Localized Linearization

Andreas Lauser

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Caches are small, memory is slow, let's do it on the GPU!

Or not?!











2 Reservoir Simulations

Field Discretization



Scalar field: Function $f : \Omega \mapsto \mathcal{R}$ were $\Omega \subseteq \mathcal{R}^n$ for some $n \in \mathcal{N}$

Usual discretization approach:

- Define the value of the field at a finite number of degrees of freedom (DOFs)
- Interpolate in-between
 - Approximates field by a vector of scalars





Partitioning Ω of into smaller sets ω_i :

- Sub-domains ω_i usually exhibit regular shape
 - Simplices, parallelepipeds, etc.
- ω_i usually called *element* or *cell*
- In the following: Cell-centered finite volume method (i.e., one DOF per cell)



Left hand side:

- Iterate over all DOFs, calculate the values f(x) and g(x) and combine them locally
- (in the following called "element based/localized approach")

Right hand side:

- Calculate all values of *f*, then all of *g*, then combine them
- (in the following called "grid based/global approach")

Consider the following:

$$egin{aligned} f(\mathbf{x}) &= |\mathbf{x}|\,,\ \star &\equiv \cdot, \end{aligned}$$

Localized:

```
for i in cells:
    x = abs(pos(i))
    f[i] = x
    g[i] = sqrt(x)
    h[i] = f[i] * g[i]
```

$$egin{aligned} g(\mathbf{x}) &= \sqrt{|\mathbf{x}|} \ h(\mathbf{x}) &= |\mathbf{x}| \sqrt{|\mathbf{x}|} \end{aligned}$$

Global:

```
for i in cells:
    f[i] = abs(pos(i))
for i in cells:
    g[i] = sqrt(abs(pos(i)))
for i in cells:
    h[i] = f[i] * q[i]
```

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¹https://poware.de/adue3aib/global-vs-local.tar.gz

- Both approaches also applicable for more complicated functions (e.g., relperms) and operators (e.g., gradients).
- Due to prefetching, performance advantage of localized approach is smaller for more expensive *f* or *g*
- Very often, function to be computed is very cheap
 - e.g. *h* = 1 − *f*
 - (Off topic: Iterative linear solvers particularly affected)

Advantages of the localized approach:

- Better cache locality
- Easier to parallelize
 - Only a single loop

Advantages of the global approach:

- More modular representation of the underlying equations
 - Simplifies/allows implementation of special purpose PDE languages like *DOLFIN*² or *Equelle*³

²http://fenicsproject.org/ ³http://equelle.org/







OPM is one of the few projects which where an "apples-to-apples" comparison can be attempted:

- opm-autodiff uses the global approach to linearization
- eWoms/ebos uses the localized approach
- Grid and deck input parameters identical/very similar

Other differences in implementation are pretty large:

- All tricks like partial relinearization and linearization rescaling are disabled
- Results should be taken with a grain of salt!

Boring Problem: SPE-1



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Large Boring Problem: Refined SPE1





BHP of the "PRODUCER" well vs. time

Performance: Apples and Bananas



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Overall time of the refined SPE1 problem. Ratio: 1.98

Performance: Apples and Bananas



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Overall time of the refined SPE1 problem. Ratio: 1.98

 Linear solver also suffers from "memory bandwith bottleneck"

Performance: Apples and Oranges



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Aggregate linearization time of the refined SPE1 problem. Ratio: 1.43

Performance: Apples and Oranges



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Aggregate linearization time of the refined SPE1 problem. Ratio: 1.43

• opm-autdiff requires a few iterations more

Performance: Green Apples and Red Apples



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Iteration adjusted linearization time for the refined SPE1 problem. Ratio: 1.26

Performance: Green Apples and Red Apples



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Iteration adjusted linearization time for the refined SPE1 problem.

Ratio: 1.26

opm-autodiff uses automatic differentiation, eWoms finite differences



- Performance comparisons are hard
- Quality of results of all simulators comparable for the unrefined and refined SPE-1 problems.
- Localized linearization approach saw better performance
 - Likely faster in principle because it reduces the "RAM bottleneck" problem



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Thank you for your attention.