The MATLAB Reservoir Simulation Toolbox

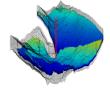
Bård Skaflestad

SIAM Geosciences, Long Beach, March 21-24, 2011

The Matlab Reservoir Simulation Toolbox (MRST)

The toolbox has the following functionality for rapid prototyping of solvers for flow and transport:

- grid structure, grid factory routines, input/processing of industry-standard formats, real-life and synthetic example grids
- petrophysical parameters and incompressible fluid models, conversion routines to/from SI and common field units, very simplified geostatistical routines
- routines for setting and manipulating boundary conditions, sources/sinks, and well models
- reservoir state (pressure, fluxes, saturations, compositions, ...)



visualisation routines for cell and face data (scalars)

Download

http://www.sintef.no/MRST/

Version 2011a was released on the 22nd of February, 2011, and can be downloaded under the terms of the GNU General Public License (GPL)

- General framework for flow and transport in porous media
- Special focus on unstructured grids and multiscale methods

Research:

- Development of new solvers and discretization schemes
- Application to new scientific/engineering problems
- Preserve know-how and promote reuse of results from previous research
- Contribute to accelerating production of new research (by our peers)
- Benchmarking compare different methods on standard test problems Students:
 - Develop students' intuition of porous media flow
 - Make it easy to test, compare, and extend existing methods
 - Textbook (with worked examples) in preparation

Why open-source and why in Matlab?

First of all, prototyping in a scripting language is much more effective than in traditional compiled languages (C/C++/FORTRAN)

- Explore alternative algorithms/implementations close to mathematics
- Gradually replace individual (or bottleneck) operations with accelerated editions callable from MATLAB
- Direct access to MATLAB environment and prototype whilst developing replacement components
- More experienced in Matlab/Octave than in e.g., Python,

Why free and open-source software:

- Research funded by Norwegian Research Council should be freely available
- Combined with publications: a means to collaborate and disseminate results, while protecting IP rights
- Support reproducible research and promote replicability

The fundamental object in MRST is the grid:

- Data structure for geometry and topology
- Several grid factory routines
- Input of industry-standard (proprietary) format(s)

Physical quantities defined as dynamic objects in MATLAB

- Properties of medium (ϕ , **K**, net-to-gross, ...)
- Reservoir fluids (ρ , μ , k_r , PVT, ...)
- Driving forces (wells, boundary conditions, sources)
- Reservoir state (pressure, fluxes, saturations, etc)

All MRST operations accept, manipulate and produce objects of these types. Physical quantities are assumed to be in SI units.

MRST core

- routines for creating and manipulating grids and physical properties
- basic flow and transport solvers (sequential splitting) for incompressible and immiscible flow

Functionality is stable and not expected to change in future releases

Modules

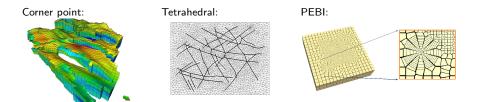
Similar to ${\rm MATLAB}{}^{\prime}{\rm s}$ toolboxes. Implements more advanced solvers and tools:

- adjoint methods, experimental multiscale, fractures, MPFA, upscaling
- black-oil models, three-phase flow, vertically integrated models, ...
- streamlines, (flow-based) coarsening, ...
- Octave support, C-acceleration, ...

Some are stable. Some are constantly changing to support ongoing research. *New modules initiated by others are much welcome*

Grids in MRST

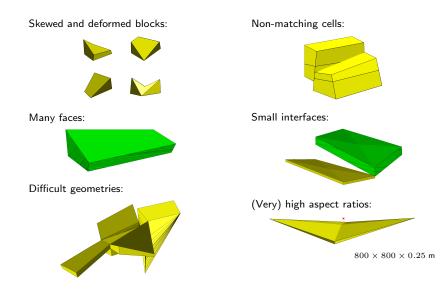
Complex reservoir geometries



- Accurate modelling of features such as faults, fractures, or erosion requires grids that are flexible with respect to geometry.
- Industry-standard grids are often nonconforming and contain complex grid-cell connectivities, as well as skewed and degenerate cells
- There is a trend towards unstructured grids with general polyhedral cells
- Standard discretization methods produce wrong results on skewed and rough cells condition numbers
- The representation of (un)structured grids affects the efficiency of numerical methods

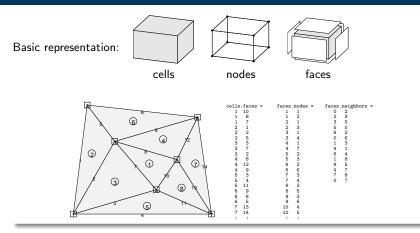
Grids in MRST

Cell geometries are challenging from a discretization point-of-view



Grids in MRST

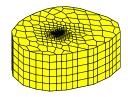
All grids are assumed to be unstructured

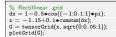


Choices in grid representation guided by utility and convenience in low-order finite-volume methods. Available geometric information typically limited to centroids, normals, areas, and volumes

Grids in MRST Examples of basic grid types

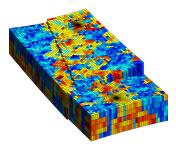
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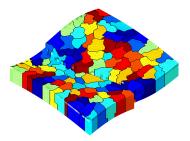
% Extrude a standard MATLAB dataset load seamount g = triangleGrid([x(:) y(:)]); P = pebi(g); V = makeLayeredGrid(P, 5); plotGrid(V), view(-40, 60), axis off % Make and read a simple Eclipse input file grdec1 = simpleGrdec1[[20, 10, 5], 0, 12] 6 6 = processGRBCE[grdec1], plotGrid(G,¹FaceAlpha¹,0.8); plotFacea(G,find(G.faces.tag>0), ¹FaceColor¹,¹red¹); view(40,40), axis off

Grids in MRST Examples of flexible gridding strategies



Hybrid grid – areal grid consisting of the following components

- radially refined grid at wells
- Cartesian along boundary
- hexahedral in interior
- polyhedral transition cells
 extruded to 3D along vertical lines



Hierarchical grid

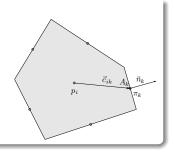
- triangular cells adapted to a curved line
- extruded to 3D, with throw along curved surface
- coarse grid constructed by collecting cells from fine grid

Discretization of flow equation General family of conservative methods for: $\nabla \cdot \vec{v} = q$, $\vec{v} = -\mathbf{K}\nabla p$

Basic formulation

$$\boldsymbol{u}_i = \boldsymbol{T}_i (\boldsymbol{e}_i p_i - \boldsymbol{\pi}_i), \qquad \boldsymbol{e}_i = (1, \dots, 1)^{\mathsf{T}}$$

 p_i – the pressure at the center of cell i u_i – the vector of outward face fluxes π_i – the vector of face pressures T_i – the one-sided transmissibilities

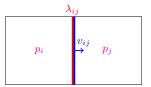


Special cases:

- The standard two-point method: $m{T}_{ii} = ec{n}_i \cdot m{K} ec{c}_i / |ec{c}_i|^2$
- Multipoint flux-approximation methods (MPFA)
- Mixed finite-element methods
- Mimetic methods

Linear system: mixed hybrid form

$$\begin{bmatrix} B & C & D \\ C^T & 0 & 0 \\ D^T & 0 & 0 \end{bmatrix} \begin{bmatrix} v \\ -p \\ \pi \end{bmatrix} = \begin{bmatrix} 0 \\ g \\ 0 \end{bmatrix},$$

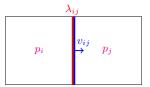


B defines an inner product. The matrix blocks read,

$$b_{ij} = \int_{\Omega} \psi_i \mathbf{T}^{-1} \psi_j \, dx, \quad c_{ik} = \int_{\Omega} \phi_k \nabla \cdot \psi_i \, dx, \quad d_{ik} = \int_{\partial \Omega} |\psi_i \cdot n_k| \, dx$$

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Positive-definite system obtained by a Schur-complement reduction

$$(D^T B^{-1} D - F^T L^{-1} F) \pi = F^T L^{-1} g,$$

 $F = C^T B^{-1} D, \quad L = C^T B^{-1} C.$

Reconstruct cell pressures and fluxes by back-substition,

$$Lp = q + F^T \pi, \qquad Bv = Cp - D\pi.$$

Herein: a mimetic method, Brezzi et al., 2005

$$Mu = (ep - \pi) \quad \longleftrightarrow \quad u = T(ep - \pi)$$

Requiring exact solution of linear flow $(p = x^{T}a + k)$:

$$MNK = C$$
 $NK = TC$

C - vectors from cell to face centroids. N - area-weighted normal vectors

Family of schemes (given by explicit formulas):

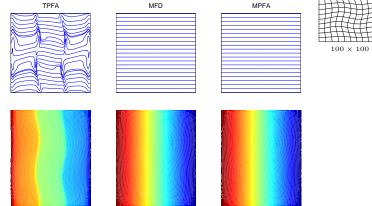
$$egin{aligned} M &= rac{1}{|\Omega_i|} oldsymbol{C} oldsymbol{K}^{-1} oldsymbol{C}^{\mathsf{T}} + oldsymbol{Q}_N^{\perp}^{\mathsf{T}} oldsymbol{S}_M oldsymbol{Q}_N^{\perp} \ T &= rac{1}{|\Omega_i|} oldsymbol{N} oldsymbol{K} oldsymbol{N}^{\mathsf{T}} + oldsymbol{Q}_C^{\perp}^{\mathsf{T}} oldsymbol{S} oldsymbol{Q}_C^{\perp} \end{aligned}$$

 Q_N^\perp is an orthonormal basis for the null space of N^\top , and S_M is any positive definite matrix. Herein, we use null-space projection

$$\boldsymbol{P}_{N}^{\perp} = \boldsymbol{Q}_{N}^{\perp} \boldsymbol{S}_{M} \boldsymbol{Q}_{N}^{\perp} = \boldsymbol{I} - \boldsymbol{Q}_{N} \boldsymbol{Q}_{N}^{\top}$$

Example: grid orientation effects

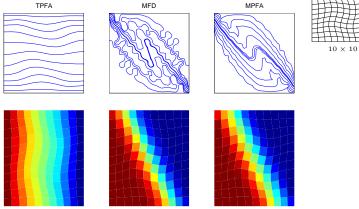
Homogeneous domain with Dirichlet boundary conditions (left,right) and no-flow conditions (top, bottom) computed with three different pressure solvers in MRST.



Homogeneous permeability with anisotropy ratio 1:1000 aligned with the grid.

Example: lack of monotonicity

Homogeneous domain with Dirichlet boundary conditions (left,right) and no-flow conditions (top, bottom) computed with three different pressure solvers in MRST.



Homogeneous permeability with anisotropy ratio 1:1000 rotated by $\pi/6$.

A (very) simple flow solver

$$\nabla\cdot\vec{v}=q,\qquad \vec{v}=-\frac{\mathbf{K}}{\mu}\big[\nabla p+\rho g\nabla z\big]$$

Vertical well and Dirichlet boundary

```
% Grid and rock parameters
nx = 20; nv = 20; nz = 10;
G = computeGeometry(cartGrid([nx, ny, nz]));
rock.perm = repmat(100 * milli*darcy, [G.cells.num, 1]);
fluid = initSingleFluid('mu', 1*centi*poise, ...
                      'rho', 1014*kilogram/meter^3);
gravity reset on
% Fluid sources and boundary conditions
    = (nx/2*ny+nx/2:nx*ny:nx*ny*nz).';
src = addSource([], c, ones(size(c)) ./ day());
bc = pside([], G, 'LEFT', 10*barsa());
% Construct components for mimetic system
S = computeMimeticIP(G, rock, 'Verbose', true);
% Solve the system and convert to bars
rSol = initResSol(G, 0);
rSol = solveIncompFlow(rSol.G.S.fluid.'src'.src.'bc'.bc);
p = convertTo(rSol.pressure(1:G.cells.num), barsa());
```

Source term and boundary condition Pressure distribution

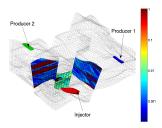
From tutorial: simpleSRCandBC.m

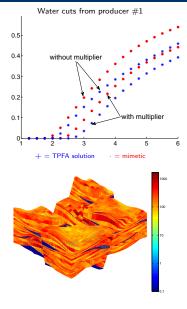
Modelling faults in consistent schemes

Faults modelled as internal boundaries, with internal jump conditions

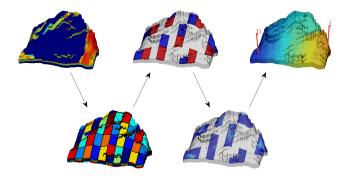
$$u_f^{\pm} = T_f(\pi_f^{\mp} - \pi_f^{\pm})$$

Gives an extended hybrid system. In addition, method to convert TPFA multipliers to fault transmissibility $T_{\rm f}$





Multiscale module: bypassing the need for upscaling?



Key idea of multiscale methods:

Local decomposition

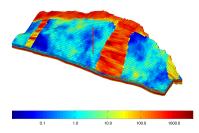
$$\begin{array}{l} \blacktriangleright \ p(\vec{x}) = \sum_{\Omega_i} p_i \phi_i(\vec{x}) \\ \hline \ \vec{v}(\vec{x}) = \sum_{\Gamma_{ij}} v_{ij} \vec{\psi}_{ij}(\vec{x}) \\ \Omega_i - \text{coarse grid block. } \Gamma_{ij} = \partial \Omega_i \cap \partial \Omega_j \end{array}$$

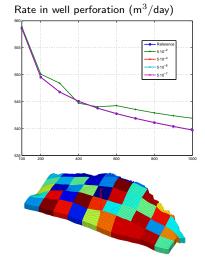
Multiscale basis functions:



Multiscale module / black-oil module: primary production

- Shallow-marine reservoir (realization from SAIGUP)
- Model size: $40 \times 120 \times 20$
- Initially filled with gas, 200 bar
- Single producer, bhp=150 bar
- Multiscale solution for different tolerences compared with fine-scale reference solution.





Coarsening module: (flow-based) coarsening by amalgamation

Amalgamation of cells

- flow-adapted grids
- simple and flexible coarsening
- adaptive gridding schemes
- efficient model reduction

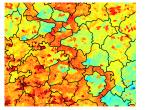
Algorithmic primitives working on a partition vector

- segment (flow) indicator into bins
- merge small cells
- refine large cells
- intersect partitions
- sanity checks to ensure connected blocks, etc
- \longrightarrow great flexibility

Cartesian grid:

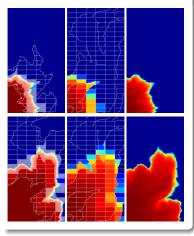


Triangular grids:

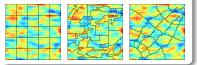


Coarsening module: (flow-based) coarsening by amalgamation

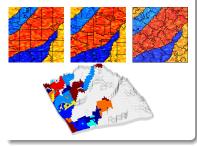
Dynamic adaption



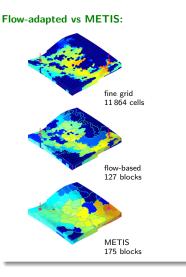
Different partitioning:



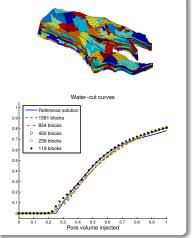
Adapting to geology



Coarsening module: (flow-based) coarsening by amalgamation

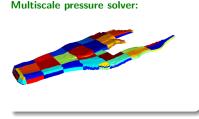


Model reduction of real-field model:

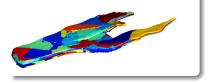


Upcoming adjoint module: production optimization

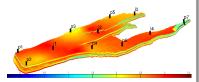
Specialized simulator: using different grids for pressure and transport



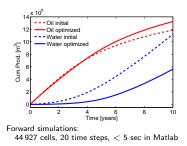
Transport on flow-adapted grid:



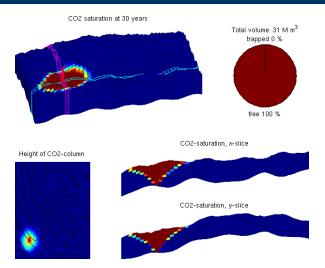
Water-flood optimization:



Reservoir geometry from a Norwegian Sea field

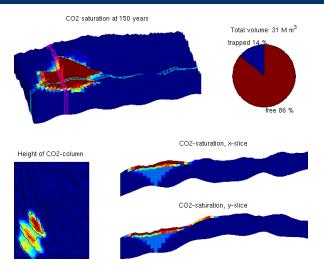


Vertically integrated module: educational and research tool



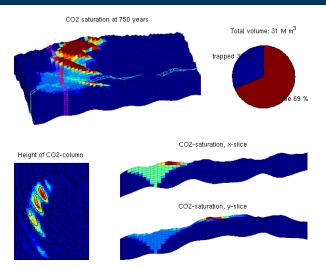
Vertically integrated module: part of a forthcoming numerical CO_2 laboratory

Vertically integrated module: educational and research tool



Vertically integrated module: part of a forthcoming numerical CO₂ laboratory

Vertically integrated module: educational and research tool



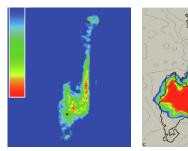
Vertically integrated module: part of a forthcoming numerical CO₂ laboratory

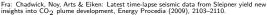
Vertically integrated module: CO2 migration at Sleipner

3D model made by Statoil to study CO_2 in the top layer of Utsira. MRST used to study physical assumptions and numerical methods:

- Simple to modify the code
- Flexible and simple upscaling (homogeneous model)
- Fast response time, simulation 2–10 min in Matlab

Seismic data (2006) / 3D simulation (tough2)







VE simulation

