

OPM Flow, status and ongoing work

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Why is OPM Flow a success?







A few highlights and numbers

How large is OPM?



Language	Files	Blanks	Comments	Lines of code	Comment/LoC
C++	2902	135842	146860	528142	0.2781*
Python	72	1529	1519	6566	0.2313
CMake	117	1502	02 2613 1323		0.1975
JSON (keywords)	1192	15	0	32973	0

* Includes license block at top of files

C++ filesizes and comments over time





Probably too imprecise measures to make dramatic claims?

Commits and PRs



Repo	Commits	PRs	Janitorial PRs
Opm-common	1661	506	39
Opm-grid	430	109	8
Opm-simulators	3469	794	61

Data from April 9, 2024 – May 22, 2025

SPE11

The OPM Team:

- Only that submitted results for all cases
- Submitted the highest resolutions
 - Can now do ~170M cells (100M last year)
- Results similar to «middle of the pack»

Thanks for your effort!



Pau-Inria





A 3D plot of the CO2 mass fraction in the liquid phase is shown below. The cells in the reporting mesh have their transparency adjusted depending on the mass fraction of CO_2 in the liquid phase, with cells that have no CO_2 being

Industrial usage at Equinor



"Equinor is using OPM Flow on one asset and is evaluating it for use on more." (Equinor 2023)

"The black-oil simulation models are now run with OPM Flow in the assets to establish input to technical decisions." (Reiso et.al 2025)

4 oil assets and 1 CO₂ storage asset discussed in paper

Paper: Lessons Learned In Using Open-source Simulation Software On Real Asset Models



Talks by TBA and Edel Reiso today, Alf Birger Rustad tomorrow!



Recent and ongoing work

Improved nonlinear solvers



4000

Mean: 391.65

Median: 320.0 Min: 0 Max: 940

Rank 3

Newton — NLDD (default) — NLDD (tuned 1) — NLDD (tuned 2)

- Nonlinear domain decompositioning (NLDD) improvements:
 - Much better logging/reporting
 - More robust, works well for most cases out of the box
 - Ongoing: group controls in parallel

Runtimes for Sleipner with Newton and NLDD

200

12

16

20

24 28 Domain ID

4000

Paper: A distributed field-scale parallel nonlinear domain decomposition solver



Timestepping: fixes and new algorithms



- New timestepping algorithm added
- Existing ones bugfixed
- Ongoing work: using convergence diagnostics

Talk by **Erik Hide Sæternes** later today!

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-o- Implemented general third order controller for adaptive time d78cf2e stepping Projects \$\$ -o- Added missing safety factor paramet for time step controller 2be953f None yet	ξ
Image: Step controller on Feb 11 Milestone	ξ

GPU-accelerated linear solvers



- Performance improvements for ILUO and DILU, autotuning by default!
- Testing facilitated by gpu-istl (easy to test new components)
 - Wrapped GPU preconditioners to be CPU preconditioners
 - Wrapped CPU preconditioners to be GPU preconditioners
 - ... But wrappers no longer *required* -> better performance!
- CPR on GPU (two-level preconditioner)
- AMG on GPU (from AMGX, Hypre, and BDA)
- Well-operators: MS wells in progress, standard wells in prototype

Papers:

- gpu-ISTL Extending OPM Flow with GPU Linear Solvers
- A Comparison of DILU And ILU(0) as GPU-Accelerated Preconditioners



Talks by Vinicius Martins and Jakob Torben tomorrow!

GPU-accelerated properties

- Can now evaluate all properties of some models on GPU
 - CO2STORE, blackoil table input all supported
 - NOT supported: endpoint scaling, hysteresis, three-phase relperm, multiplexed pvt properties
 - Initial prototype target: SPE11-like cases
- Minimal changes to existing code
 - Typical change: templatize on storage (std::vector -> template argument)
- Almost identical code runs on CPU and GPU!

```
template<class TraitsT>
class PiecewiseLinearTwoPhaseMaterialParams
template <class TraitsT, class VectorT = std::vector<typename TraitsT::Scalar>>
class PiecewiseLinearTwoPhaseMaterialParams
template <class TraitsT, template <class> class Vector = std::vector>
class PiecewiseLinearTwoPhaseMaterialParams
```

Talk by **Tobias Meyer Andersen** tomorrow!

Local Grid Refinement (LGR)

- More flexible *parent cell sets*
 - not necessary block-shaped
 - not necessary all-active cells
 - LGRs sharing faces/corners is supported now
- MPI parallel CpGrid with LGRs in progress
- Wells aware of LGRs
- Initial output file support (EGRID)

Talks by Antonella Ritorto and Artur Castiel tomorrow!

Geomechanics

- Coupled flow and geomechanics, fractures
- Improved grid processing
 - Geomech (and thermal) grids must include zero-porosity cells
 - NNCs for flow mostly meaningless for mechanics

Fracture surface embedded in corner point geometry

Talk by Halvor Møll Nilsen later today!

Parallel performance

- Distributed multisegment wells
- Improved MPI partitioning
 - Revised graph: each well's connected cells are one single node
- Lock-free OpenMP/thread parallelization
 - Used for general assembly, tracer, some output
 - Makes high --threads-per-process more feasible

Assembly time same for 10 processes and 1 thread each and 1 process with 10 threads

Talks by Lisa Julia Nebel and Michal Tóth tomorrow!

Number of MPI processes:	1				
Threads per MPI process:	10				
Setup time:	0.68 s				
Deck input:	0.32 s				
Number of timesteps:	331				
Simulation time:	95.31 s				
Assembly time:	4.52	(Wasted:	0.0	s;	0.0%)
Well assembly:	0.70 s	(Wasted:	0.0	s;	0.0%)
Linear solve time:	42.57 s	(Wasted:	0.0	s;	0.0%)
Linear setup:	19.35 s	(Wasted:	0.0	s;	0.0%)
Props/update time:	4.88 s	(Wasted:	0.0	s;	0.0%)
Pre/post step:	42.19 s	(Wasted:	0.0	s;	0.0%)
Output write time:	0.27 s				
Overall Linearizations:	1374	(Wasted:		0;	0.0%)
Overall Newton Iterations:	1043	(Wasted:		0;	0.0%)
Overall Linear Iterations:	1919	(Wasted:		0;	0.0%)

Number of MPI processes:	10			
Threads per MPI process:	1			
Setup time:	1.00 s			
Deck input:	0.50 s			
Number of timesteps:	333			
Simulation time:	28.62 s			
Assembly time:	4.47 5	(Wasted:	0.0 s;	0.2%)
Well assembly:	0.40 s	(Wasted:	0.0 s;	0.2%)
Linear solve time:	9.79 s	(Wasted:	0.1 s;	0.8%)
Linear setup:	4.22 s	(Wasted:	0.0 s;	0.2%)
Props/update time:	2.77 s	(Wasted:	0.0 s;	0.1%)
Pre/post step:	8.40 s	(Wasted:	0.0 s;	0.0%)
Output write time:	1.78 s			
Overall Linearizations:	1387	(Wasted:	2;	0.1%)
Overall Newton Iterations:	1055	(Wasted:	2;	0.2%)
Overall Linear Iterations:	2202	(Wasted:	34;	1.5%)

CO₂ simulation

- Advanced modeling capabilities, f.ex:
 - Salt precipitation
 - Upcaling of convecting mixing
- Good performance and scaling
- Successful SPE11 participation!

Example papers:

New Sub-grid Model for Convective Mixing in Field-Scale Storage Simulation

Impact of Intermittency on Salt Precipitation During CO2 Injection

Talks by **Tor Harald Sandve** today and **Trine Mykkeltvedt** tomorrow!

Brine density from SPE11C, 101-million cell simulation

Robustness, wells, groups++

- Improved PVT extrapolation at low pressures
 - Avoid unphysical behaviour
- Improved initialization of wells and network
- Revised algorithms for finding solution before full Newton iteration
 - Dozens of small changes
 - A lot of ongoing work in this area
- Completing features (example: GSATPROD)
- Robustness and performance go together!

Well cells from an OLYMPUS case

Compositional simulation

- General compositional simulation!
 - Arbitrary number of components
 - Additional water phase
 - Multiple equations of state (EOS) available
 - Standard well model implemented (some limitations for now)
- Output to VTK and Eclipse formats

1D validation of compositional behaviour

Talk by Kai Bao later today!

... and much more!

- Input/Output improvements
 - Much improved consistency checks for saturation functions
 - Support updated restart format (E100 version 2022.4)
 - Some region support for UDQs
 - UDQ checks and validation
- Reservoir coupling
 - Prototype available
- Integrated machine learning
 - can replace arbitrary relation with ML model (some assembly required...)
- Python coupling, API and tools
 - PET, Pycopm, plopm, pyopmspe11++, python docs

Talks by **Rolf Johan Lorentzen** later today and **David Landa Marban** tomorrow!

Technology for a better society