



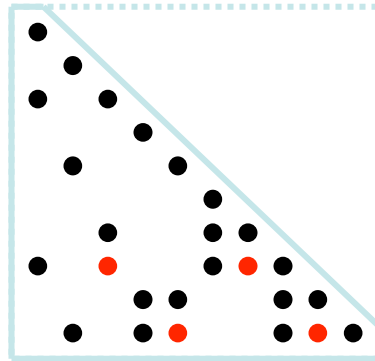
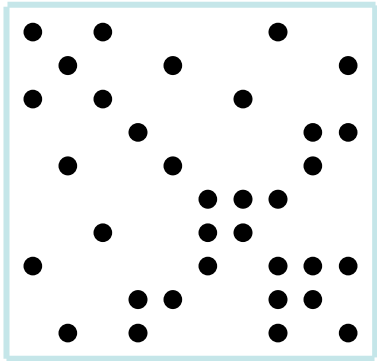
Empirical Complexity of Laplacian Linear Solvers: Discussion

Erik Boman, Sandia National Labs
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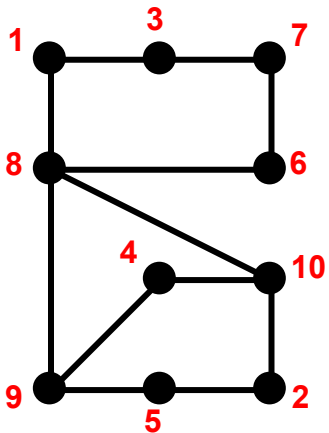
Simons Institute Workshop on
Fast Algorithms via Spectral Methods
December 2, 2014

Support: Intel, Microsoft, DOE Office of Science, NSF

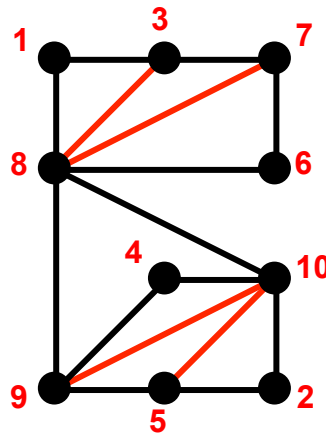
Graphs and sparse matrix computation (philosophical digression)



Symmetric Gaussian elimination:
for $j = 1$ to n
add edges between j 's
higher-numbered neighbors



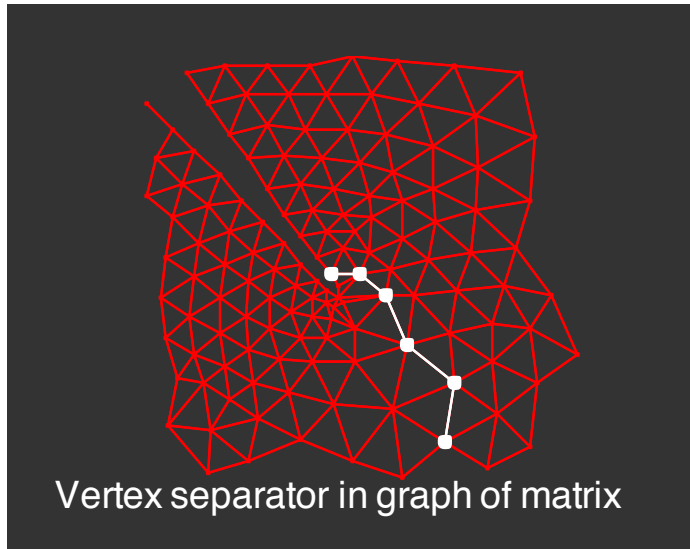
A



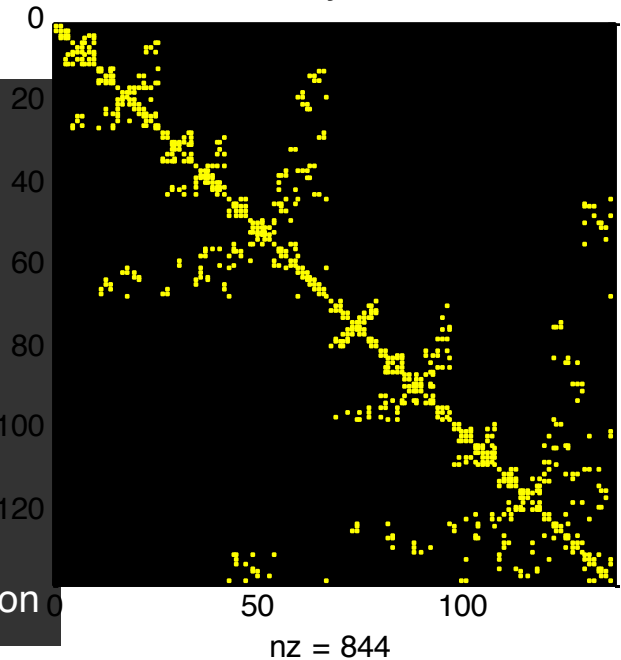
L (chordal)

Nested dissection and graph partitioning

[George 1971, then many papers]



Matrix reordered by nested dissection



- Find a small vertex separator, number it last, recurse on subgraphs
- Approx optimal separators \Rightarrow approx optimal fill & flop count

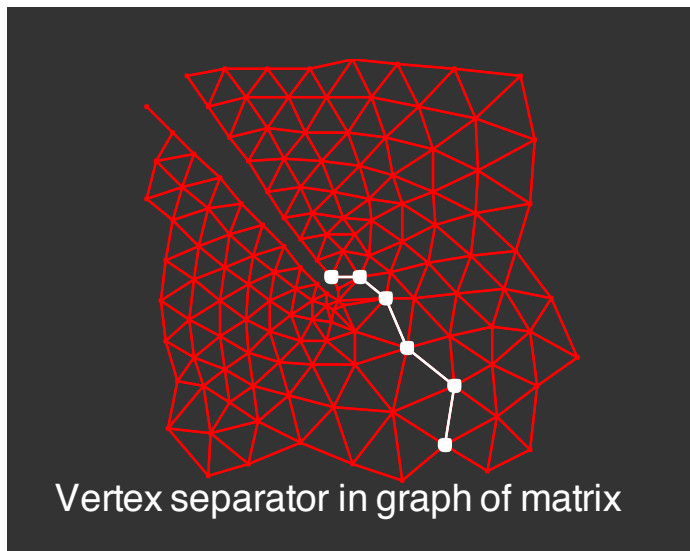
Graph algorithms in sparse matrix computation

Many, many graph algorithms have been used, invented, and implemented at large scale for sparse matrix computation:

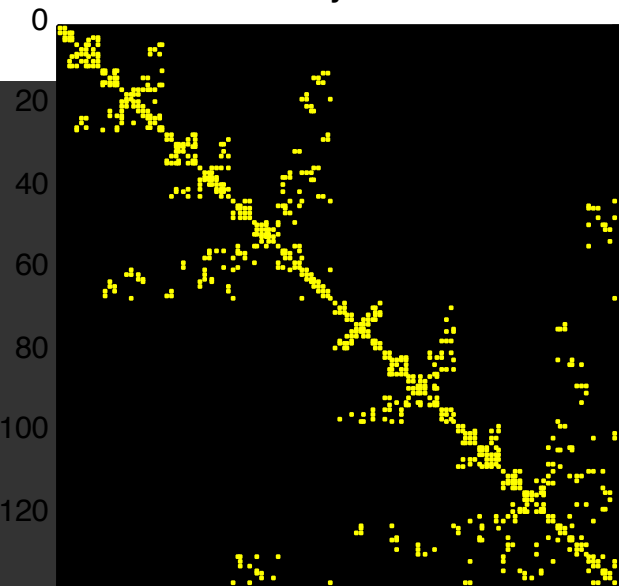
- Symmetric problems: elimination tree, nonzero structure prediction, sparse triangular solve, sparse matrix-matrix multiplication, min-height etree, ...
- Nonsymmetric problems: sparse topological solve, bipartite matching (weighted and unweighted), Dulmage-Mendelsohn decomposition / strong components, ...
- Iterative methods: graph partitioning again, independent sets, low-stretch spanning trees, ...

Nested dissection and graph partitioning

[George 1971, then many papers]



Matrix reordered by nested dissection



nz = 844

- Find a small vertex separator, number it last, recurse on subgraphs
- Approx optimal separators \Rightarrow approx optimal fill & flop count
- **It took more than 20 years for nested dissection to become the method of choice for sparse GE *in practice*.**

Combinatorial Laplacian algorithms

(partial list)

- Vaidya 1990: $O(n^{1.75})$
- Spielman/Teng 2004: $O(n \log^c n)$
- Koutis/Miller/Peng 2010: $O(n \log n \log \log n)$
- Kelner/Orecchia/Sidford/Zhu 2013: $O(n \log^2 n \log \log n)$
(for sparse graphs, fixed ϵ)

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Okay, it's been more than 20 years now ...

The Laplacian World Championships



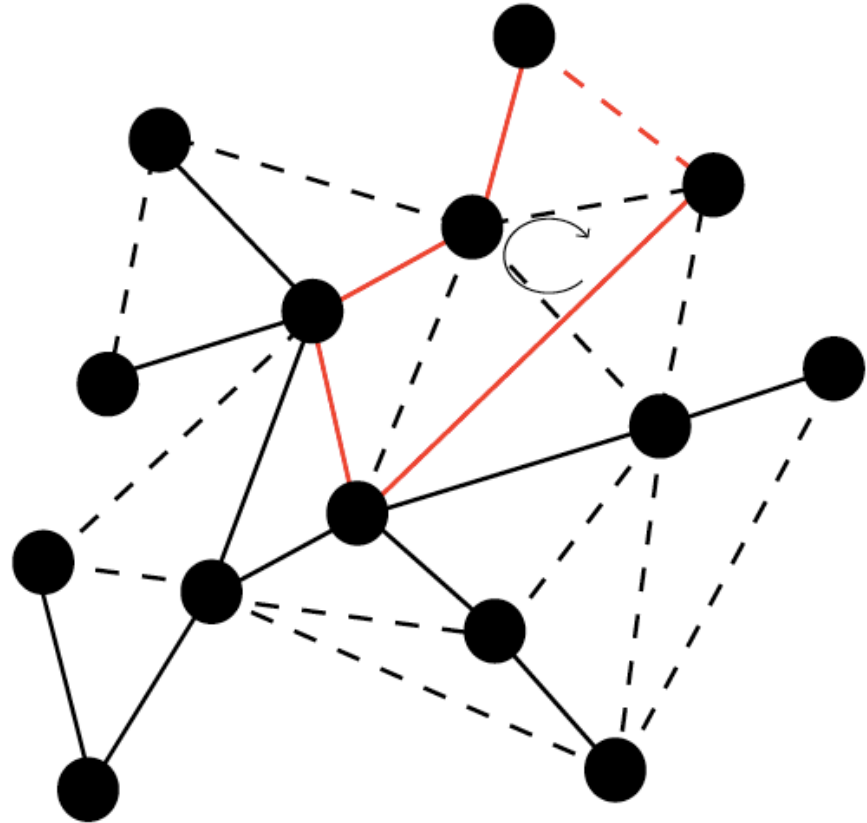
- Why should you participate?
- What would make you more likely to?
- How can we incent team leaders, members, HPC experts, ... ?
- What are the right metrics?
Single-core time; parallel time;
anything not based on timings?
- Can “you implement, we measure”
work for anything besides just
single-core time?
- Thoughts on test graphs?
Collections, generators, graphs
supplied by contestants?
- Should code from entries be
available for wider use?
- What’s the right schedule?
- What’s the right venue for the
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- Other comments? Questions?
- How can we make this succeed?

Comparing empirical complexity to theory

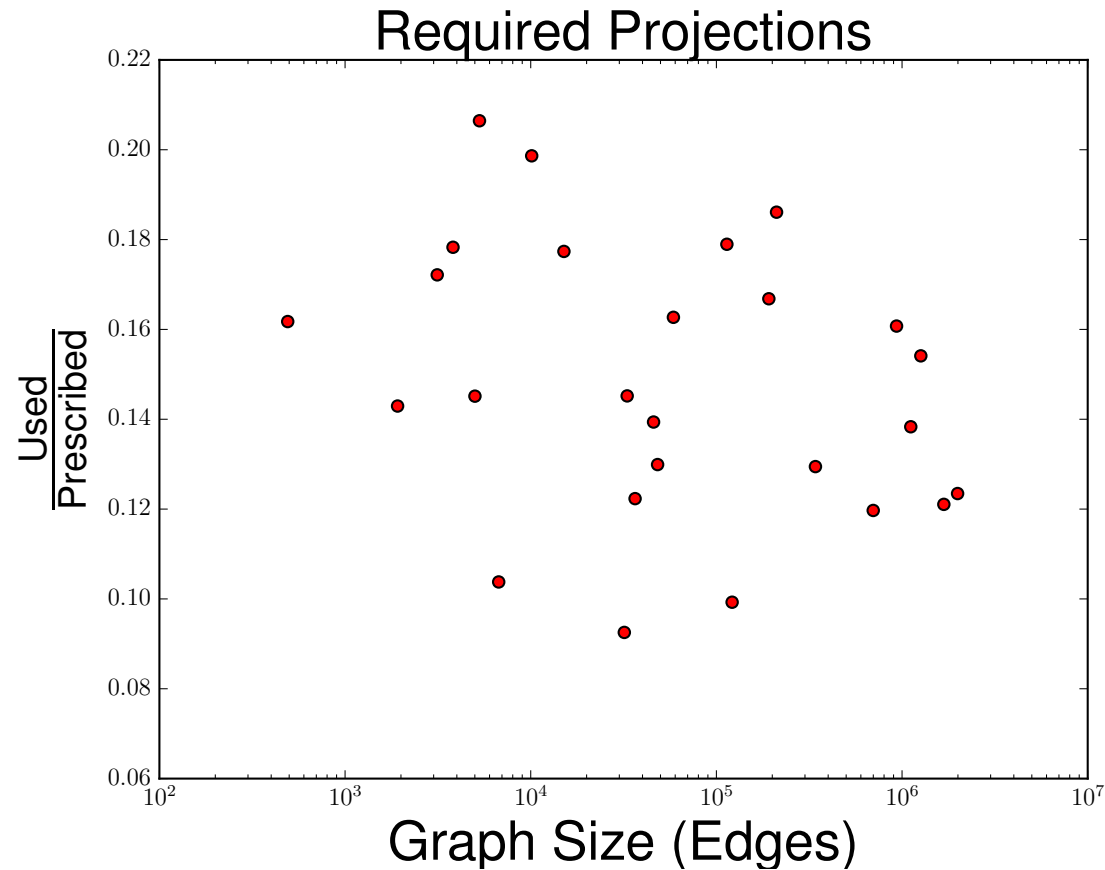
- Just counting things, not measuring time
 - a stepping stone between $O(\)$ and running time
- Here: K/O/S/Z SimpleSolver (henceforth “RK”)
 - $\log n$ off of their FullSolver
- A whole bunch of sample matrices
 - UF collection, DIMACS, BTER, etc.

Kelner/Orecchia/Sidford/Zhu algorithm

- Select cycle (with probability proportional to stretch) from a fundamental cycle basis.
- Update flows around cycle.

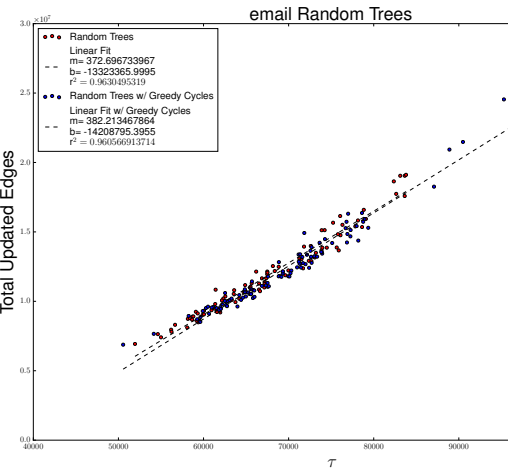


Number of projections to convergence

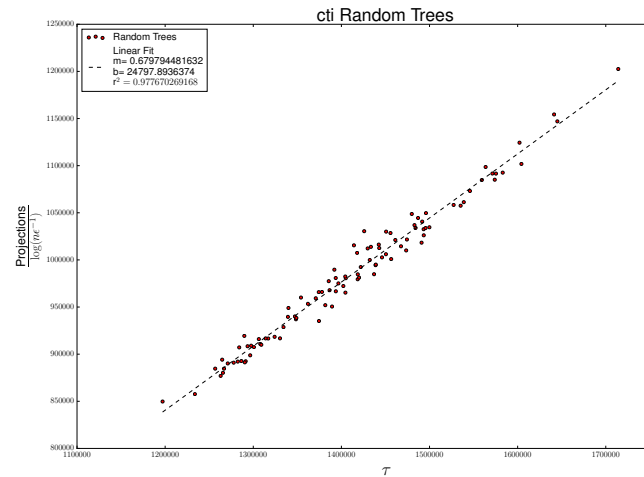


- More optimistic than the bound by a factor of 5 or 10.

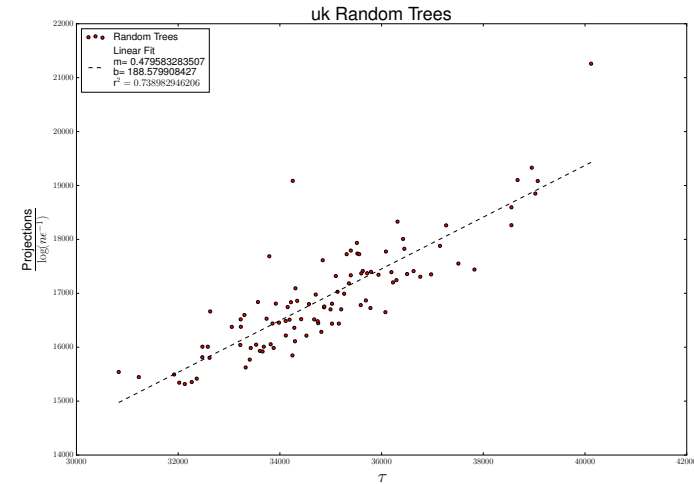
Scaling of work with tree stretch



email interactions



physical model



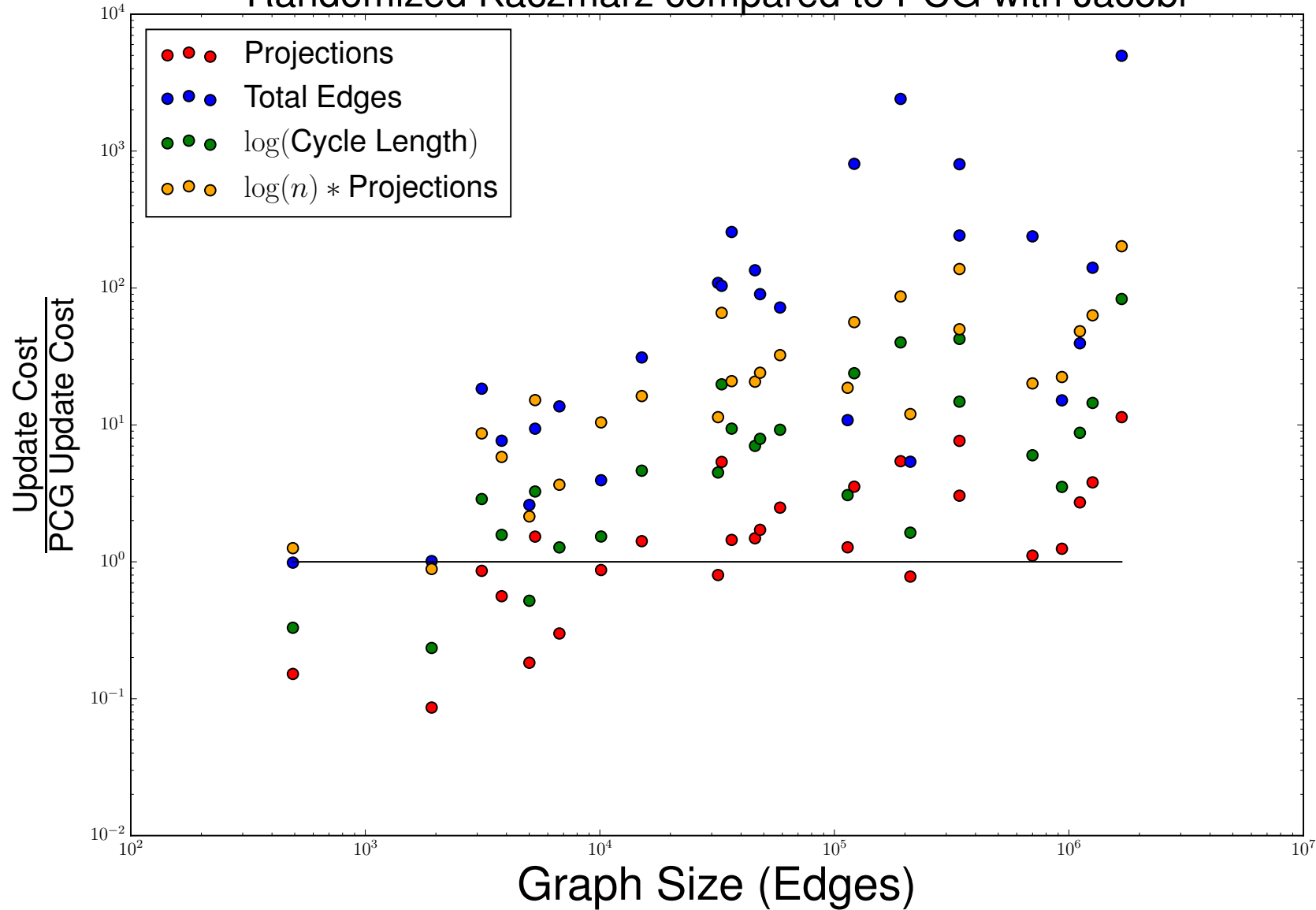
road network

- Good correlation, better for some types of graphs than others.

Empirical complexity of RK compared to PCG

- PCG = conjugate gradient preconditioned by Jacobi
- PCG does about $\text{iters} * (m+n)$ “edge touches”
- Count the cost of an RK projection in four different ways:
 - cycle length (naive)
 - $\lg n$ (fast fundamental cycle data structure)
 - $\lg (\text{cycle length})$ (unwarranted optimism)
 - 1 (surely a lower bound)

Randomized Kaczmarz compared to PCG with Jacobi



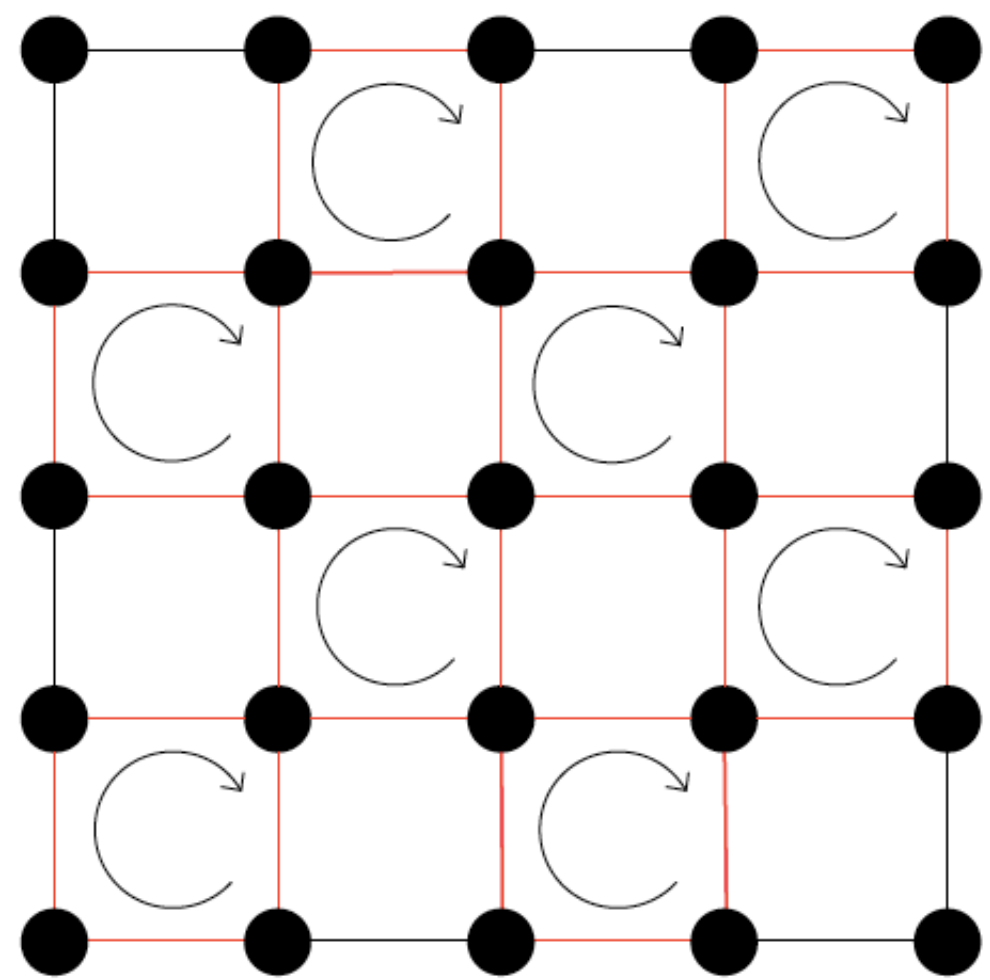
How about expanding the available projections?

- Add some cycles to basis
- Potentially gain parallelism in updates
- Maybe add flexibility with non-fundamental cycles

- More of the projection cost metrics are fantasy now, since we don't have fast update algorithms in general.

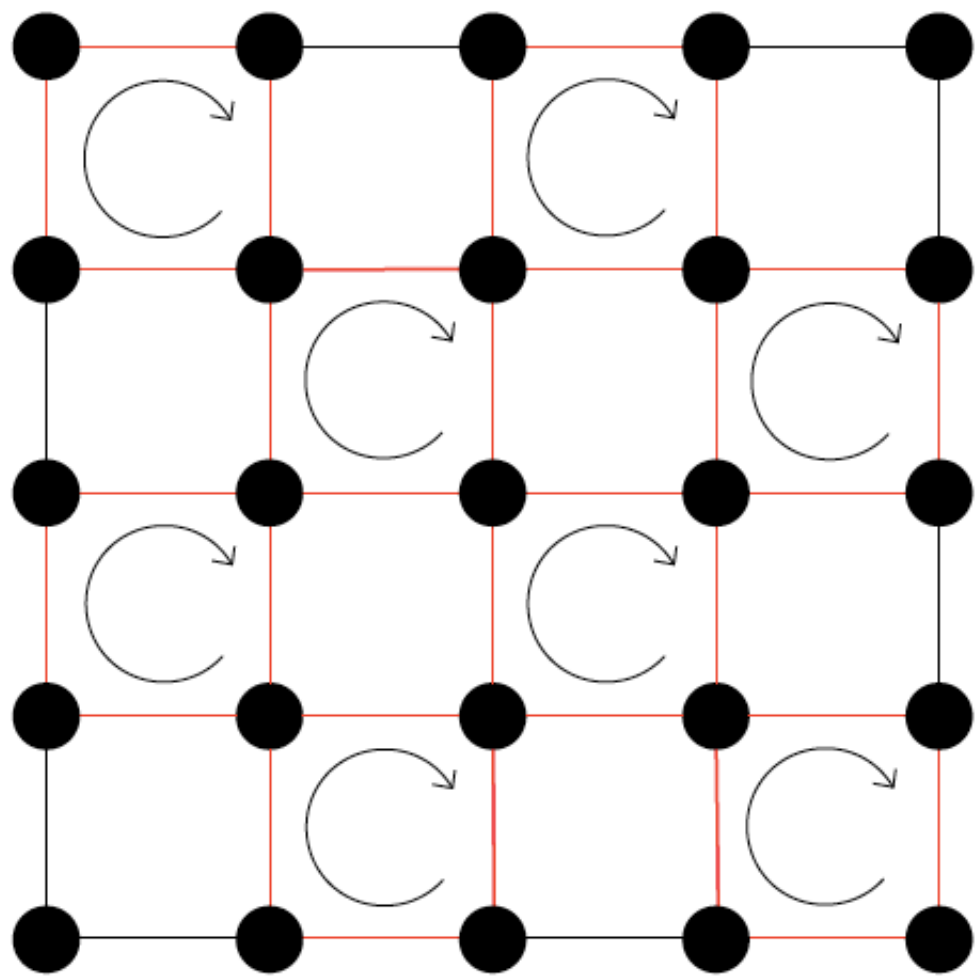
Example: non-fundamental cycles on a square grid

- Edge-disjoint cycles can be updated in parallel.

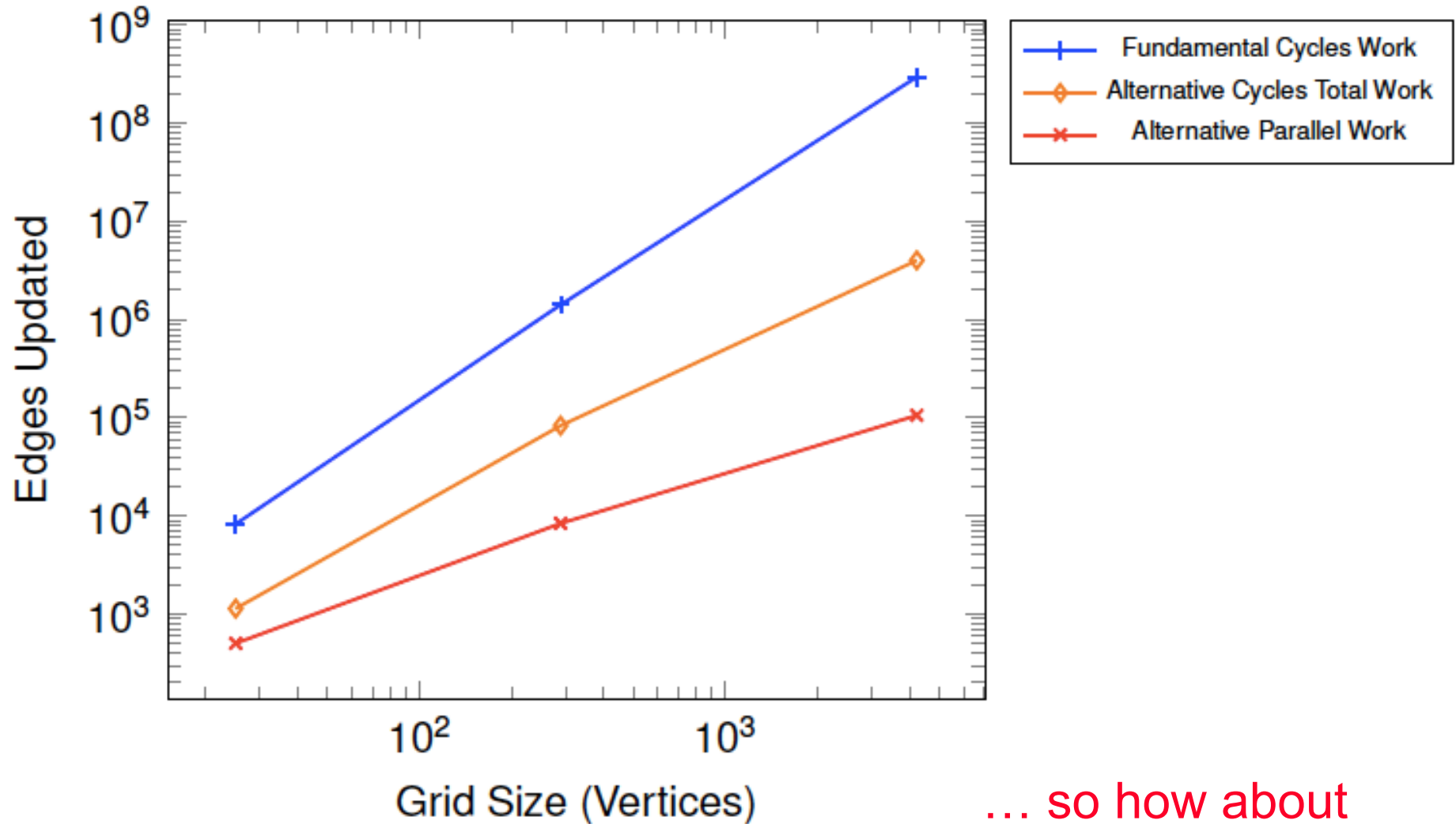


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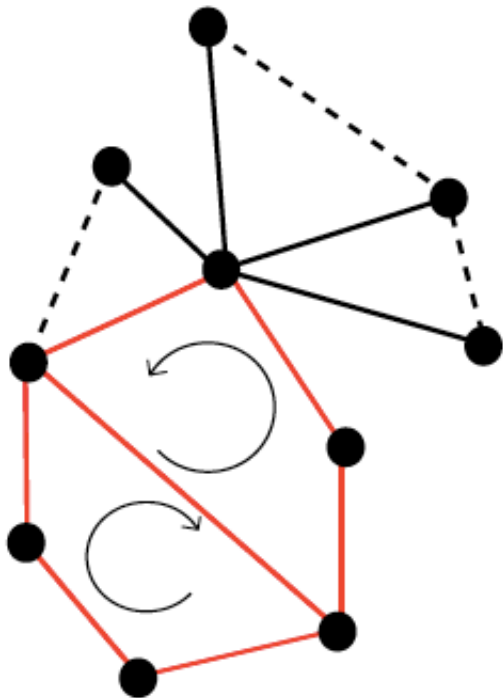


Work and span for square grid example



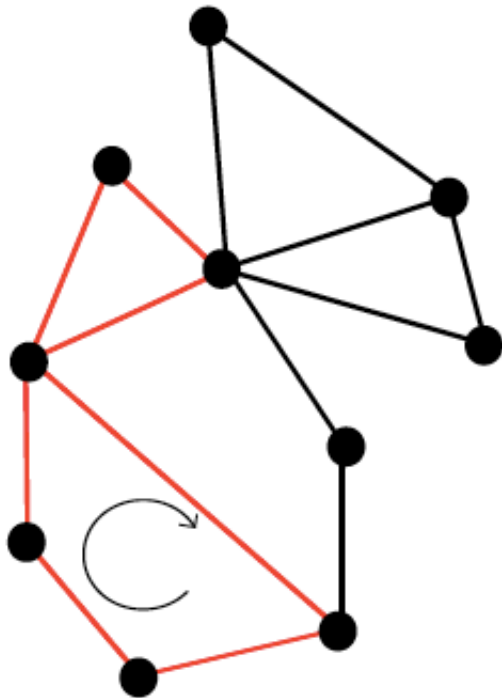
... so how about
general graphs?

Tree shortcut cycles



- Select an off-tree edge.
- Search for a shortcut between endpoints of this edge.
 - Only search edges closer to root of tree.
 - Truncate search to control cost.
- Replace tree cycle with shortcut cycle.
 - Cycle space dimension remains the same.

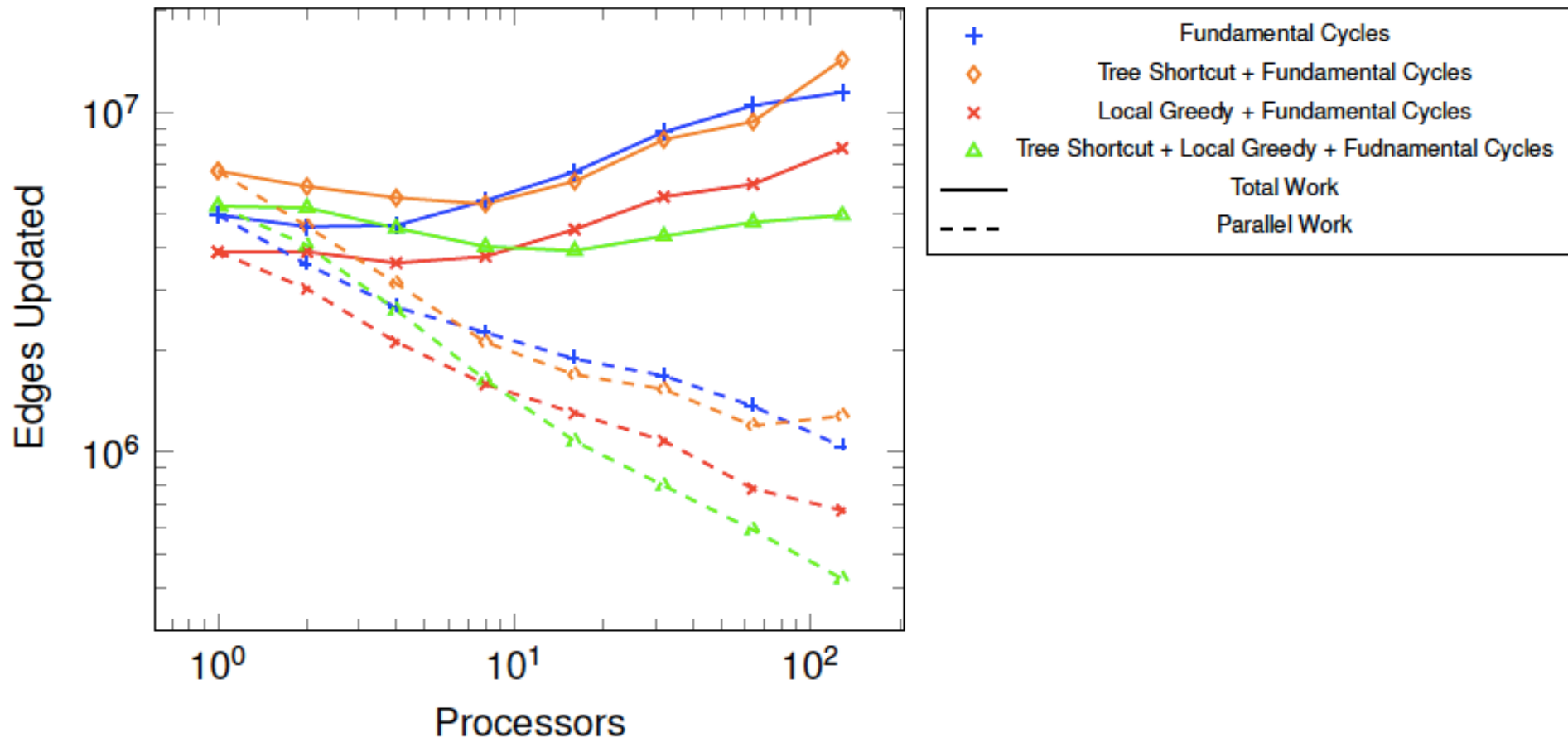
Local greedy cycles



- Select an unmarked edge.
- Find smallest cycle containing this edge.
 - Truncate search to control cost.
 - If found mark all the used edges.
- This set is not guaranteed to span the cycle space.

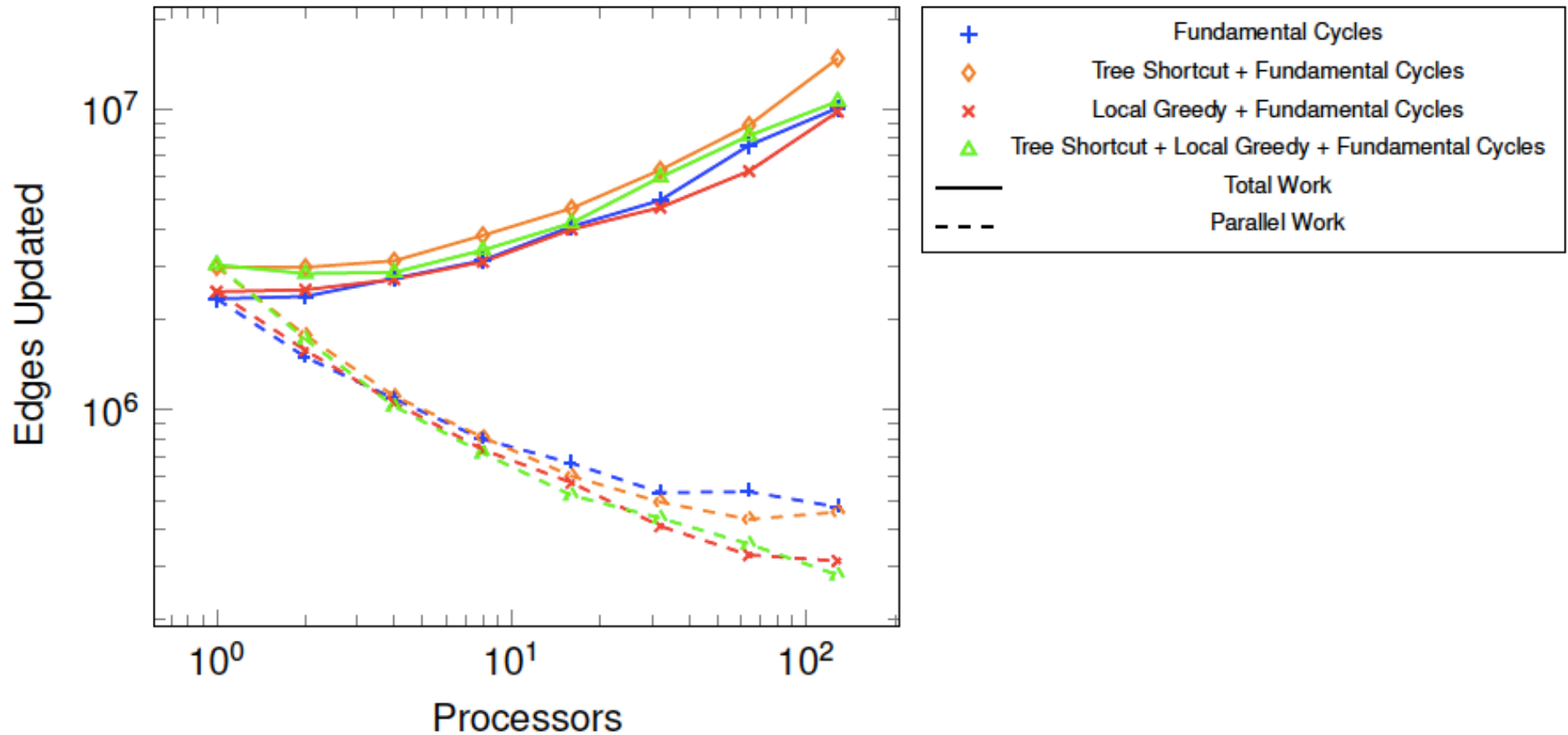
Comparing cycle sets: work and parallel work

USpowerGrid



Comparing cycle sets: work and parallel work

email



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