Filtering: A Method for Solving Graph Problems in MapReduce

Benjamin Moseley UIUC

Silvio Lattanzi Google

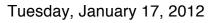
Google

Siddharth Suri Yahoo!

Sergei Vassilvitski Yahoo!



Y



Overview

- Introduction to MapReduce model
- Our settings
- Our results
- Open questions

Introduction to the MapReduce model

XXL Data

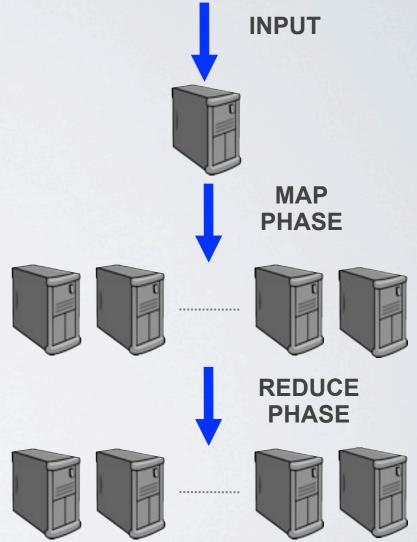
- Huge amount of data
- Main problem is to analyze information quickly

XXL Data

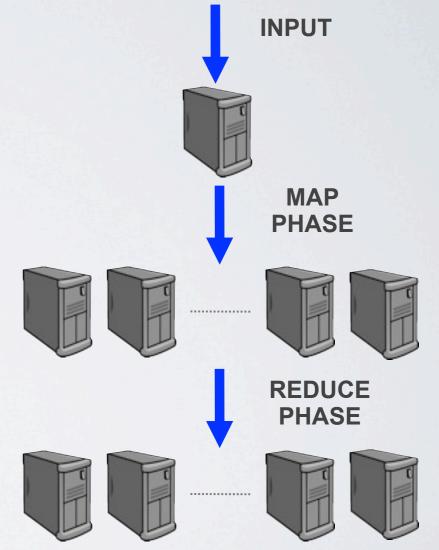
- Huge amount of data
- Main problem is to analyze information quickly
- New tools
- Suitable efficient algorithms



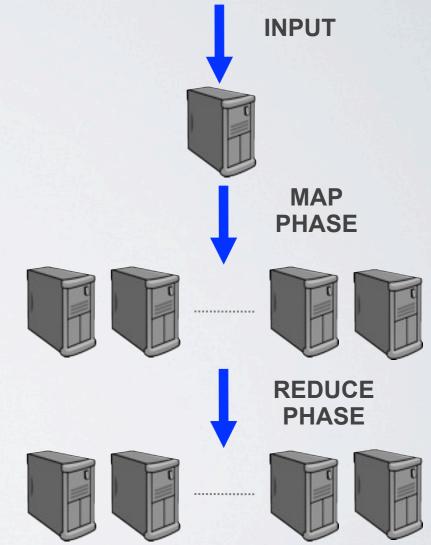
 MapReduce is the platform of choice for processing massive data



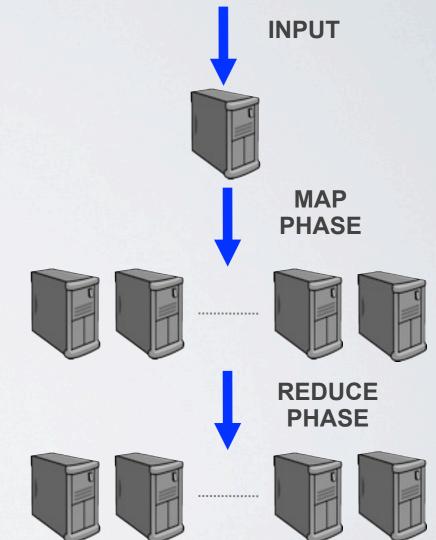
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 < key, value >



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- Data are represented as tuple
 < key, value >
- Mapper decides how data is distributed
- Reducer performs non-trivial computation locally



How can we model MapReduce?

PRAM

- No limit on the number of processors
- Memory is uniformly accessible from any processor
- No limit on the memory available

How can we model MapReduce?

STREAMING

- Just one processor
- There is a limited amount of memory
- No parallelization

[Karloff, Suri and Vassilvitskii]

• N is the input size and $\epsilon > 0$ is some fixed constant

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 \mathcal{MRC}^{i} : problem that can be solved in $O(\log^{i} N)$ rounds

Combining map and reduce phase

Mapper and Reducer work only on a subgraph

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Time is constrained by the number of rounds

Algorithmic challenges

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Algorithmic challenges

No machine can see the entire input

- No communication between machines during each phase
- Total memory is $N^{2-2\epsilon}$

MapReduce vs MUD algorithms

In MUD framework each reducer operates on a stream of data.

In MUD, each reducer is restricted to only using polylogarithmic space.

We study the Karloff, Suri and Vassilvitskii model

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• We focus on class \mathcal{MRC}^0

• We assume to work with dense graph $m = n^{1+c}$, for some constant c > 0

Our settings

• We assume to work with dense graph $m = n^{1+c}$, for some constant c > 0

 Empirical evidences that social networks are dense graphs [Leskovec, Kleinberg and Faloutsos]

Dense graph motivation

[Leskovec, Kleinberg and Faloutsos]

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•Lowest value of c founded .08 and four graphs have c > .5

Our results

Results

Constant rounds algorithms for

- Maximal matching
- Minimum cut
- 8-approx for maximum weighted matching
- 2-approx for vertex cover
- 3/2-approx for edge cover

Notation

- G = (V, E) : Input graph
- n: number of nodes
- *m* : number of edges
- η : memory available on each machine
- N: input size

Filtering

 Part of the input is dropped or filtered on the first stage in parallel

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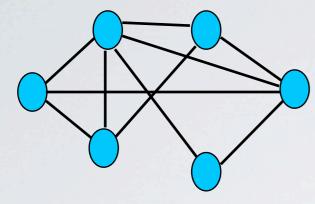
Next some computation is performed on the filtered input

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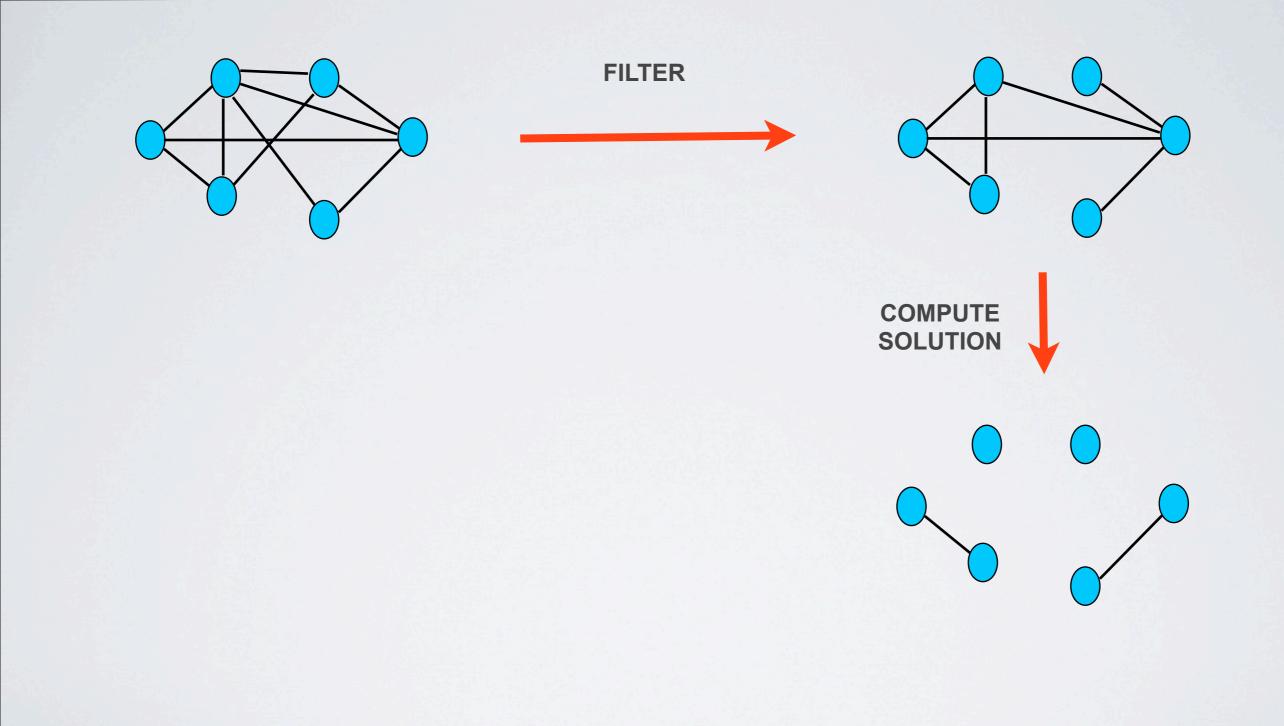
Finally some patchwork is done to ensure a proper solution



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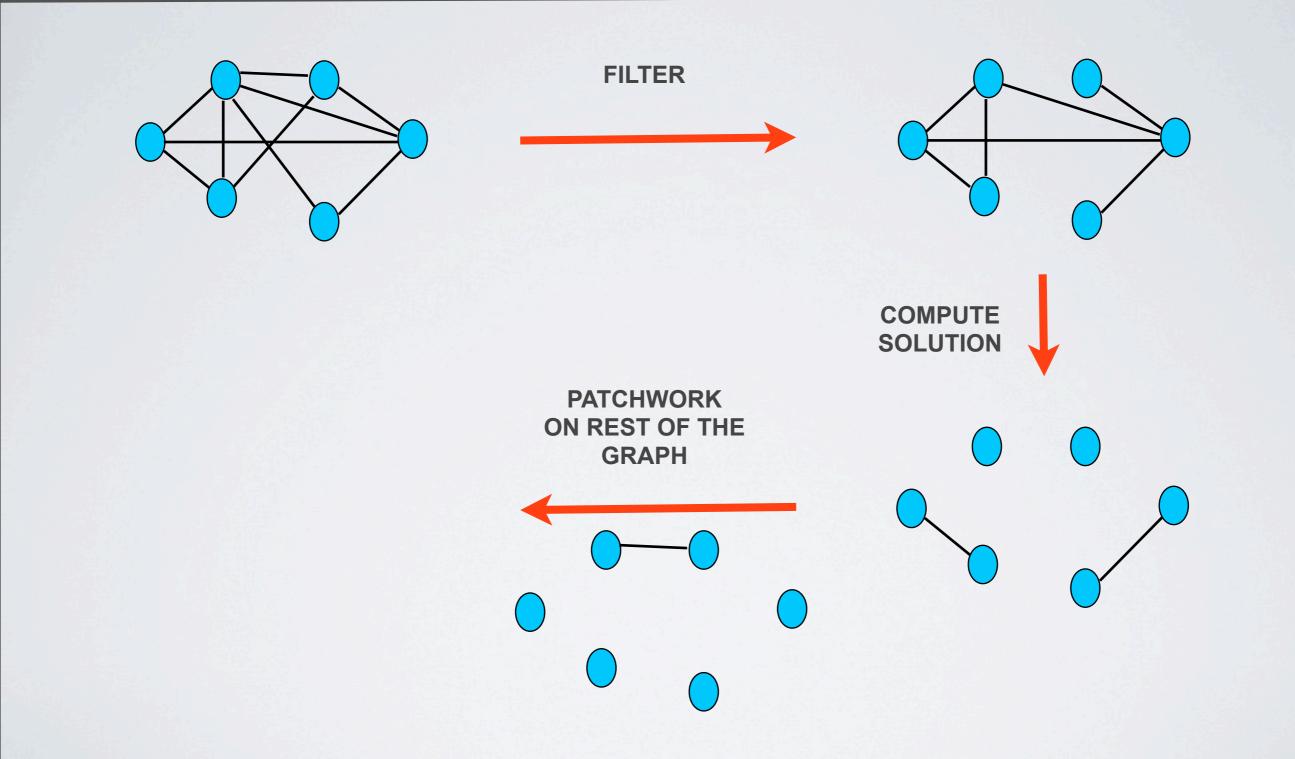
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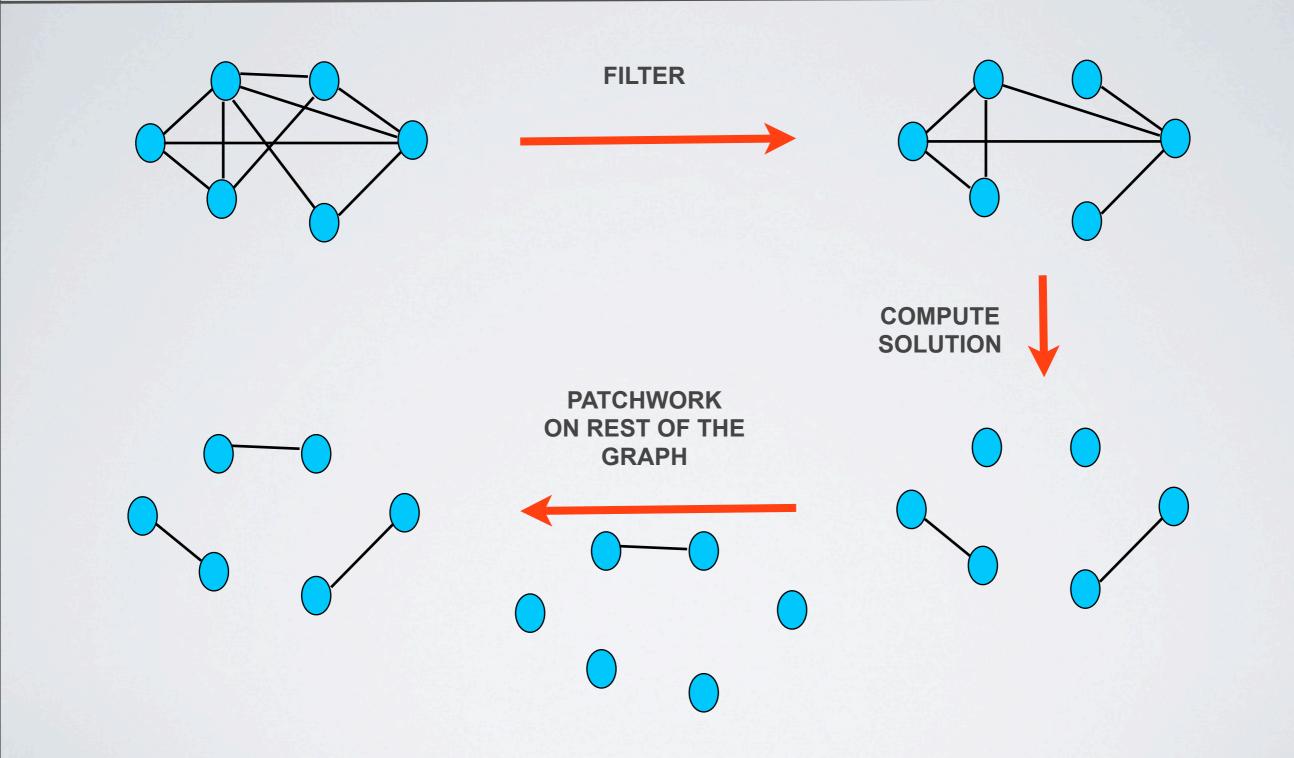
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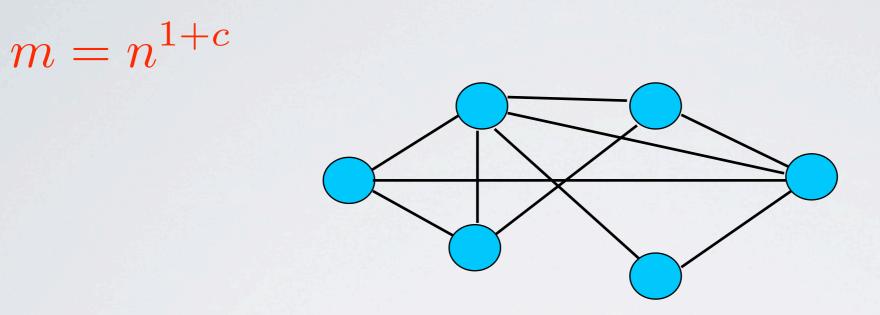
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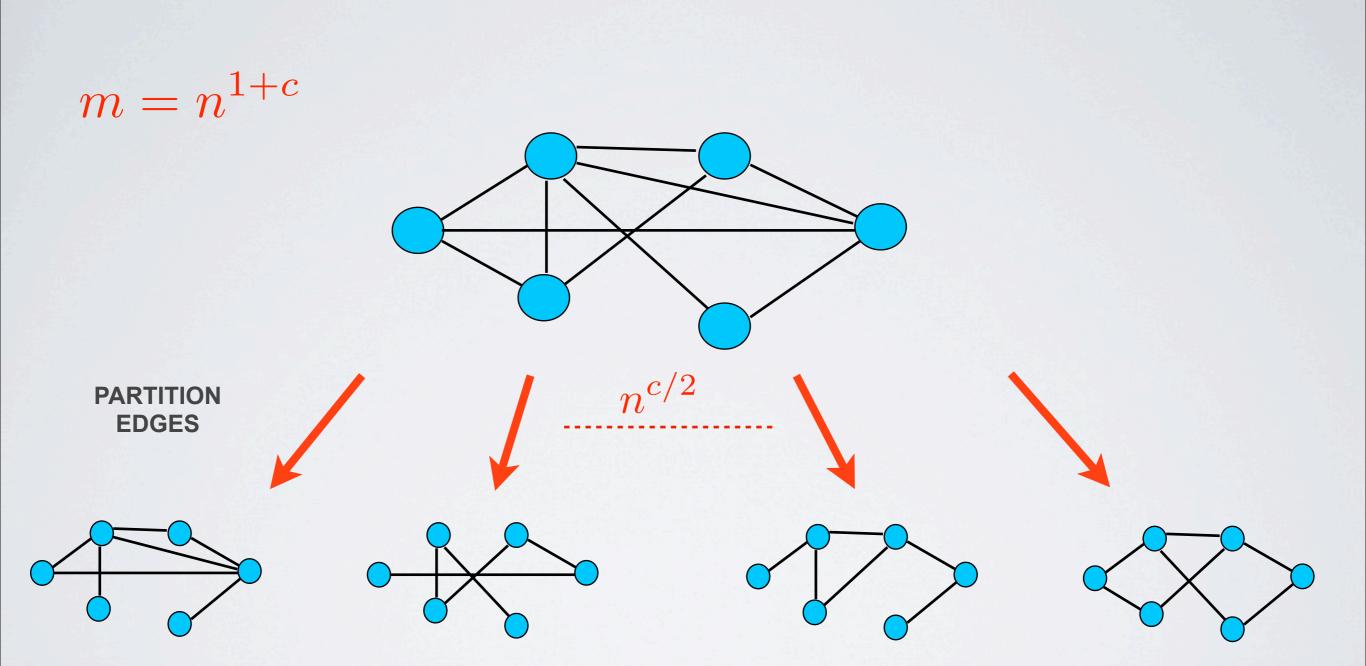
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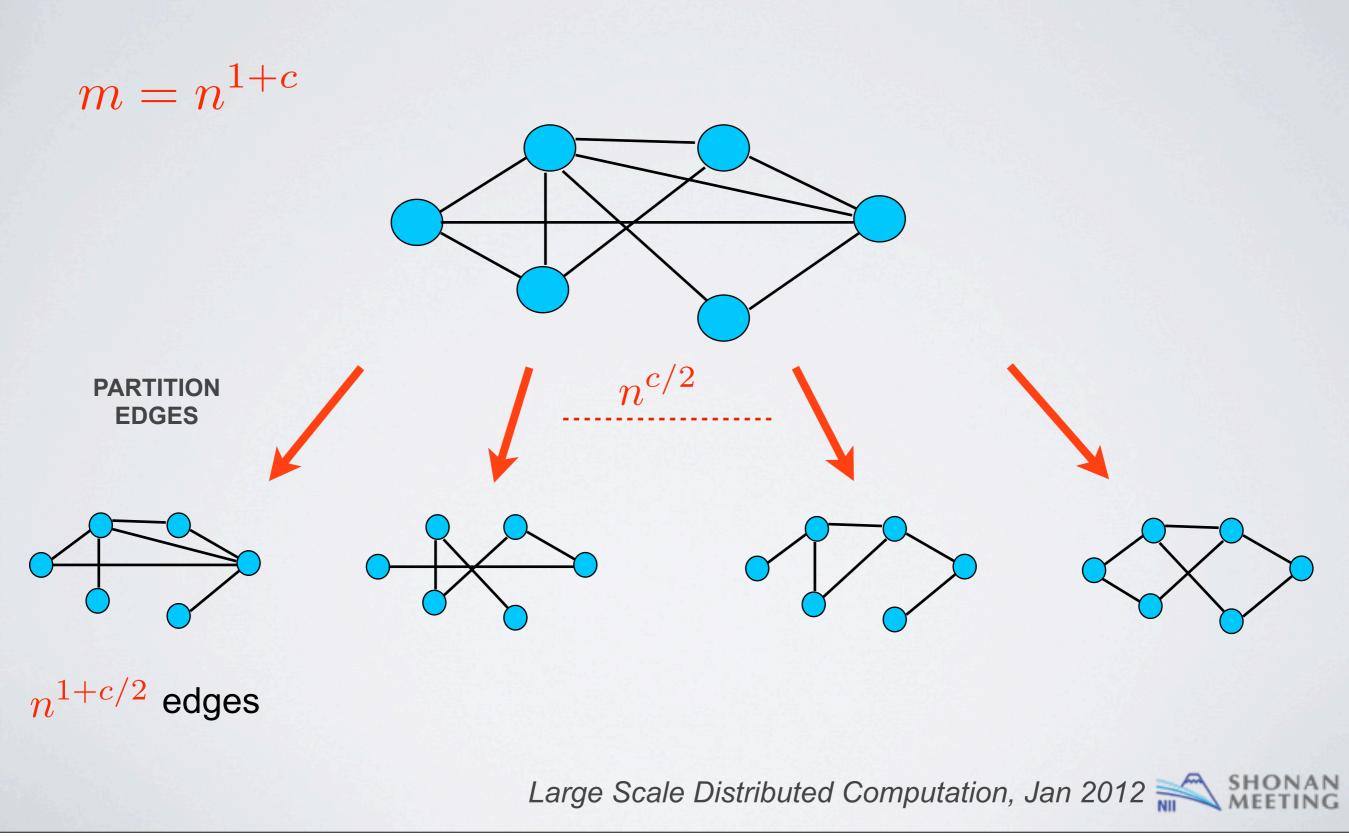


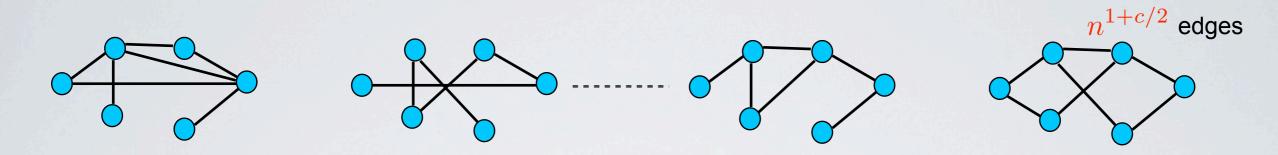
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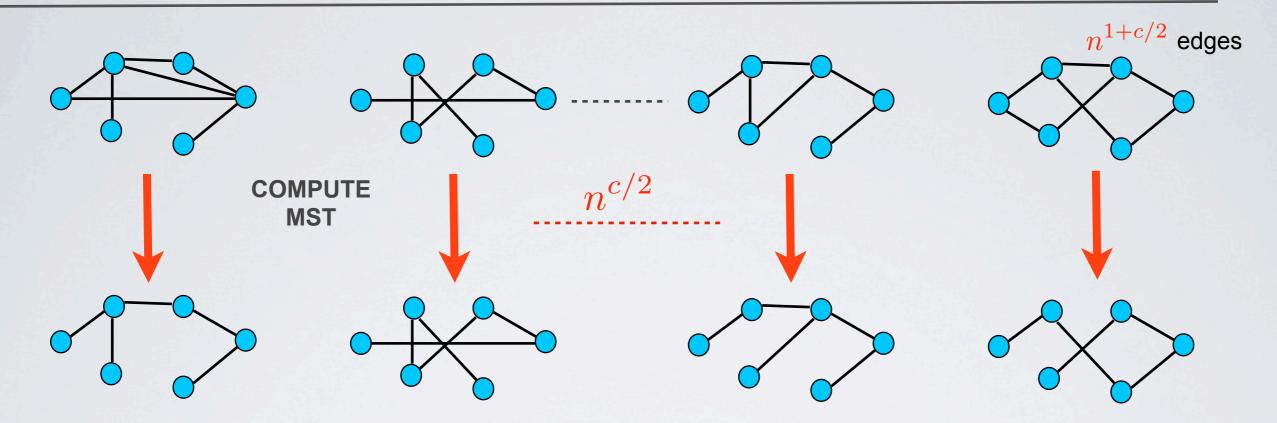
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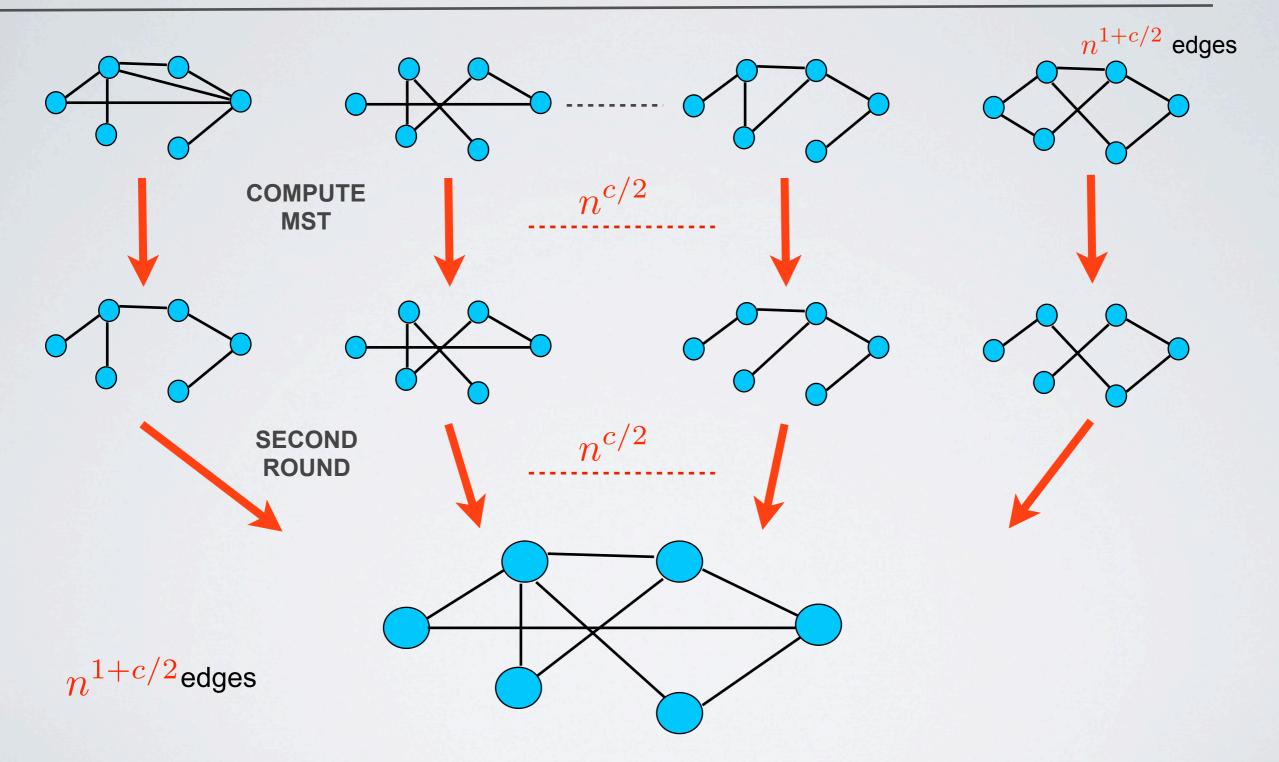












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- No edge in the final solution is discarded in partial solution
- The algorithm runs in two rounds
- No more than $O\left(n^{\frac{c}{2}}\right)$ machines are used
- No machine uses memory greater than $O(n^{1+c/2})$

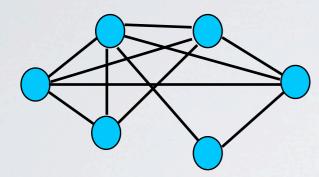
Maximal matching

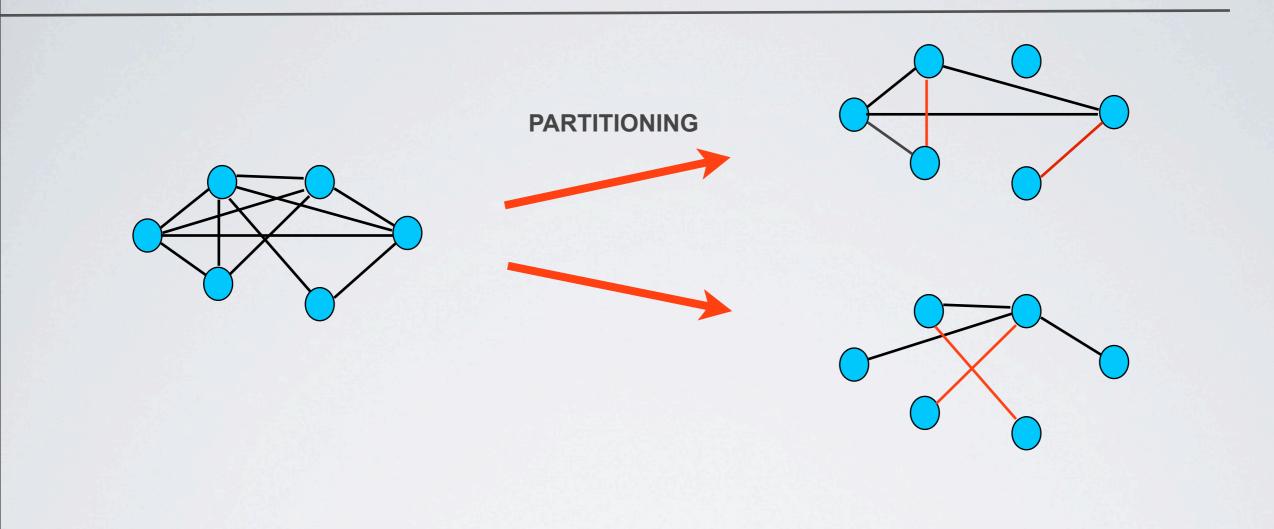
 Algorithmic difficulty is that each machine can only see edges assigned to the machine

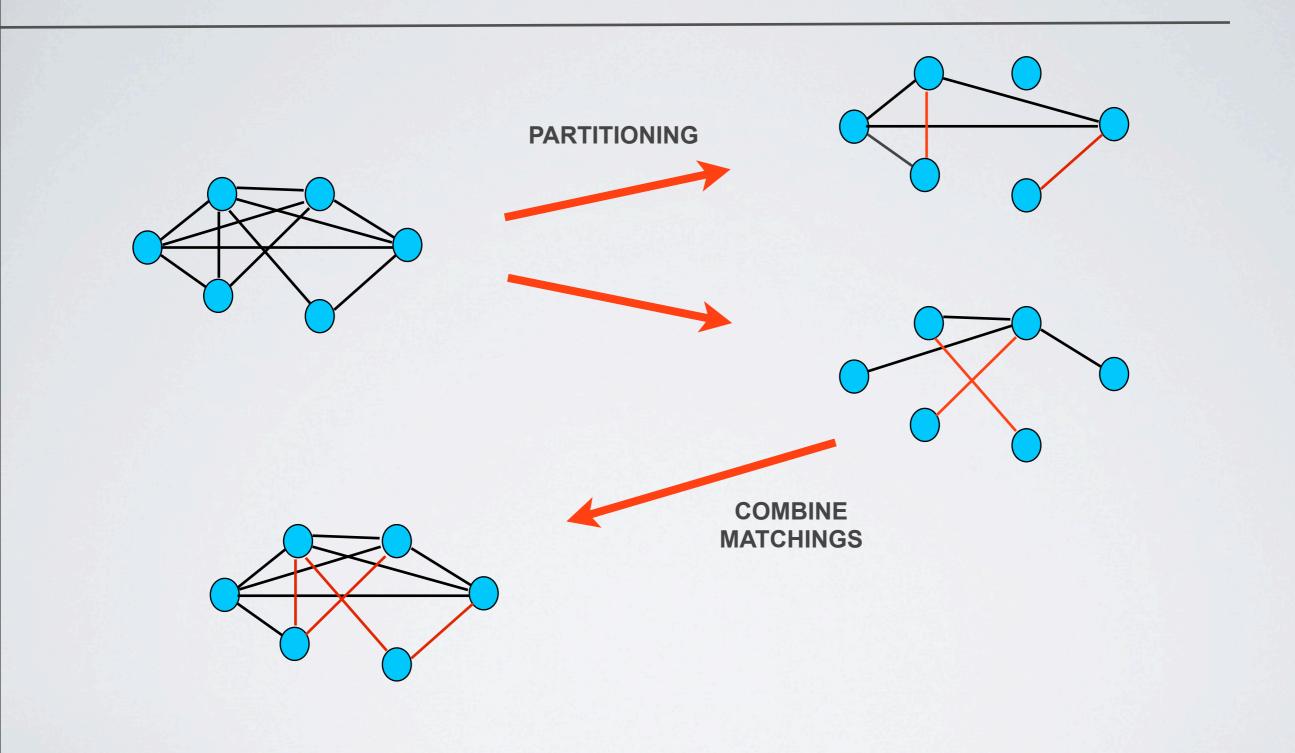
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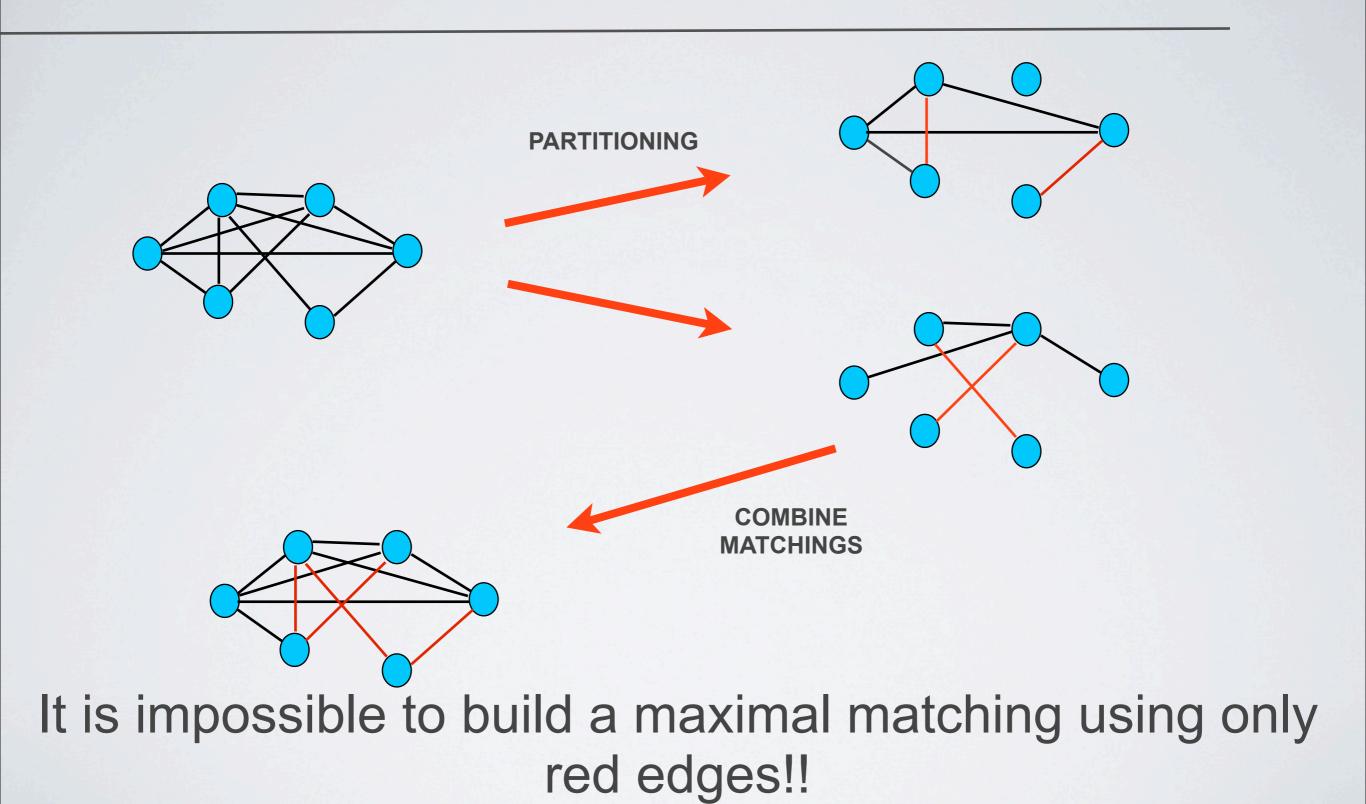
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Is a partitioning based algorithm feasible?

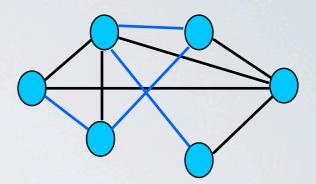






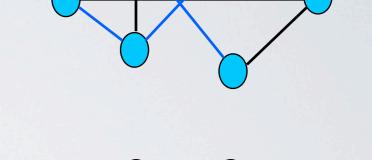


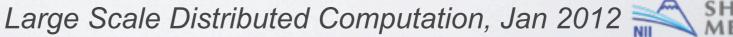
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•Let M' be a maximal matching on G[E']

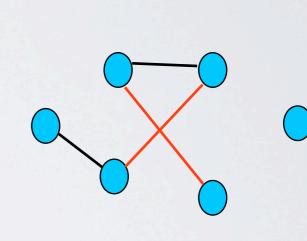


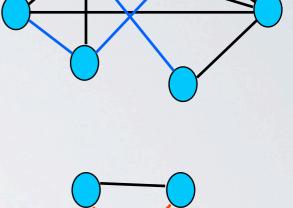


• Consider any subset of the edges E'

-Let M' be a maximal matching on G[E']

• The unmatched vertices form a independent set





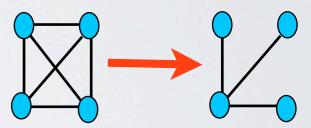
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Find a matching on a sample and then find a matching on unmatched vertices

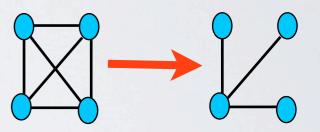
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- For dense portions of the graph, many edges should be sampled



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Sparse portions of the graph are small
 and can be placed on a single machine

• Sample each edge independently with probability $p = \frac{10 \log n}{n^{c/2}}$

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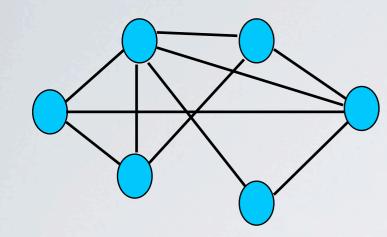
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Consider the induced subgraph on unmatched vertices

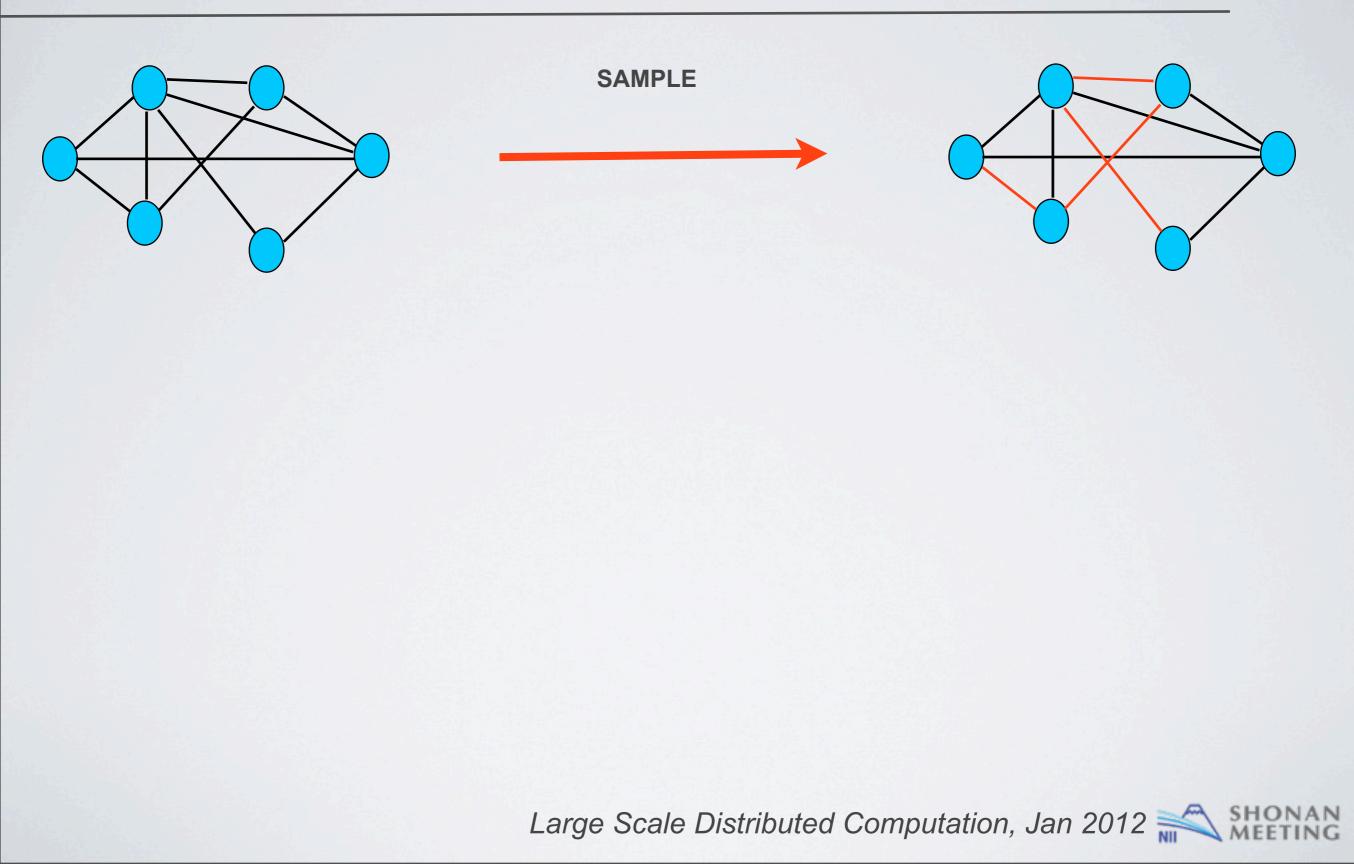
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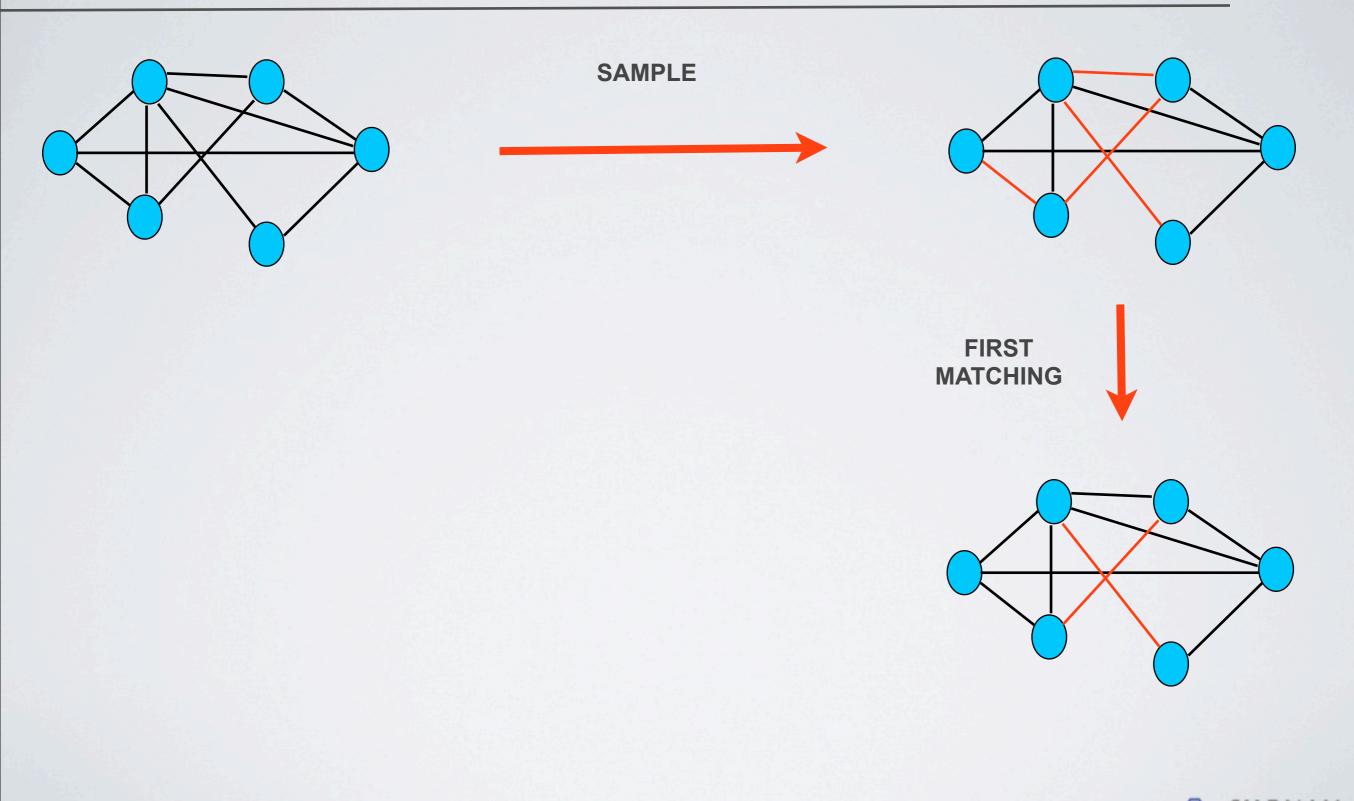
- Find a matching on the sample
- Consider the induced subgraph on unmatched vertices
- Find a matching on this graph and output the union



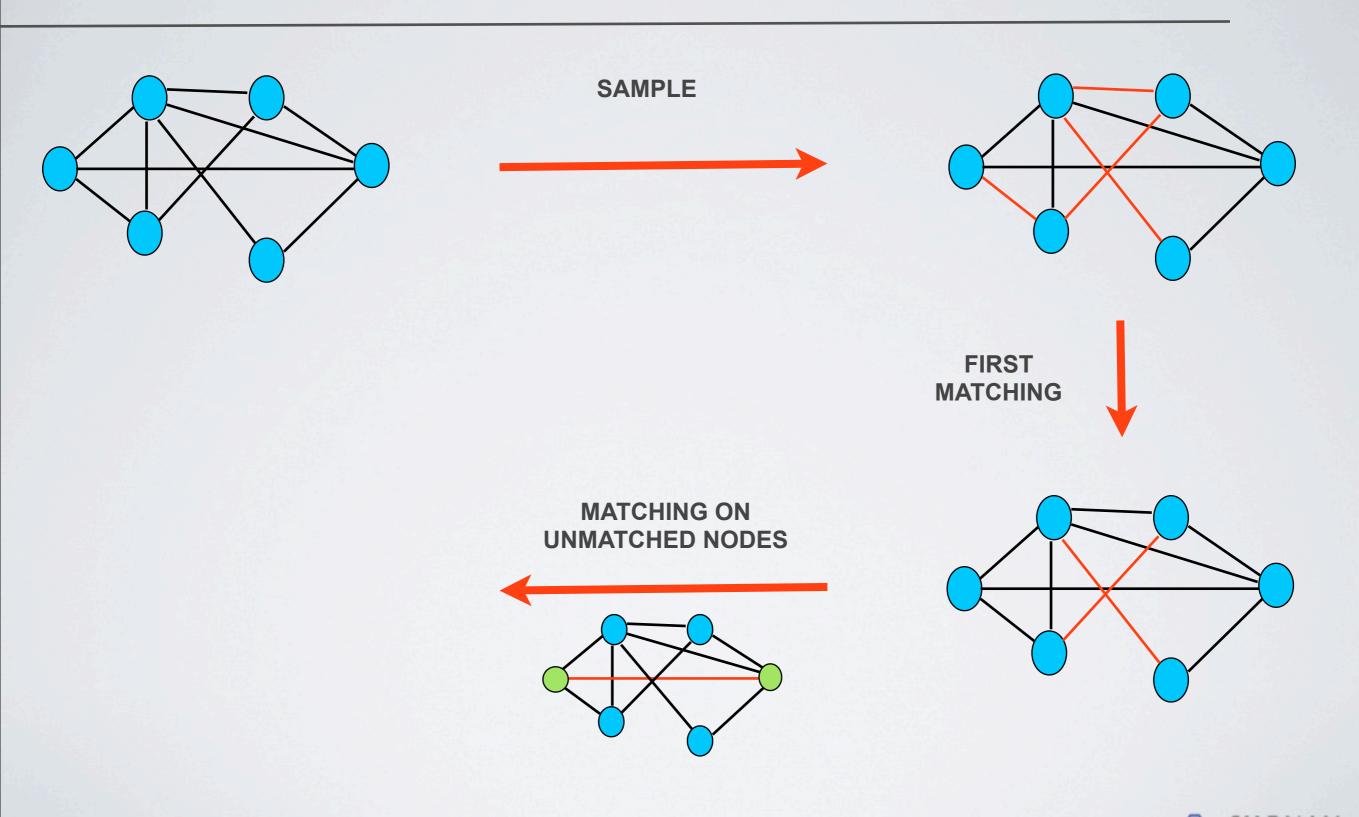
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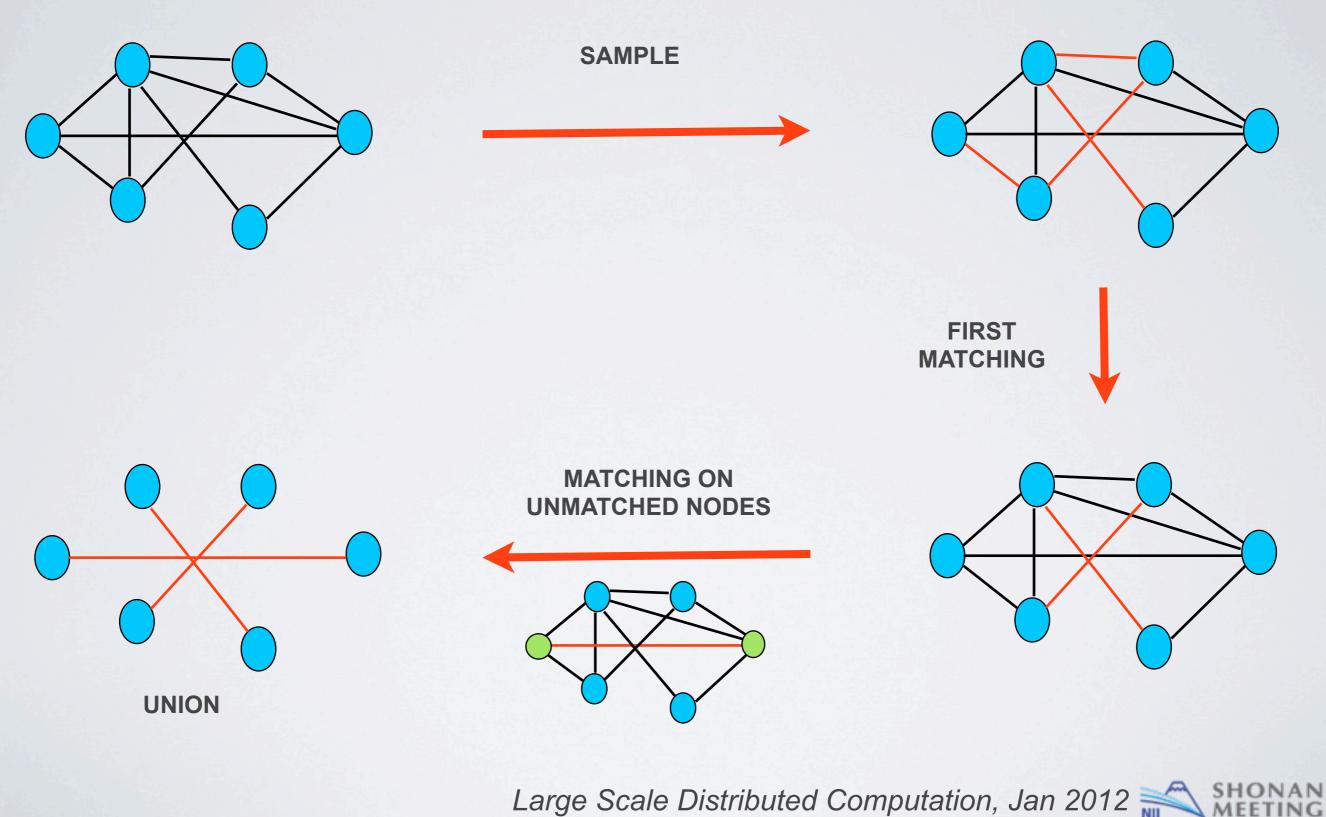




Algorithm



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Correctness

Consider the last step of the algorithm

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All unmatched vertices are placed on a machine

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Consider the last step of the algorithm

All unmatched vertices are placed on a machine

 All unmatched vertices or are matched at the last step or have only matched neighbors

Bounding the rounds

Three rounds:

Sampling the edges

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 Find a matching on a single machine for the sampled edges

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Three rounds:

- Sampling the edges
- Find a matching on a single machine for the sampled edges
- Find a matching for the unmatched vertices

- Each edge was sampled with probability $p = rac{10\log n}{n^{c/2}}$

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- Using Chernoff the sampled graph has size $\tilde{O}(n^{1+c/2})$ with high probability
- Can we bound the size of the induced subgraph on the unmatched vertices?

• For a fixed induced subgraph with at least $n^{1+c/2}$ edges the probability an edge is not sampled is: $\left(1 - \frac{10\log n}{n^{c/2}}\right)^{n^{1+c/2}} \leq exp(-10n\log n)$

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• Union bounding over all 2^n induced subgraphs shows that at least one edge is sampled from every dense subgraph with probability

$$1 - exp(-10\log n) \ge 1 - \frac{1}{n^{10}}$$

• Can we run the algorithm with less then $\tilde{O}(n^{1+c/2})$ memory?

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• We can amplify our sampling technique!

Sample as many edges as possible that fits in memory

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Find a matching

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- Sample as many edges as possible that fits in memory
- Find a matching
- If the edges between unmatched vertices fit onto a single machine, find a matching on those vertices
- Otherwise, recurse on the unmatched nodes
- With $n^{1+\epsilon}$ memory each iteration a factor of n^{ϵ} edges are removed resulting in $O(c/\epsilon)$ rounds

Parallel computation power

 Maximal matching algorithm does not use parallelization

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 Maximal matching algorithm does not use parallelization

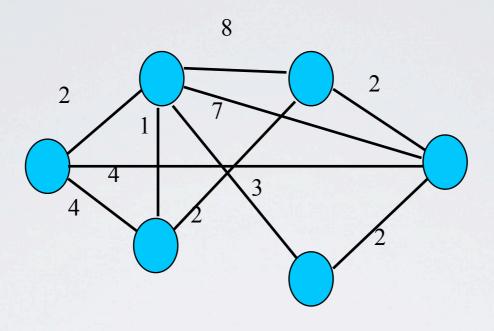
We use a single machine in every step

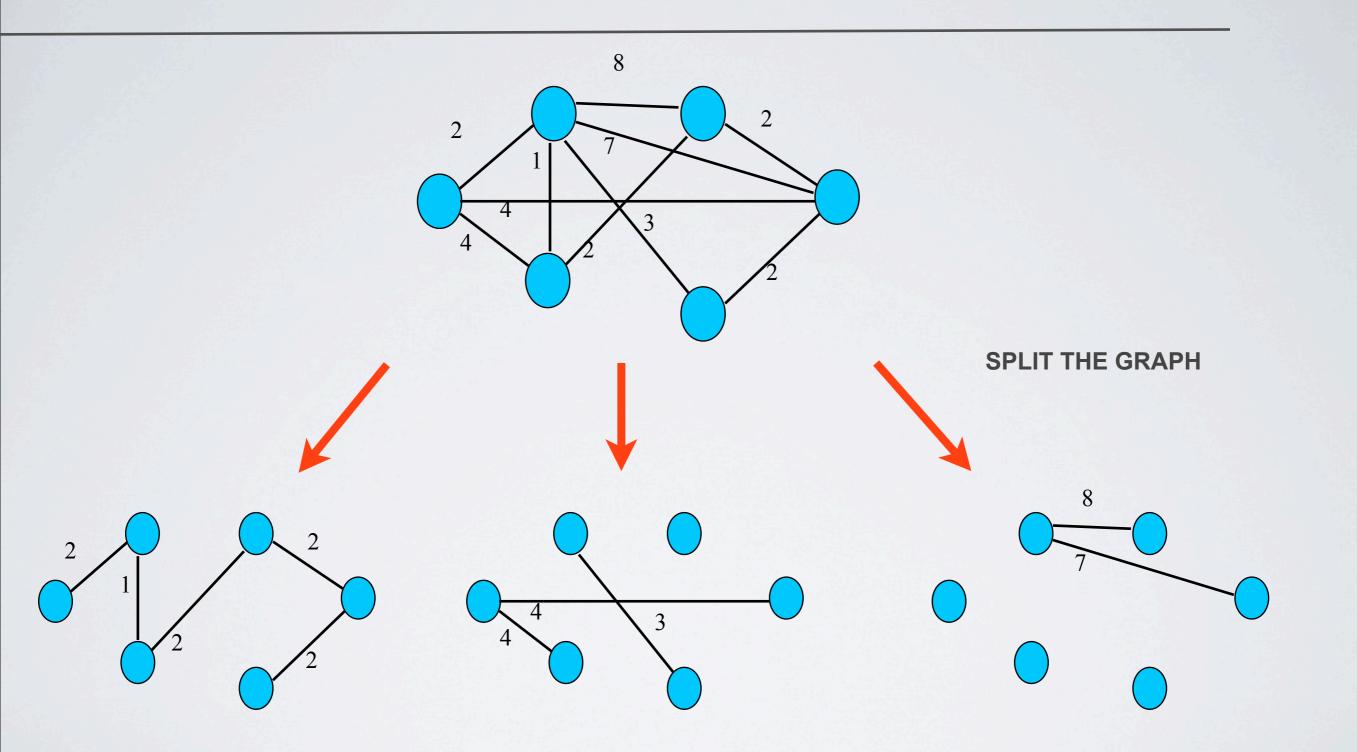
Parallel computation power

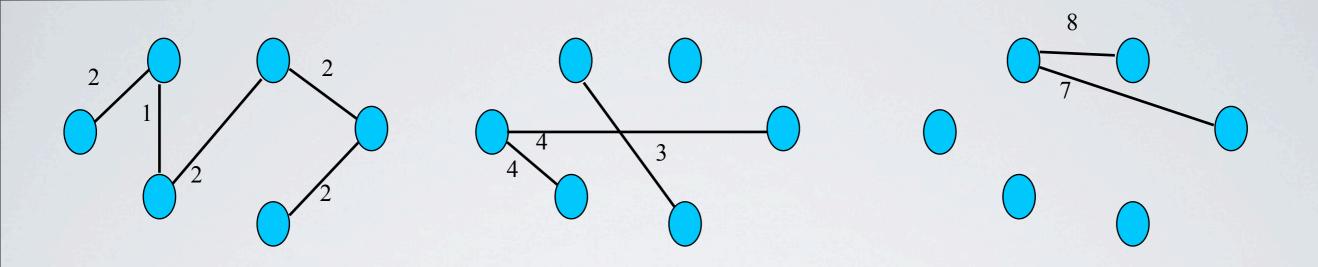
 Maximal matching algorithm does not use parallelization

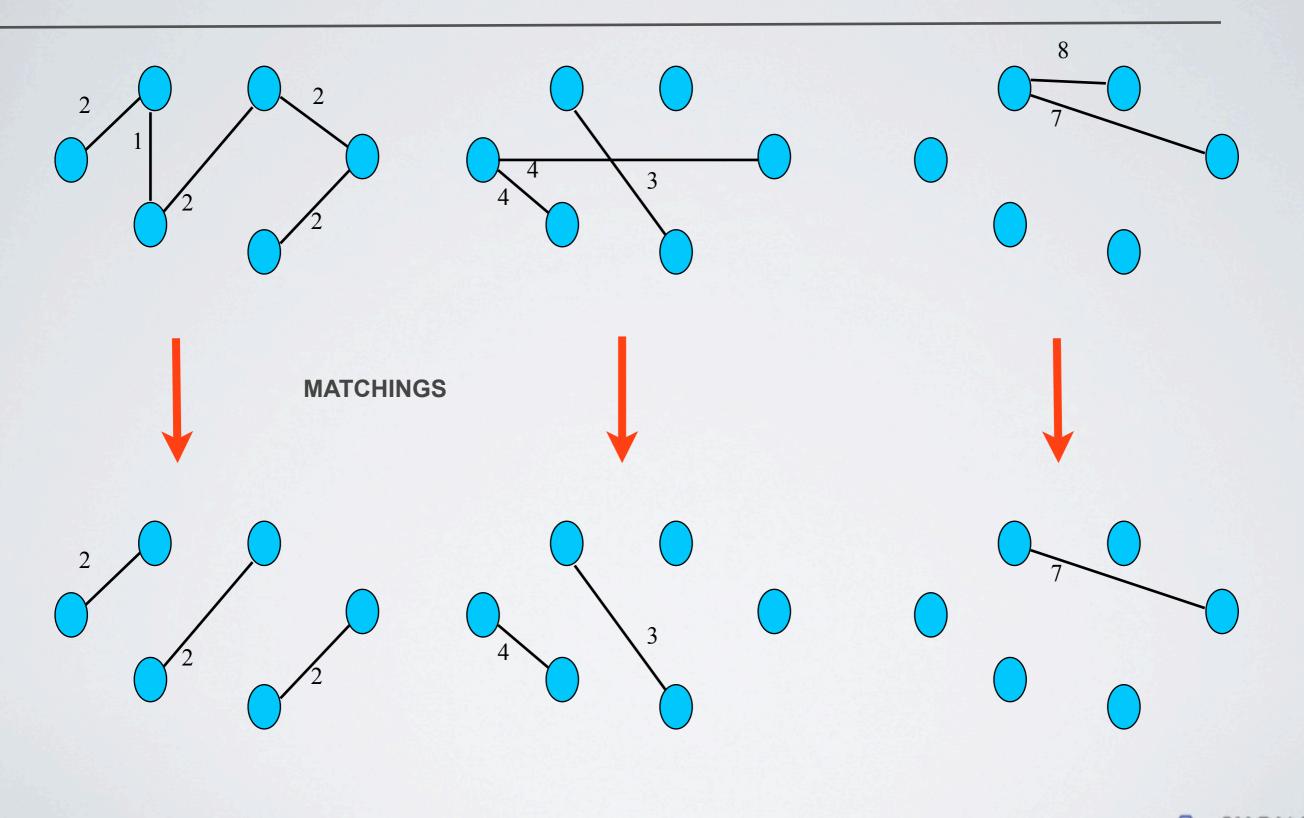
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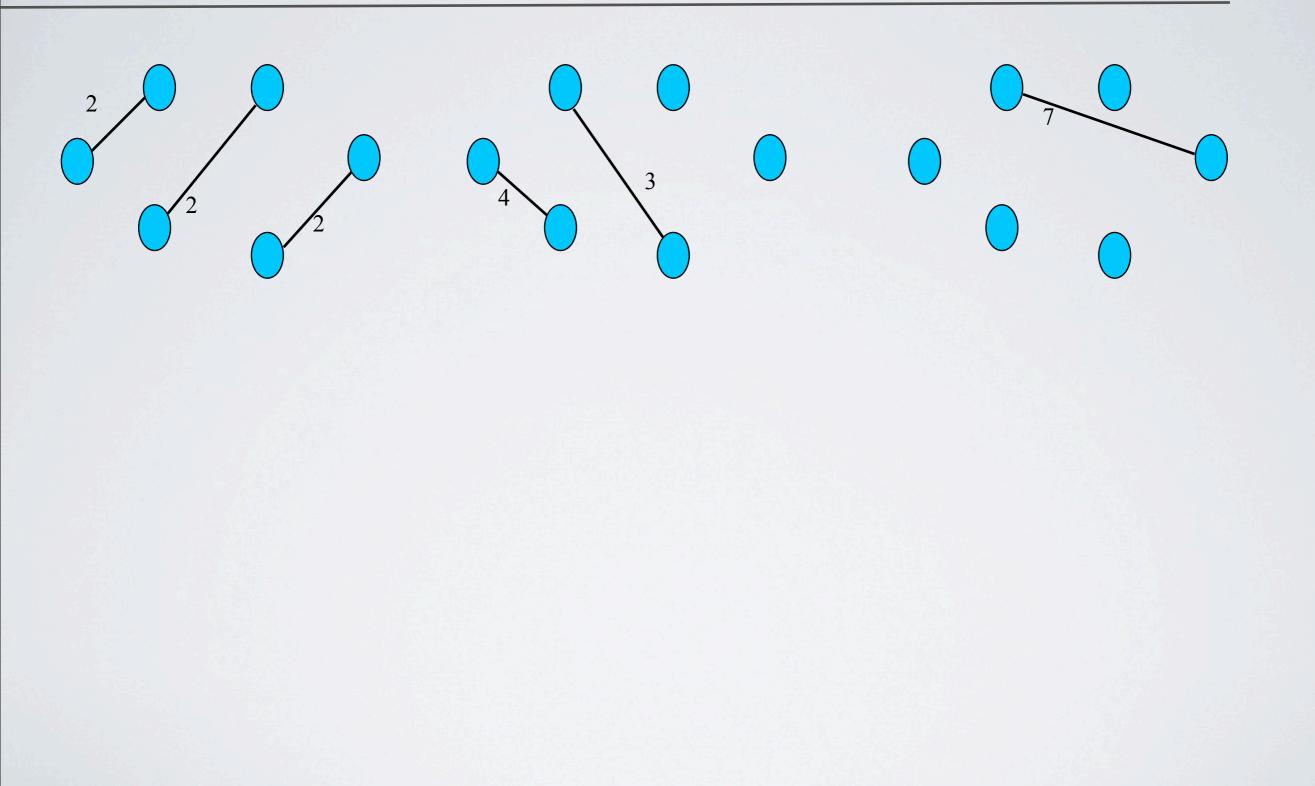
When parallelization is used?

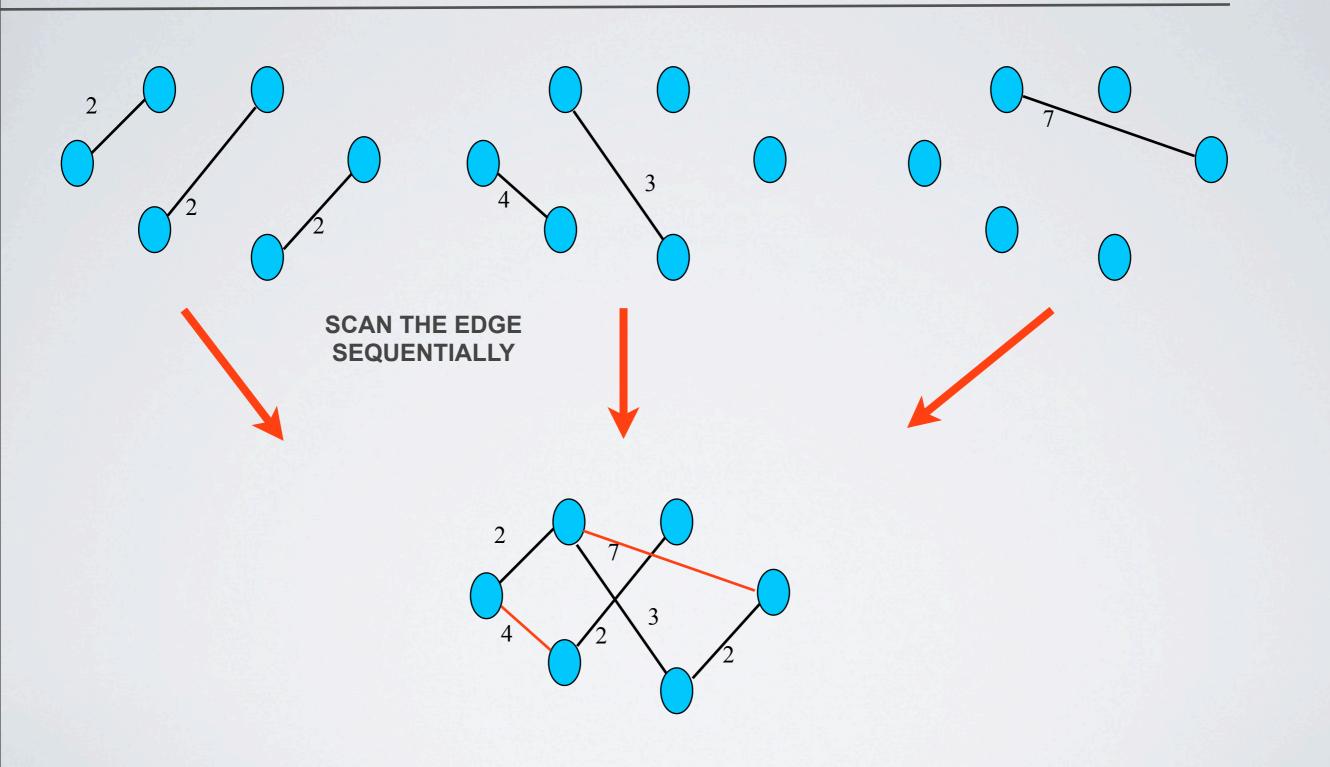












Bounding the rounds and memory

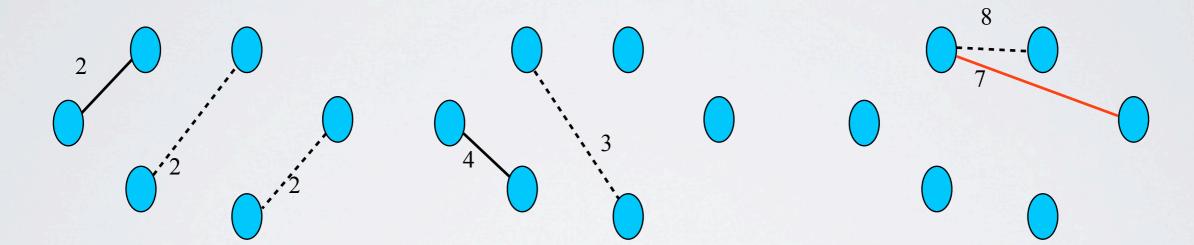
Four rounds:

•Splitting the graph and running the maximal matching: 3 rounds and $\tilde{O}(n^{1+c/2})$ memory

• Compute the final solution: 1 round and $O(n \log n)$ memory

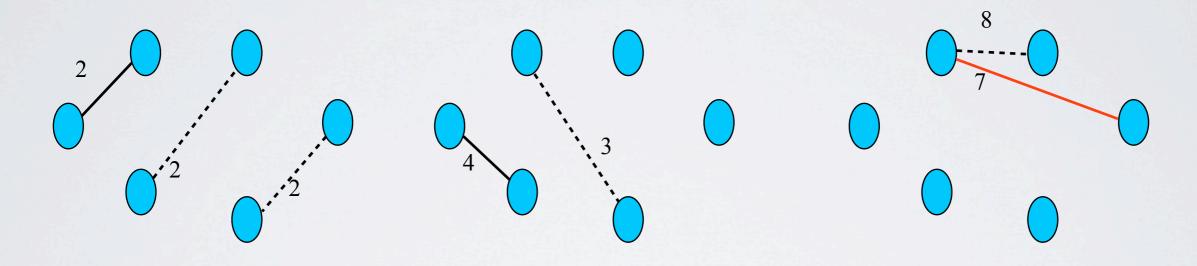
Approximation guarantee (sketch)

 An edge in the solution can block at most 2 edges in each subgraph of smaller weight



Approximation guarantee (sketch)

 An edge in the solution can block at most 2 edges in each subgraph of smaller weight



 We loose a factor or 2 because we do not consider the weight

Other algorithms

Based on the same intuition:

2-approx for vertex cover

3/2-approx for edge cover

Minimum cut

Partition does not work, because we loose structural informations

Minimum cut

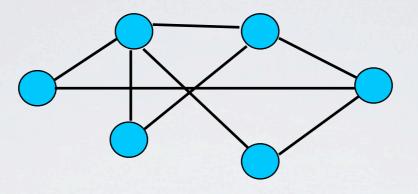
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Sampling does not seem to work either

Minimum cut

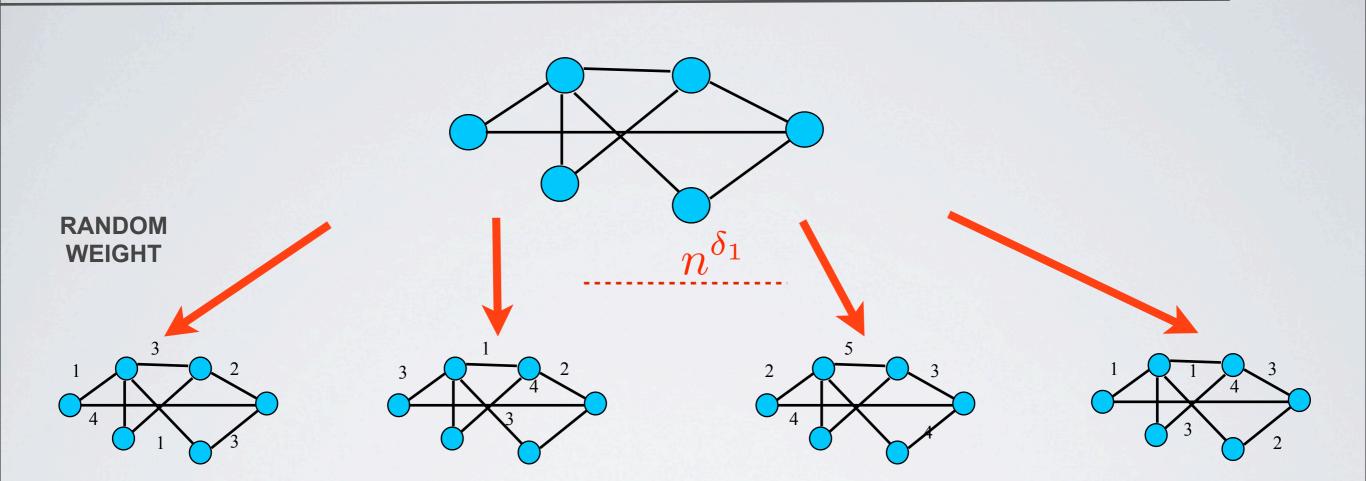
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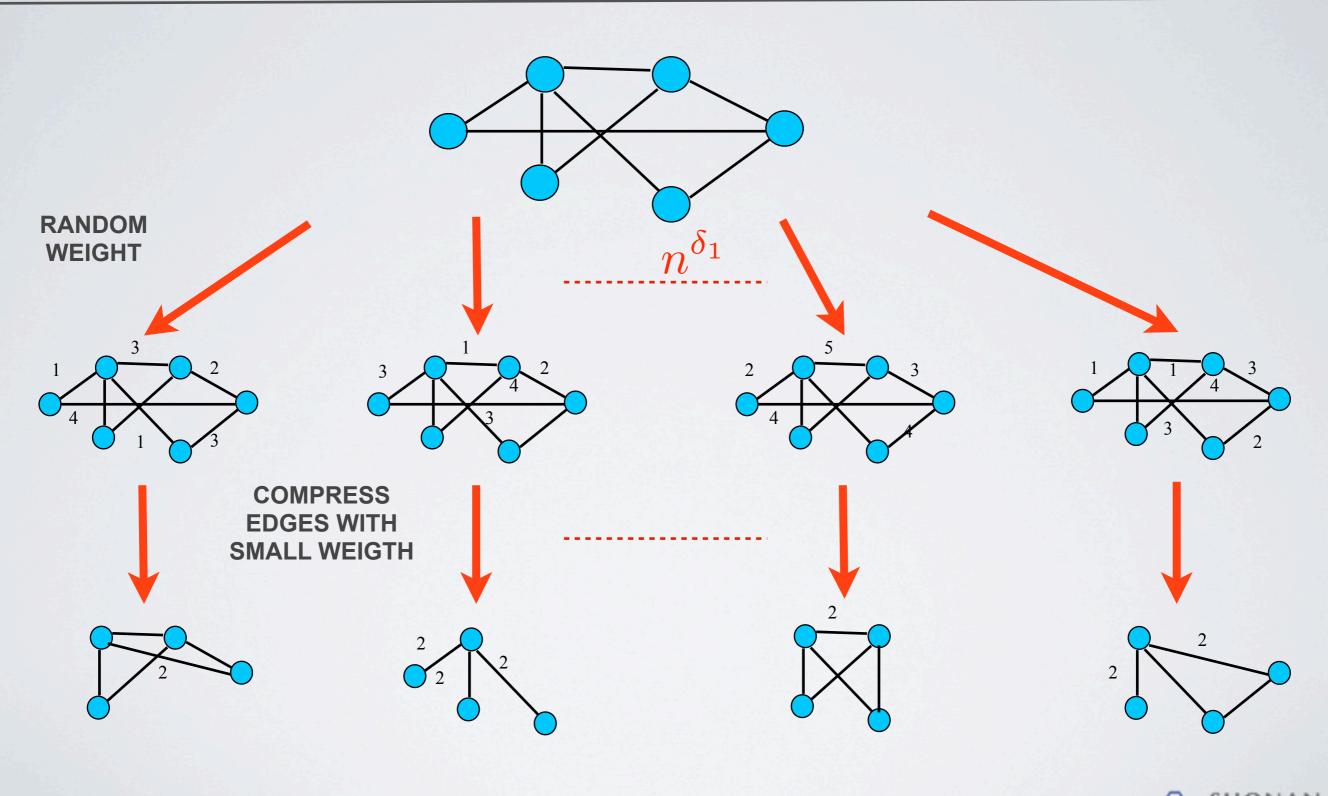
- Partition does not work, because we loose structural informations
- Sampling does not seem to work either
- We can use the first steps of Karger algorithm as a filtering technique
- The random choices made in the early rounds succeed with high probability, whereas the later rounds have a much lower probability of success

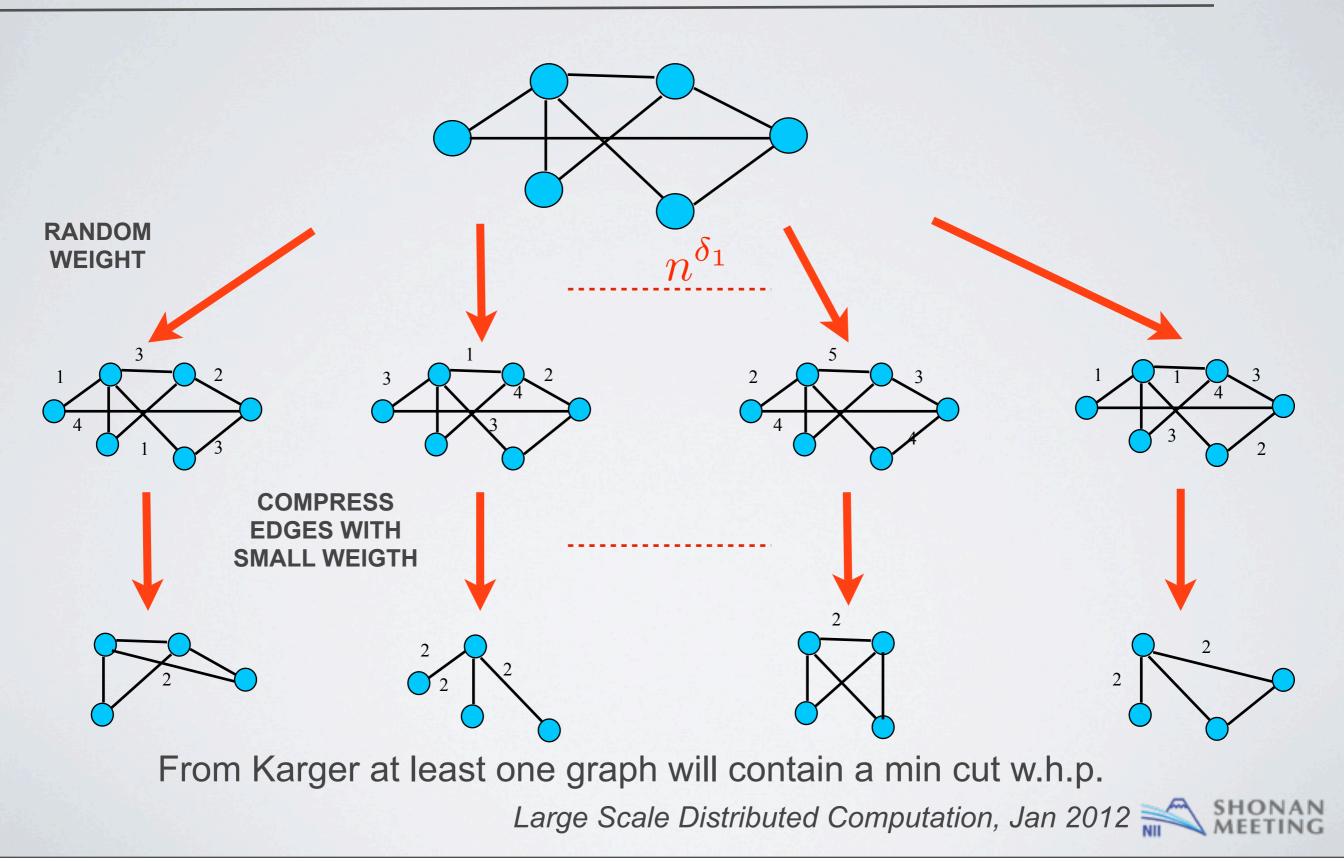


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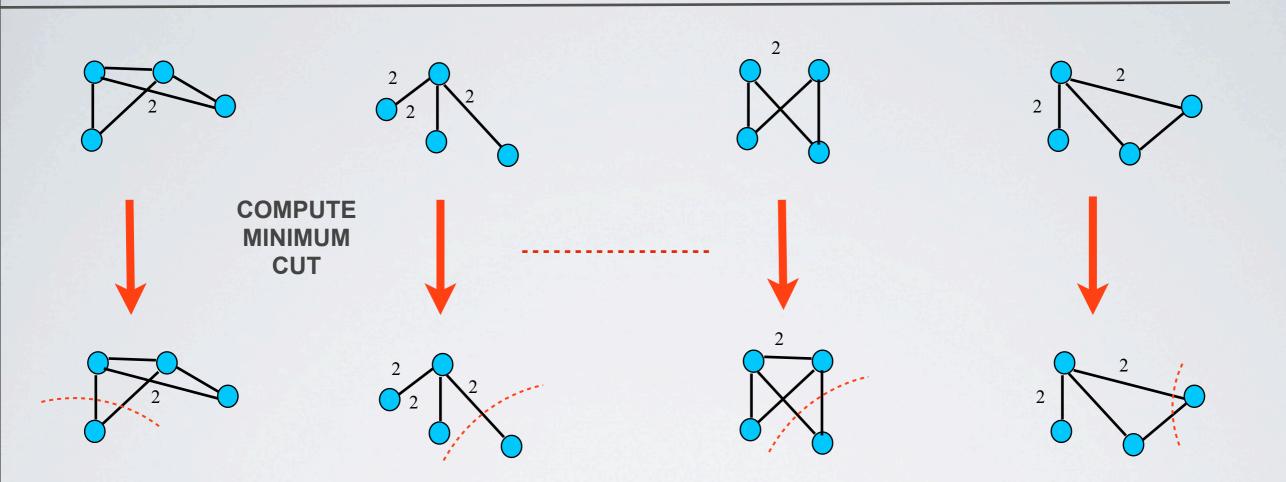






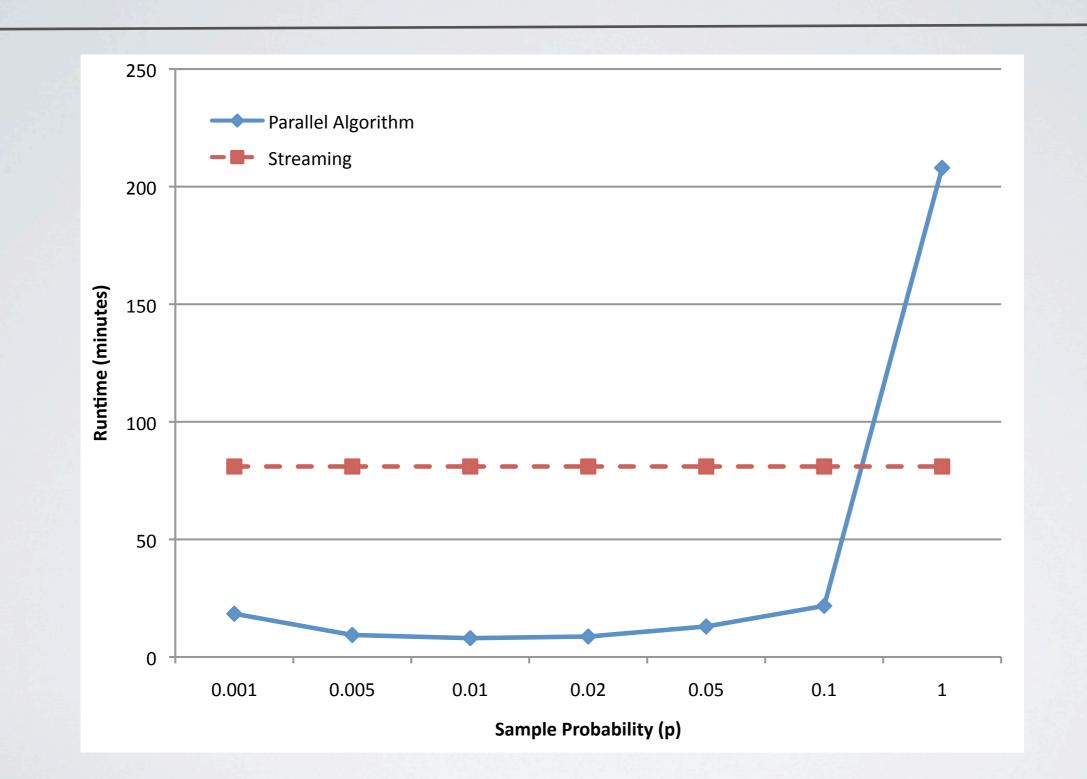
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We will find the minimum cut w.h.p.

Empirical result (matching)



Open problems

Open problems 1

Maximum matching

Shortest path

Dynamic programming

Open problems 2

Algorithms for sparse graph

Does connected components require more than 2 rounds?

Lower bounds

Thank you!

