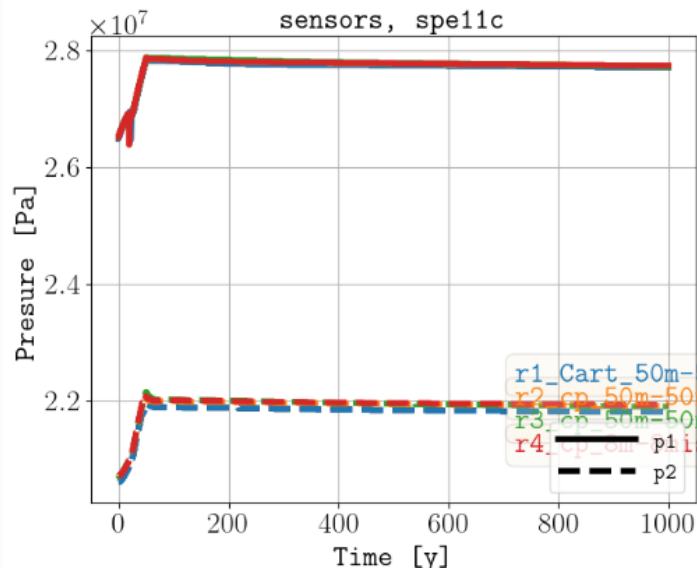


r4 cp [8m,8mish,8mish]

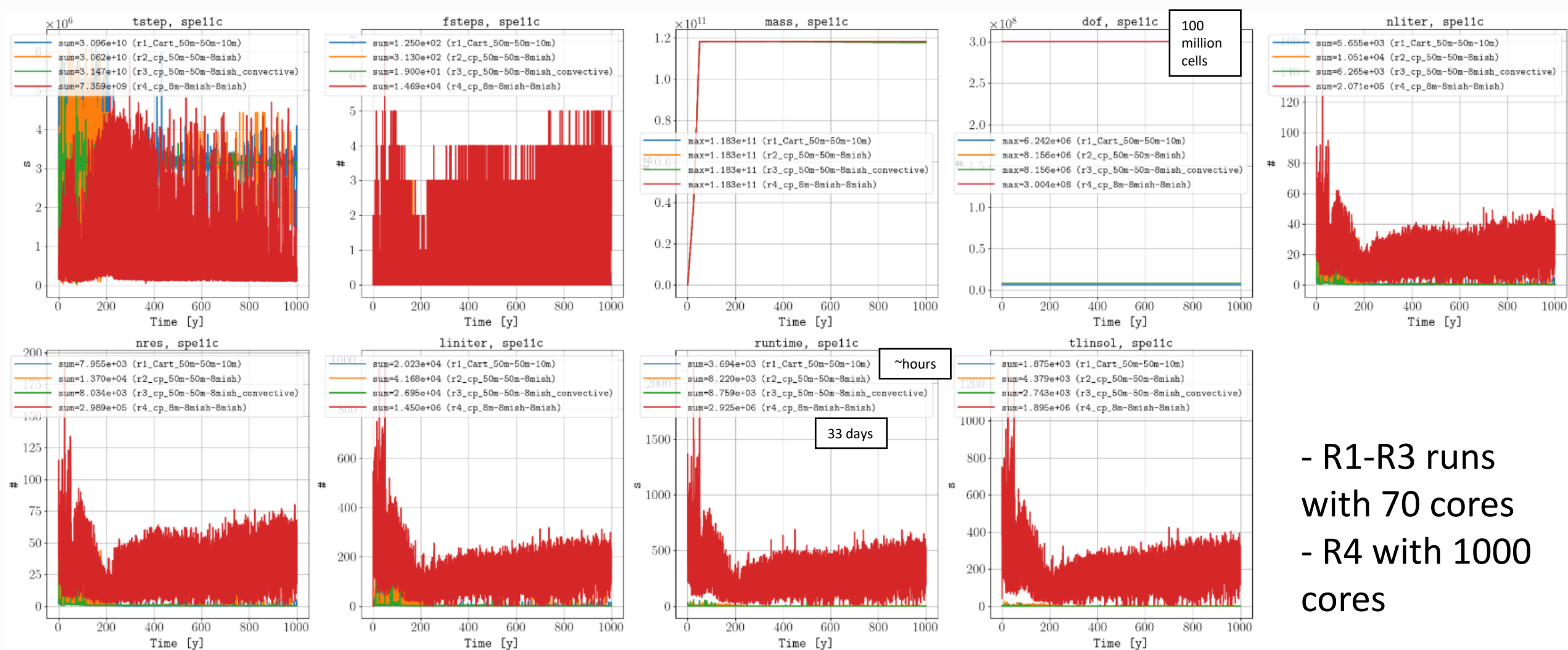


R3: "New sub-grid model for convective mixing in field-scale CO₂ storage simulation". Mykkeltvedt et al. In press TIMP 2024

SPE11 C



SPE11 C



- R1-R3 runs
with 70 cores
- R4 with 1000
cores

Summary

- [Pyopmspe11](#)
- CO2 simulations in OPM Flow using CO2STORE
- OPM Flow HPC simulations
- SGM for convective mixing (DRSDTCON)

The OPM team:

CLIMIT



GASSNOVA



CSSR



equinor

TNO innovation
for life



SINTEF

NORCE

OPM-OP AS