

# *Seriously Equivalent + Scalable Parallel Machine Learning*

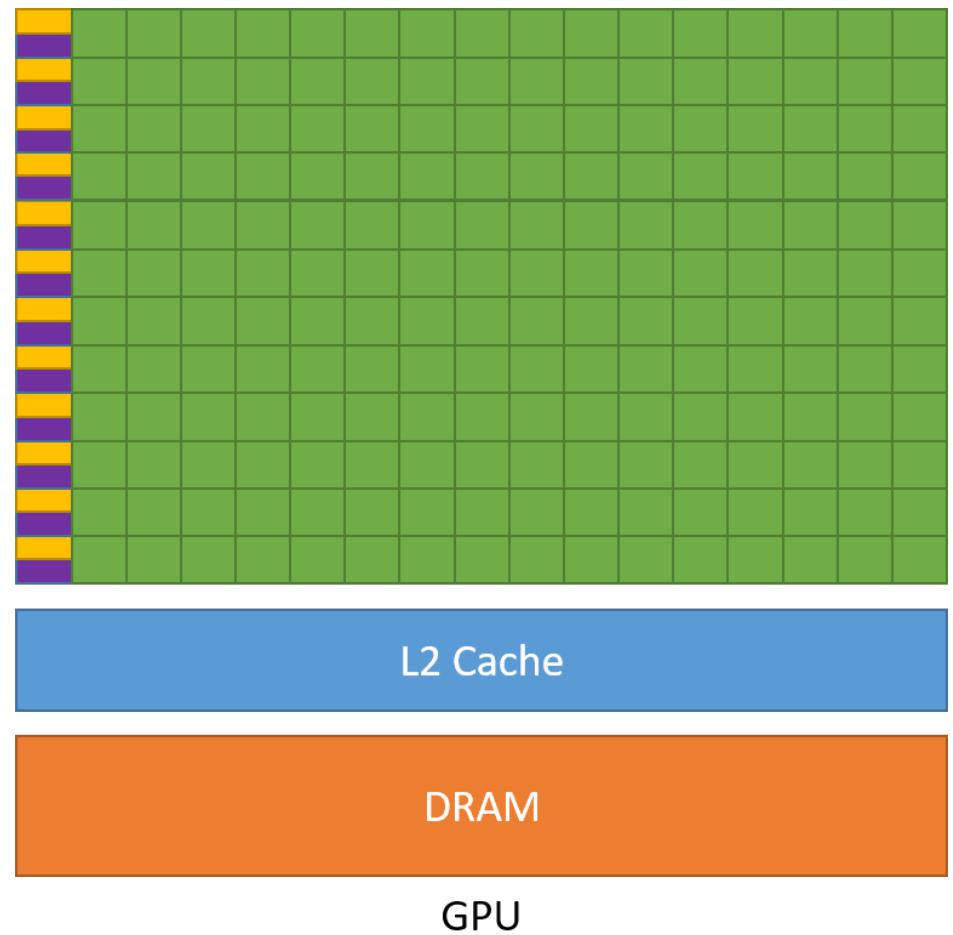
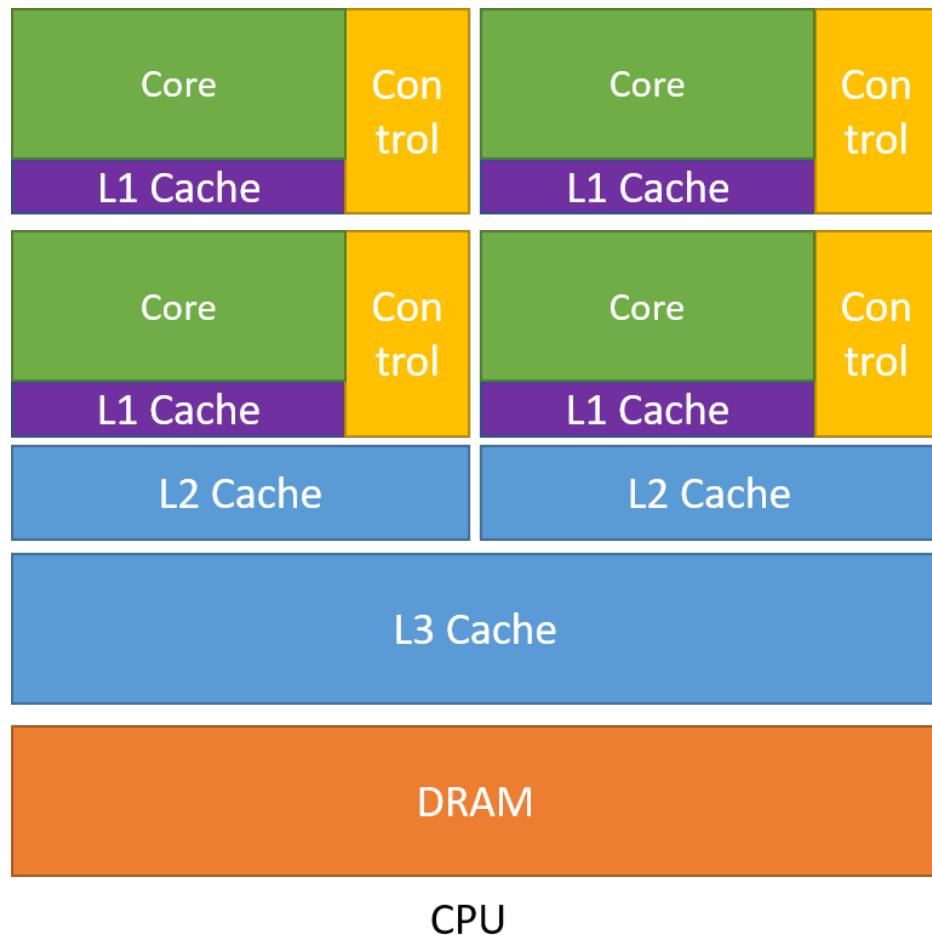


Dimitris Papailiopoulos

# Today

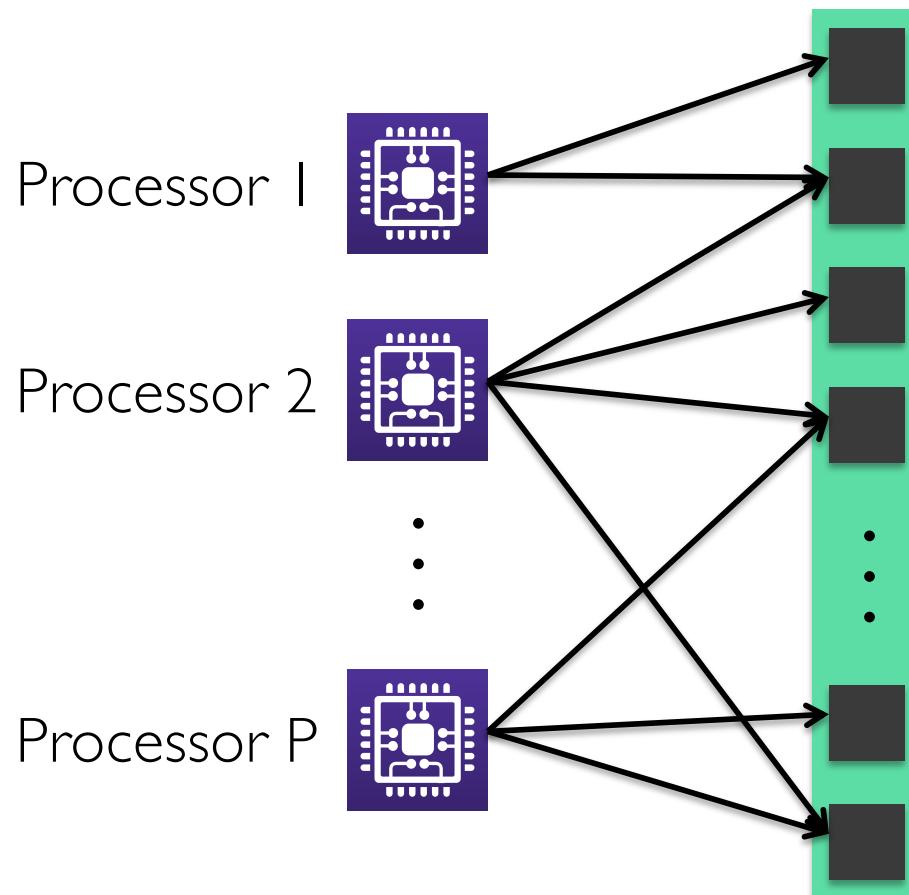
- Serial Equivalence
- Beyond Hogwild
- How much asynchrony is possible?
- Open Problems

# Single Machine, Multi-core



# Today

## (Maximally) Asynchronous



# Today

## Serial Equivalence

$$A_{\text{serial}}(S, \pi) = A_{\text{parallel}}(S, \pi)$$

For all Data sets S

For all data order  $\pi$  (data points can be arbitrarily repeated)

### Main advantage:

- we only need to “prove” speedups
- Convergence proofs inherited directly from serial

### Main Issue:

- Serial equivalence too strict
- Cannot guarantee any speedups in the general case

# The Stochastic Updates Meta-algorithm

# Stochastic Updates

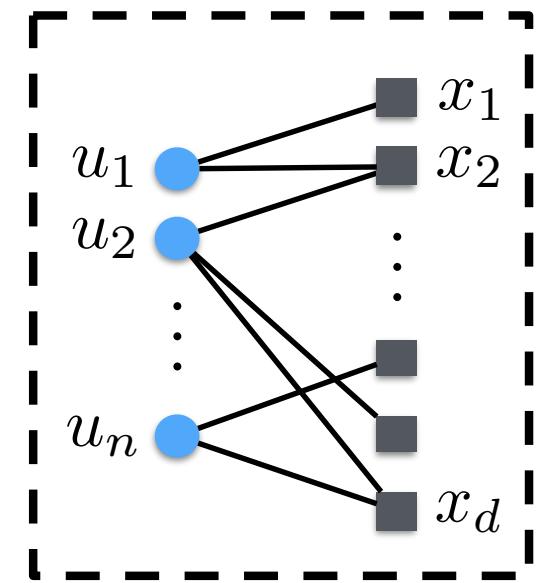
---

## Algorithm 1 Stochastic Updates pseudo-algorithm

---

```
1: Input:  $\mathbf{x}$ ;  $f_1, \dots, f_n$ ;  $u_1, \dots, u_n$ ;  $\mathcal{D}$ ;  $T$ .  
2: for  $t = 1 : T$  do  
3:   sample  $i \sim \mathcal{D}$   
4:    $\mathbf{x}_{\mathcal{S}_i} = u_i(\mathbf{x}_{\mathcal{S}_i}, f_i)$  //update global model on  $\mathcal{S}_i$   
5: Output:  $\mathbf{x}$ 
```

---



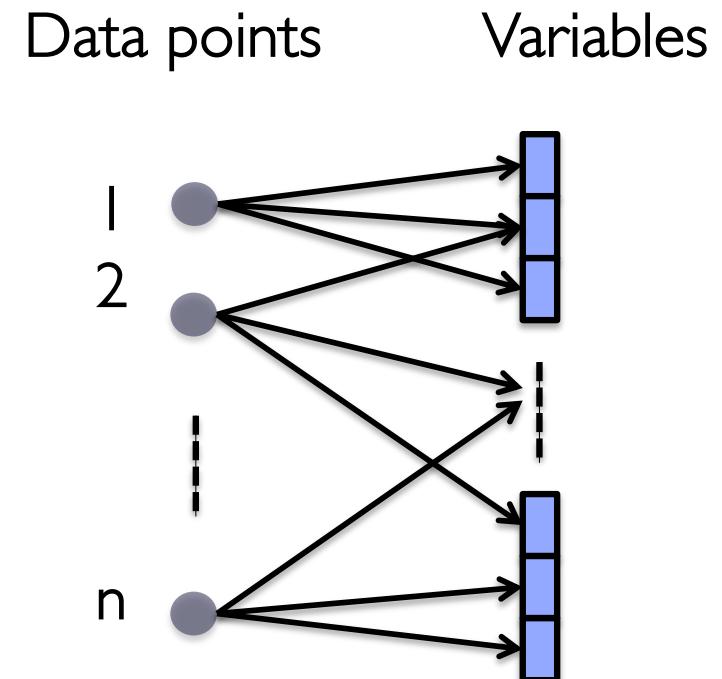
What does this solve?

# Stochastic Updates: A family of ML Algorithms

Many algorithms with sparse access patterns:

- SGD
- SVRG / SAGA
- Matrix Factorization
  - word2vec
  - K-means
- Stochastic PCA
- Graph Clustering

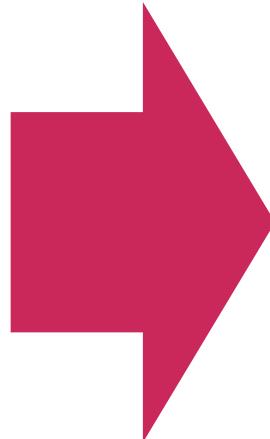
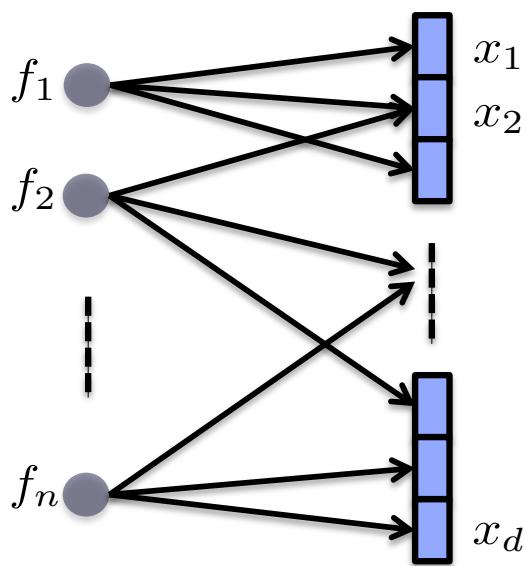
...



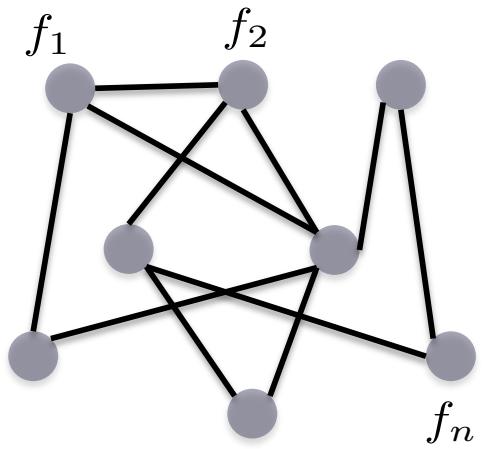
Can we parallelize under Serial Equivalence?

# A graph view of Conflicts in Parallel Updates

# The Update Conflict Graph



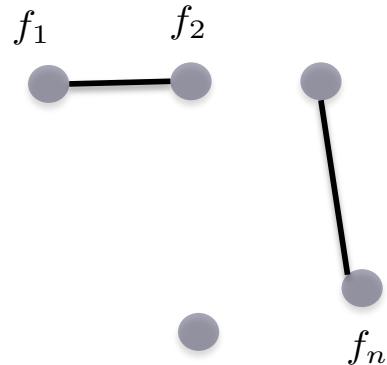
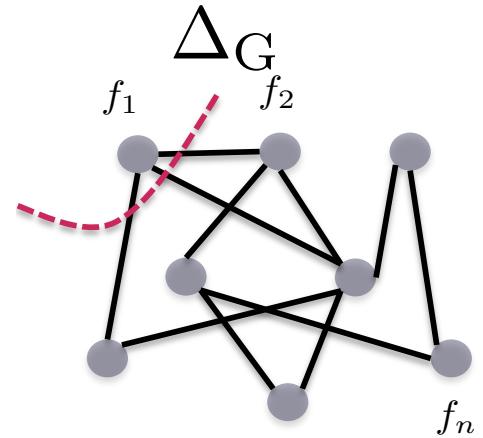
conflict graph



An edge between 2 updates if they overlap

# The Theorem [Krivelevich'14]

conflict graph

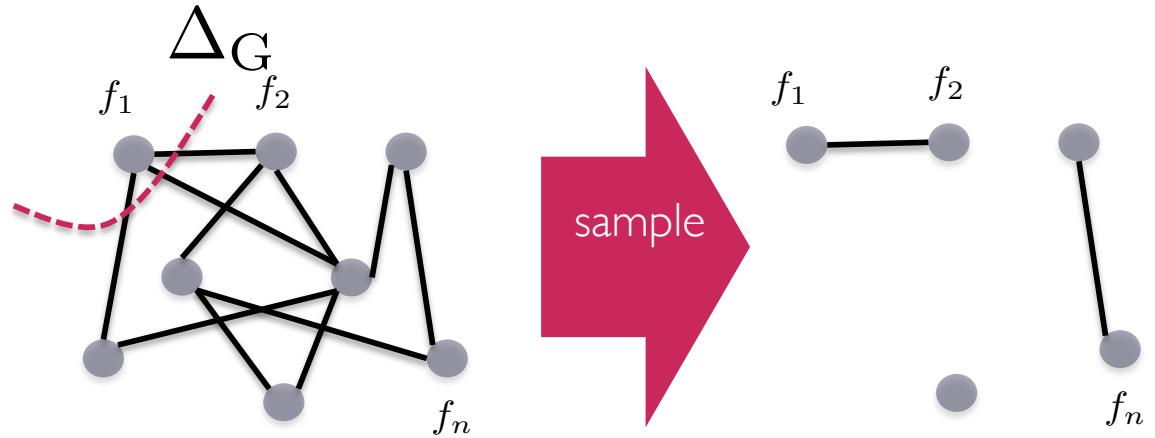


Lemma:

Sample less than  $P \leq (1 - \epsilon) \frac{n}{\Delta_G}$  vertices (with/without replacement)

# The Theorem [Krivelevich'14]

conflict graph



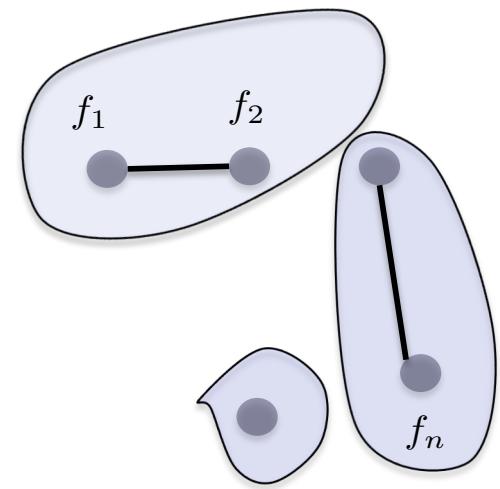
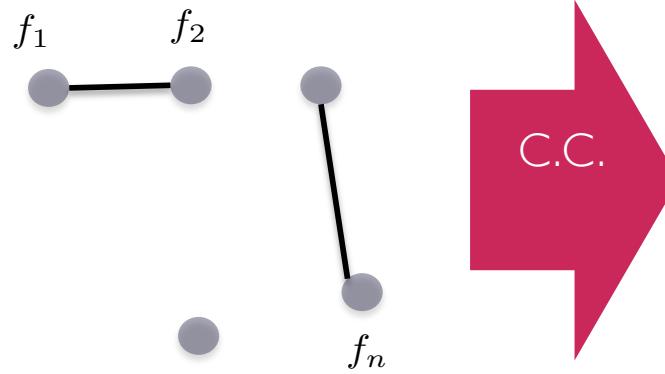
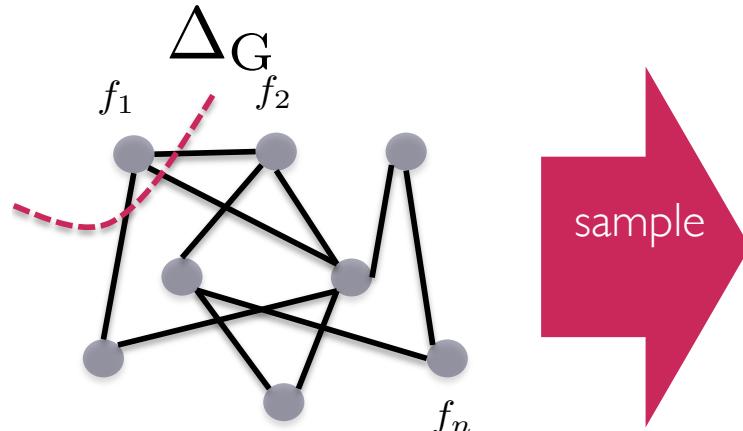
Lemma:

Sample less than  $P \leq (1 - \epsilon) \frac{n}{\Delta_G}$  vertices (with/without replacement)

Then, the induced sub-graph shatters

# The Theorem [Krivelevich'14]

conflict graph



Lemma:

Sample less than  $P \leq (1 - \epsilon) \frac{n}{\Delta_G}$  vertices (with/without replacement)

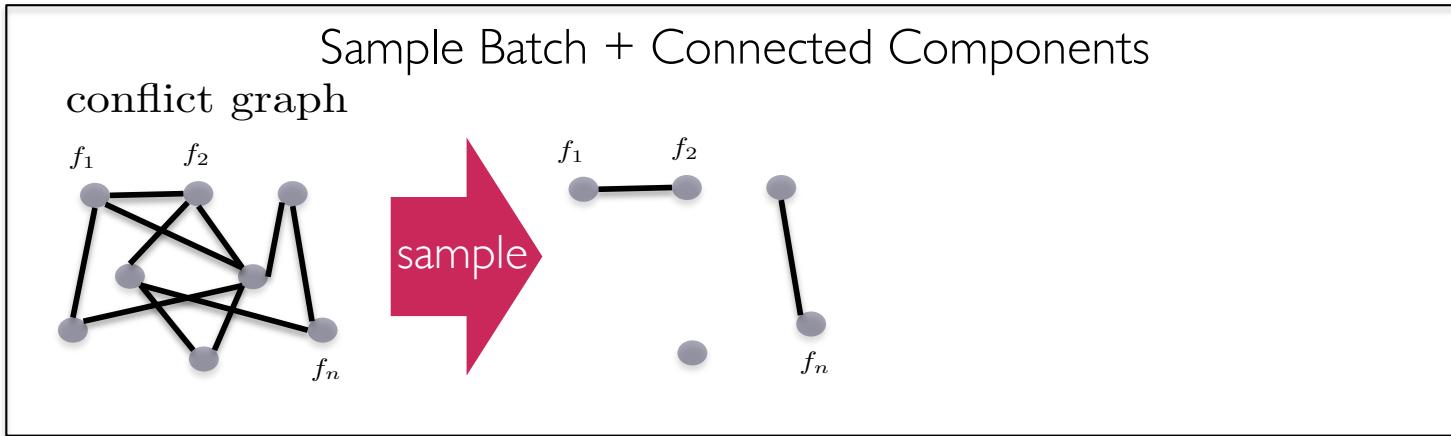
Then, the induced sub-graph shatters,  
The largest connected component has size

$$O\left(\frac{\log n}{\epsilon^2}\right)$$

Even if the Graph was a Single Huge Conflict Component!

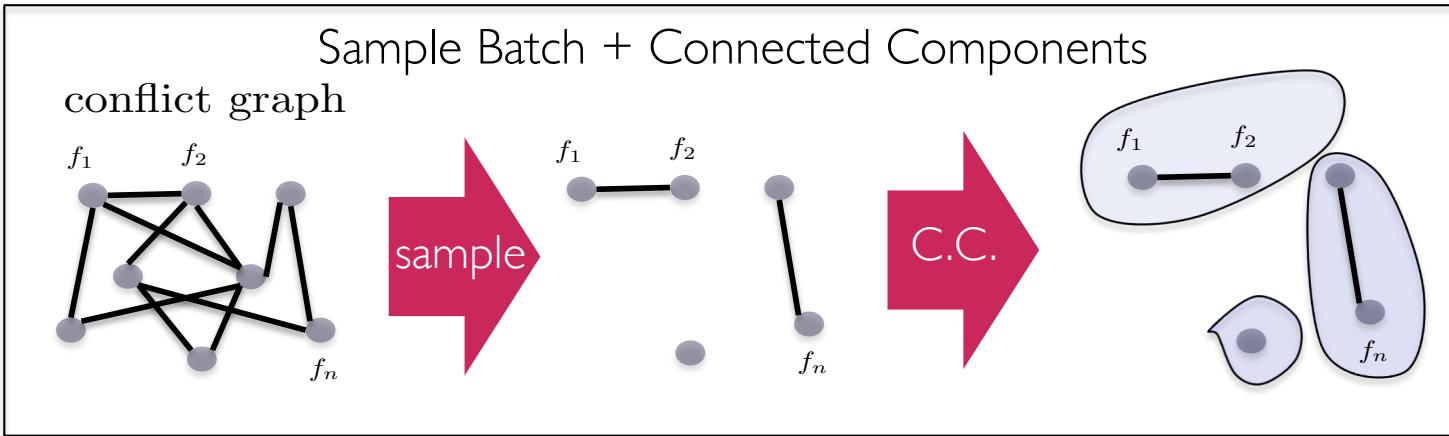
# Building a Parallelization Framework out of a Single Theorem

Phase I



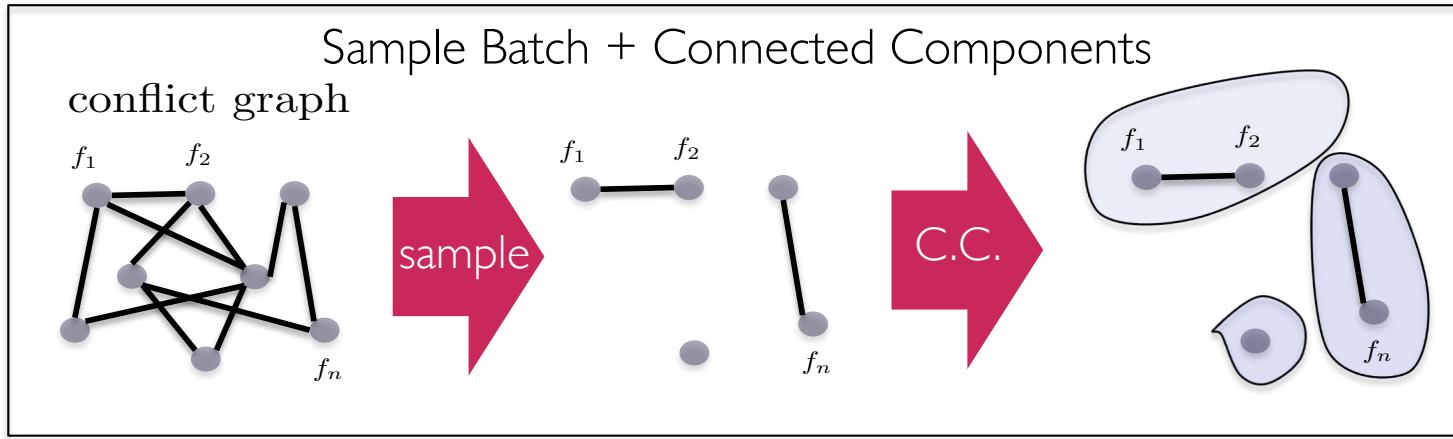
Sample  $B = (1 - \epsilon) \cdot \frac{n}{\Delta}$  vertices

Phase I



$$\text{Sample } B = (1 - \epsilon) \cdot \frac{n}{\Delta} \text{ vertices}$$

Phase I



Sample  $B = (1 - \epsilon) \cdot \frac{n}{\Delta}$  vertices

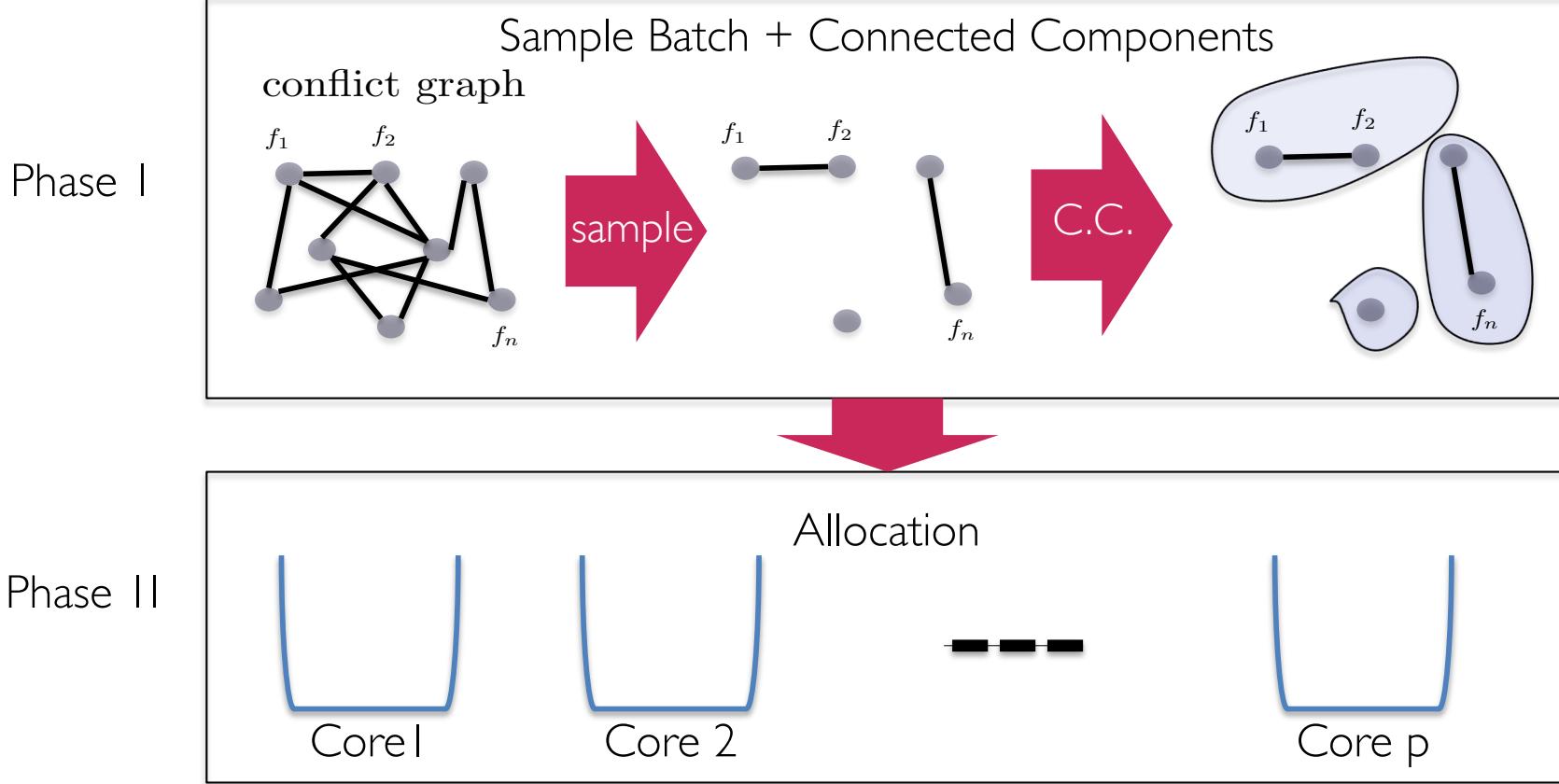
Compute Conn. Components

NOTE: No conflicts **across** groups!

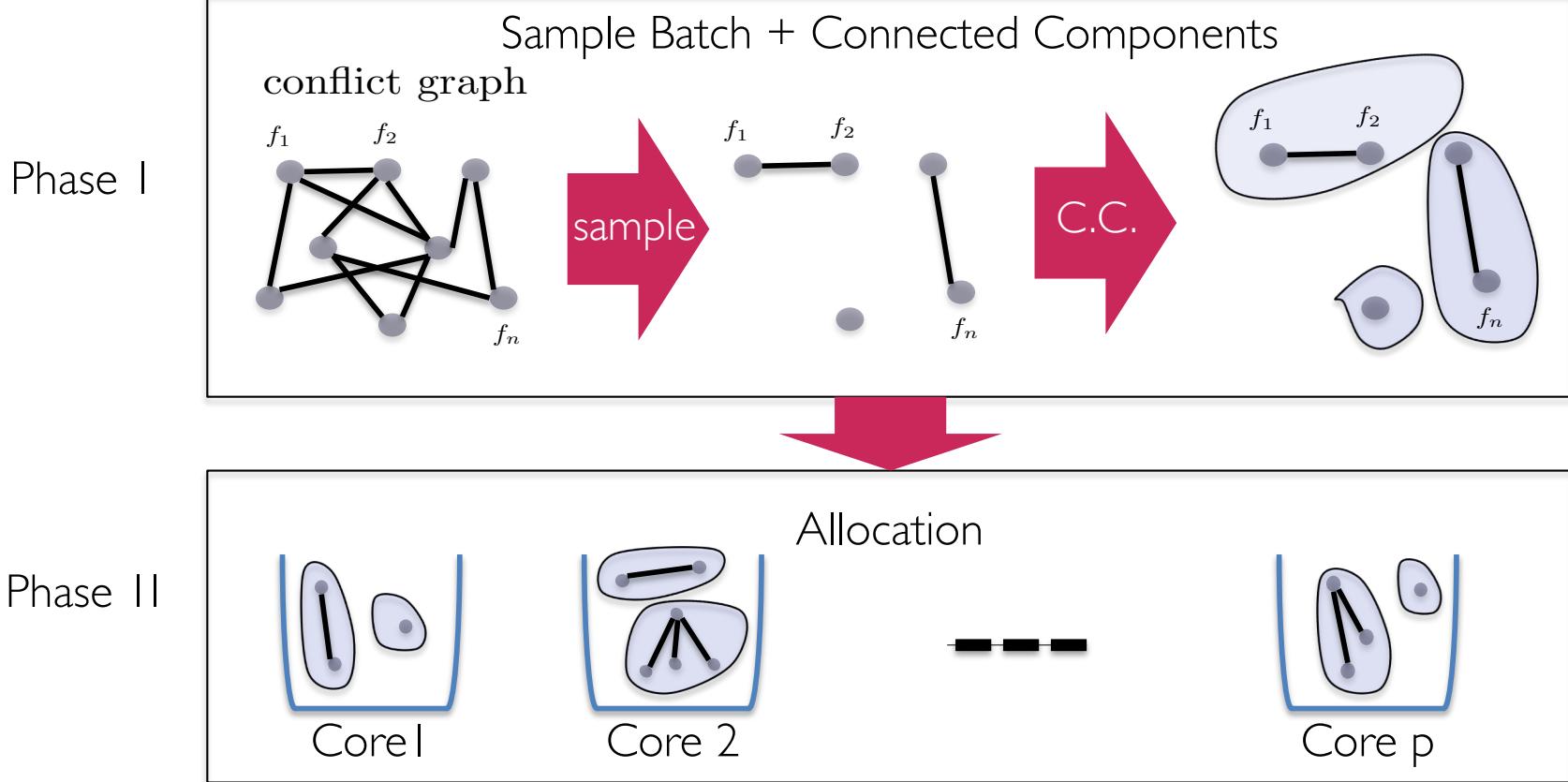
Max Conn. Comp =  $\log n \Rightarrow n/(\Delta \log n)$  tiny components

Yay! Good for parallelization

No conflicts across groups = we can run Stochastic Updates on each of them in parallel!

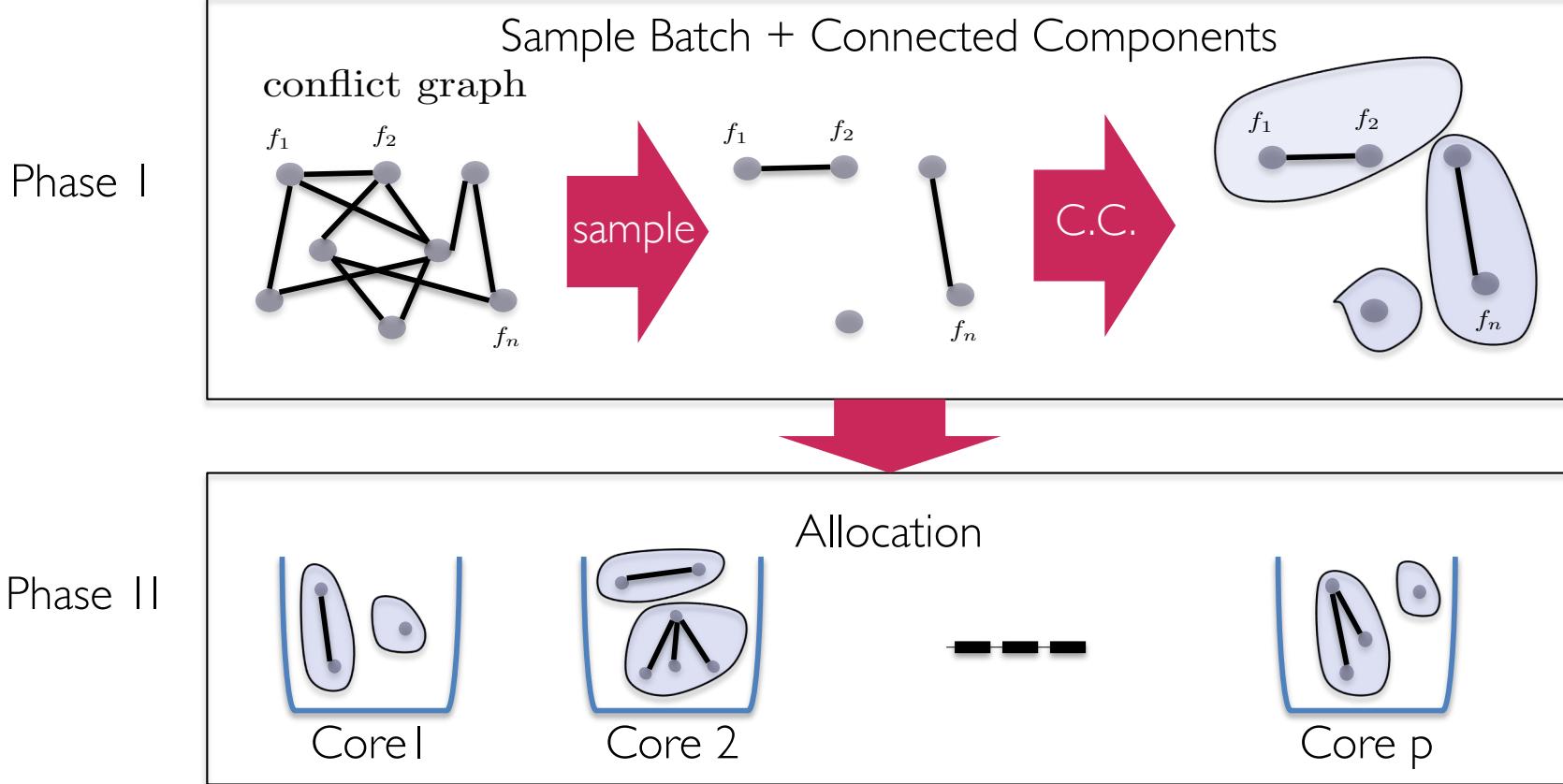


$$\text{Cores} < \text{Batch size} / \log n = n / (\Delta \log n)$$



$$\text{Cores} < \text{Batch size} / \log n = n / (\Delta \log n)$$

A Single Rule: Run the updates serially inside each connected component  
Automatically Satisfied since we give each **conflict group** to a single core.

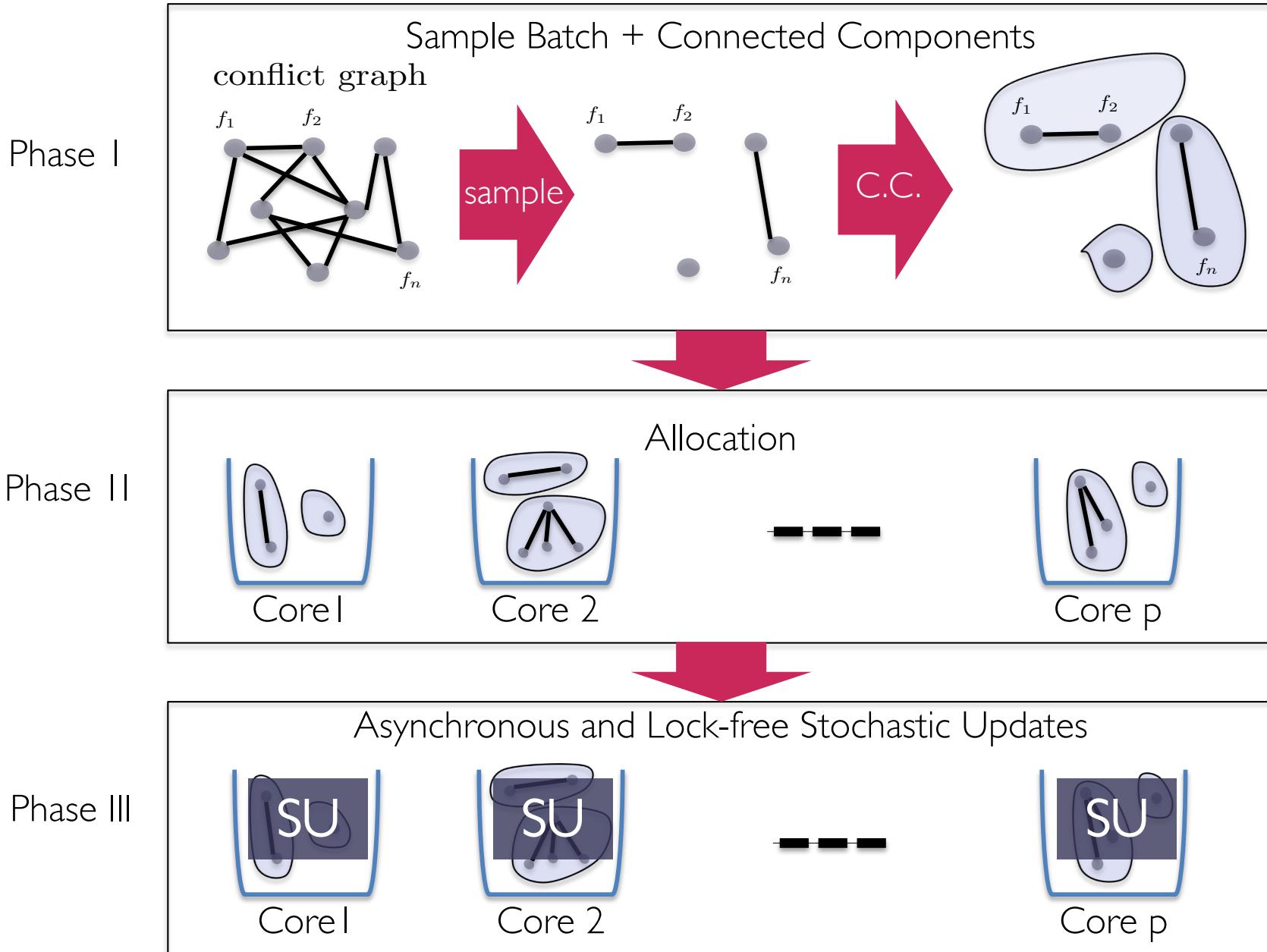


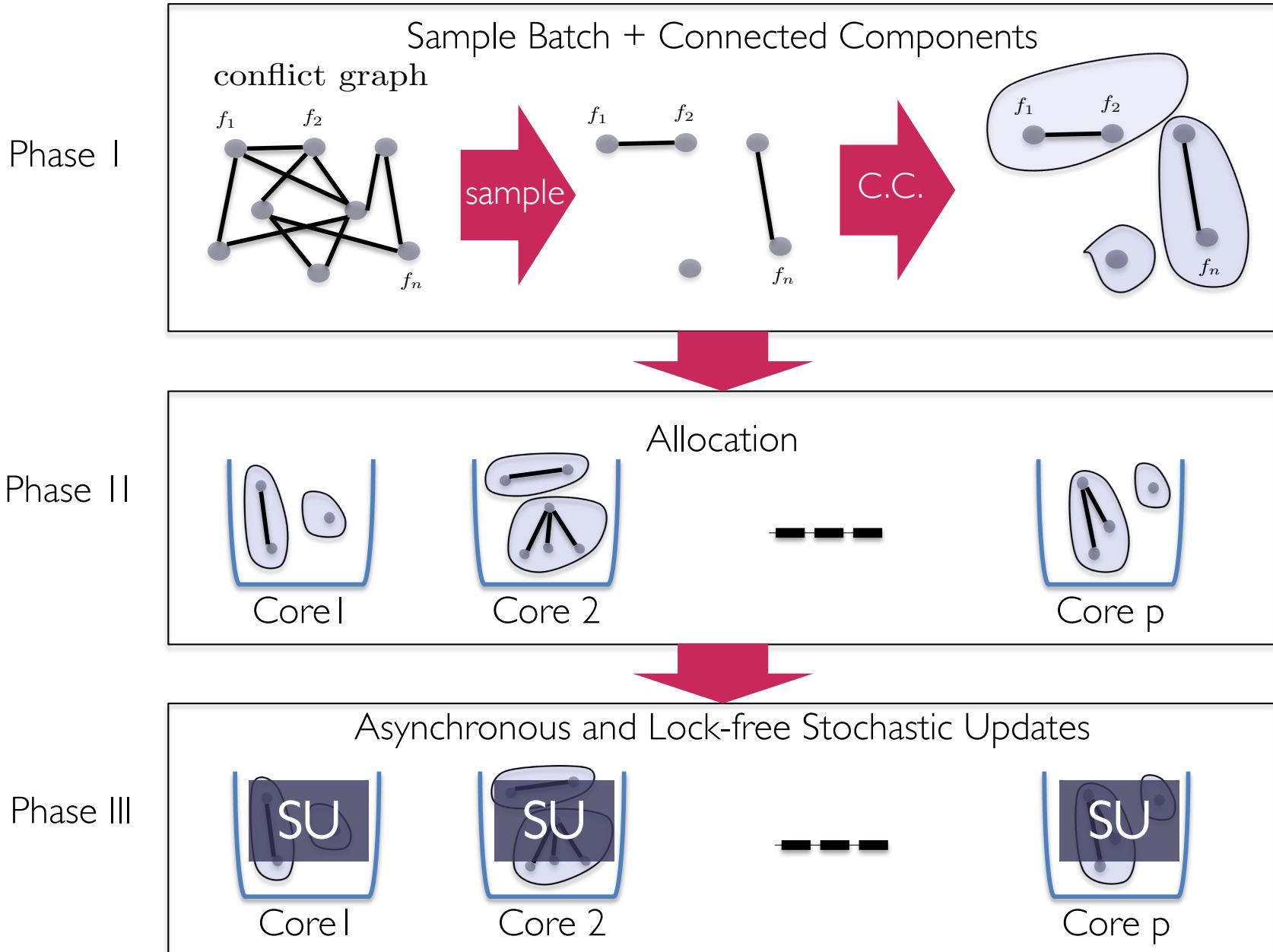
Policy I: Random allocation, good when cores  $\ll$  Batch size

Policy II: Greedy min-weight allocation

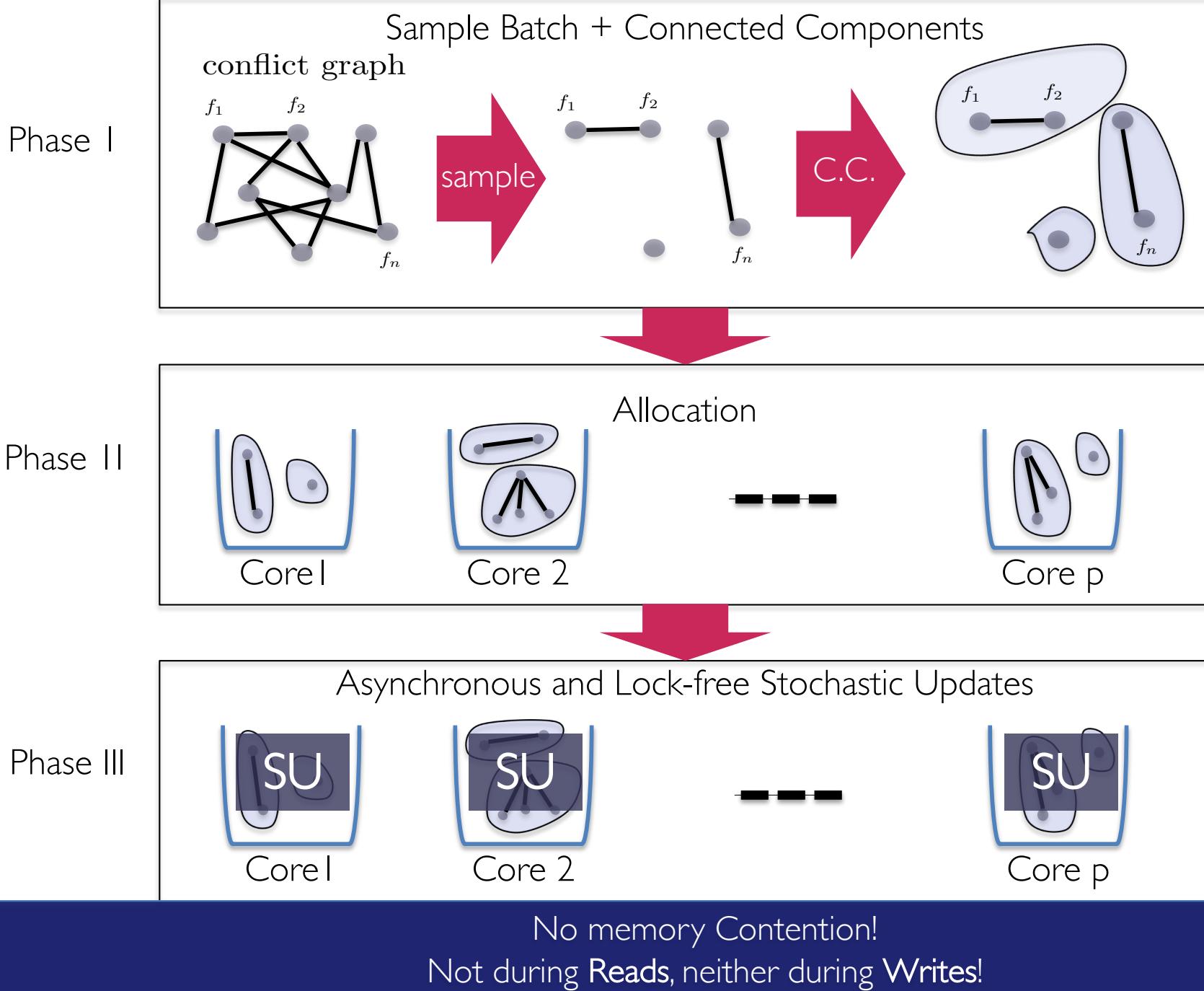
*(80% as good as optimal (which is NP-hard))*

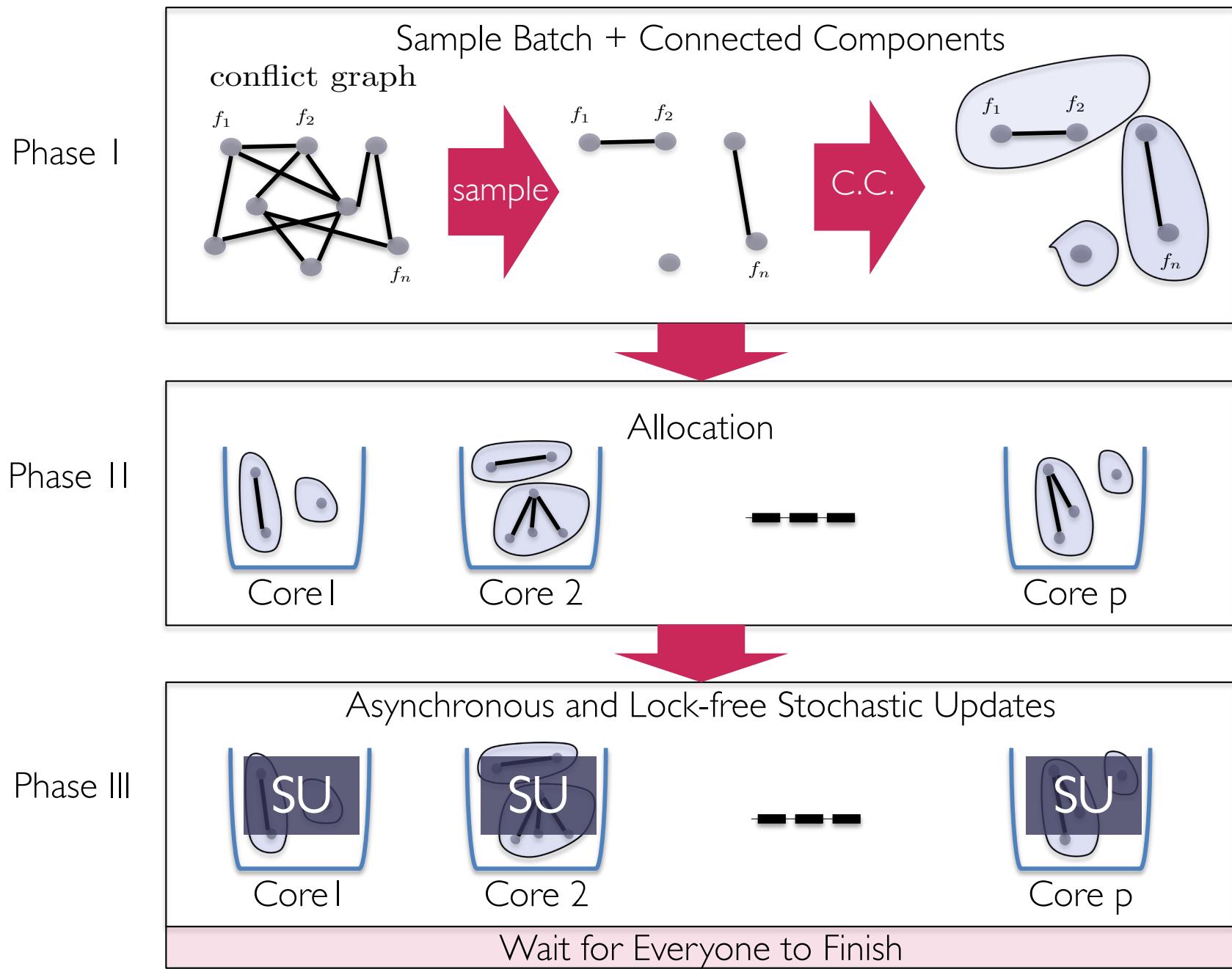
A Single Rule: Run the updates serially inside each connected component  
Automatically Satisfied since we give each **conflict group** to a single core.

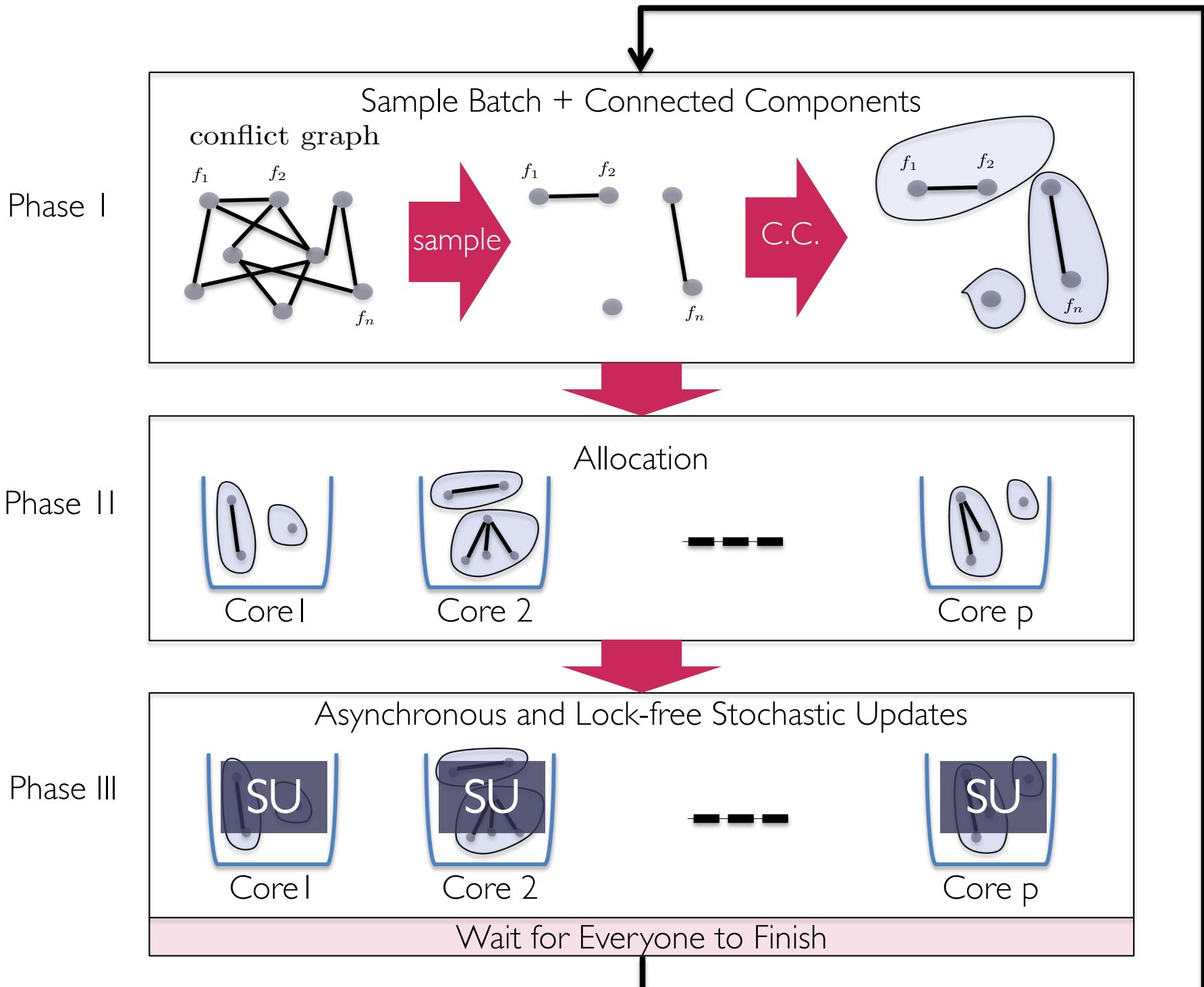




Each core runs Asynchronously and Lock-free! (No communication!)







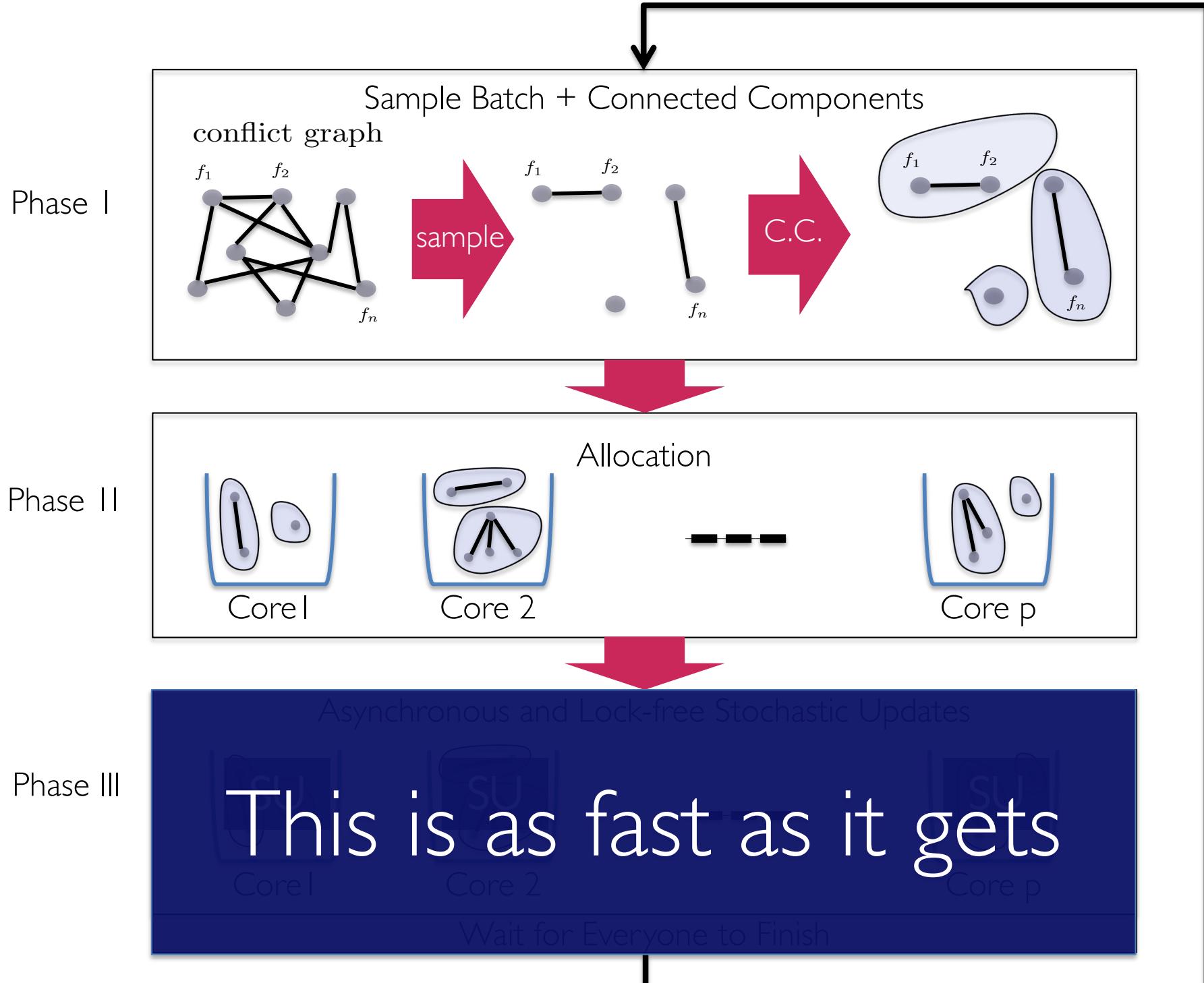
---

## Algorithm 2 CYCLADES

---

- 1: **Input:**  $G_u, T, B$ .
  - 2: Sample  $n_b = T/B$  subgraphs  $G_u^1, \dots, G_u^{n_b}$  from  $G_u$
  - 3: Cores compute in parallel CCs for sampled subgraphs
  - 4: **for** batch  $i = 1 : n_b$  **do**
  - 5:     Allocation of  $\mathcal{C}_1^i, \dots, \mathcal{C}_{m_i}^i$  to  $P$  cores
  - 6:     **for** each core in parallel **do**
  - 7:         **for** each allocated component  $\mathcal{C}$  **do**
  - 8:             **for** each update  $j$  (in order) from  $\mathcal{C}$  **do**
  - 9:                  $\mathbf{x}_{\mathcal{S}_j} = u_j(\mathbf{x}_{\mathcal{S}_j}, f_j)$
  - 10:     **Output:**  $\mathbf{x}$
-

This guarantees Serially Equivalence  
But does it guarantee speedups?



Phase I

Sample Batch + Connected Components  
conflict graph



Serial Cost:  $O\left(\frac{n}{\Delta} \cdot \text{sparsity} \cdot \log n\right)$

If Phase I and II are fast  
We are good.

Phase II

Allocation



Serial Cost:

$O\left(\frac{n}{\Delta} \log n\right)$

Phase III

Asynchronous and Lock-free Stochastic Updates

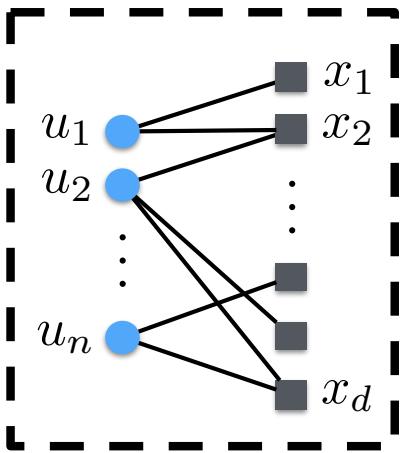


This is as fast as it gets

Wait for Everyone to Finish



# Main Theorem



Note:

$$\text{Serial Cost} = E_u \cdot \kappa$$

$\kappa$  = cost / coordinate update

**Theorem 4.** Let us assume any given update-variable graph  $G_u$  with average, and max left degree  $\bar{\Delta}_L$  and  $\Delta_L$ , such that  $\frac{\Delta_L}{\bar{\Delta}_L} \leq \sqrt{n}$ , and with induced max conflict degree  $\Delta$ . Then, CYCLADES on  $P = O(\frac{n}{\Delta \cdot \bar{\Delta}_L})$  cores, with batch sizes  $B = (1 - \epsilon) \frac{n}{\Delta}$  can execute  $T = c \cdot n$  updates, for any constant  $c \geq 1$ , selected uniformly at random with replacement, in time

$$\mathcal{O}\left(\frac{E_u \cdot \kappa}{P} \cdot \log^2 n\right), \quad \rightarrow \quad \text{Speedup} = \frac{P}{\log^2 n}$$

with high probability.

Assumptions:

- 1) Not too large max degree (approximate “regularity”)
- 2) Not too many cores
- 3) Sampling according to the “Graph Theorem”

Phase I

Identical performance as serial for

- SGD

- SVRG / SAGA

Phase II

- Sparse Network training

- Matrix Factorization

- Word2Vec

- Matrix Completion

- Greedy Clustering

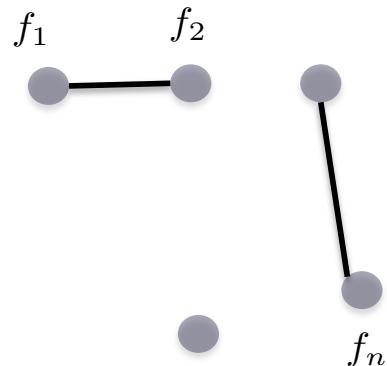
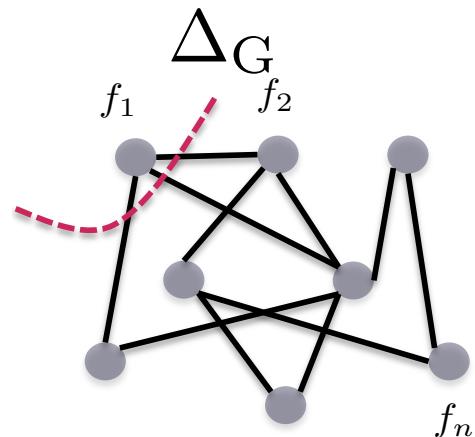
Phase III

Wait for Everyone to Finish

# Fast Connected Components

# Fast Connected Components

conflict graph

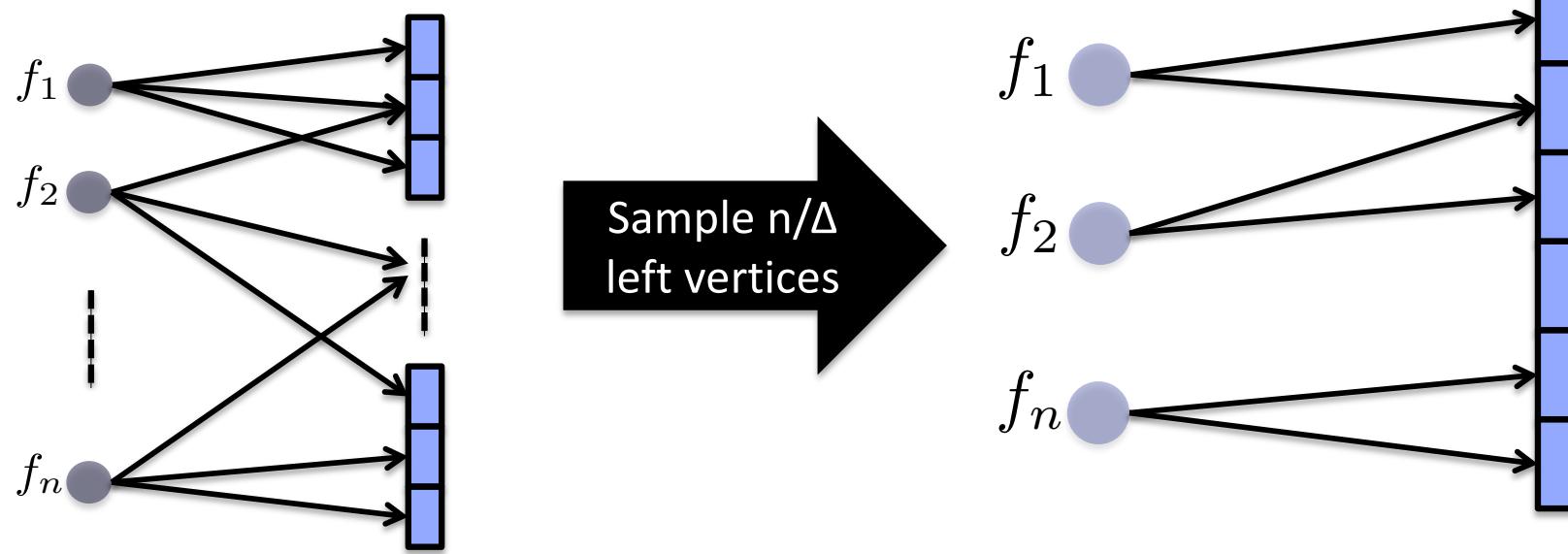


If you have the conflict graph CC is easy...  $O(\text{Sampled Edges})$

Building the conflict graph requires  $n^2$  time...

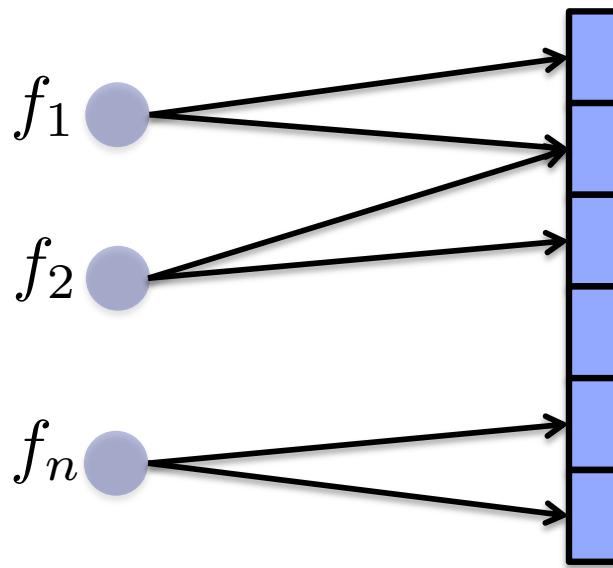
No, thanks.

# Fast Connected Components



Sample on the Bipartite, not on the Conflict Graphs

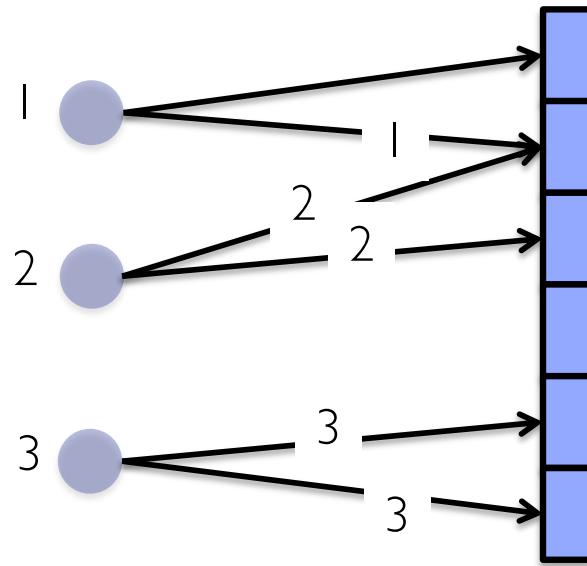
# Fast Connected Components



## Simple Message passing Idea

- Gradients send their IDs
  - Coordinates Compute Min and Send Back
  - Gradients Compute Min and Send Back
- Iterate till you're done

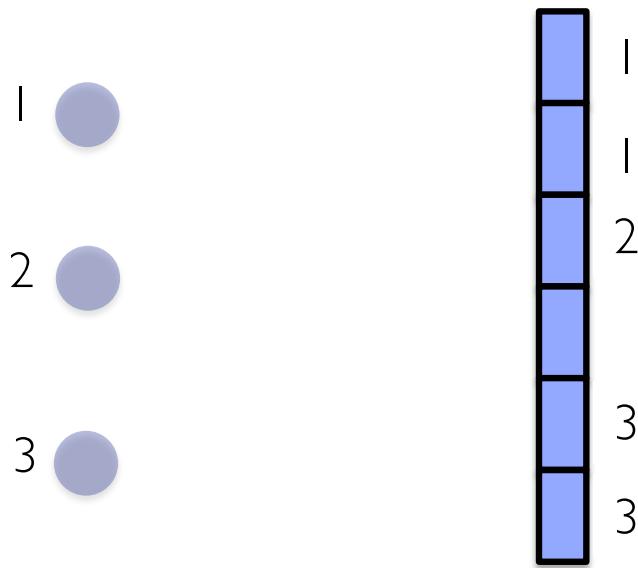
# Fast Connected Components



## Simple Message passing Idea

- Gradients send their IDs
  - Coordinates Compute Min and Send Back
  - Gradients Compute Min and Send Back
- Iterate till you're done

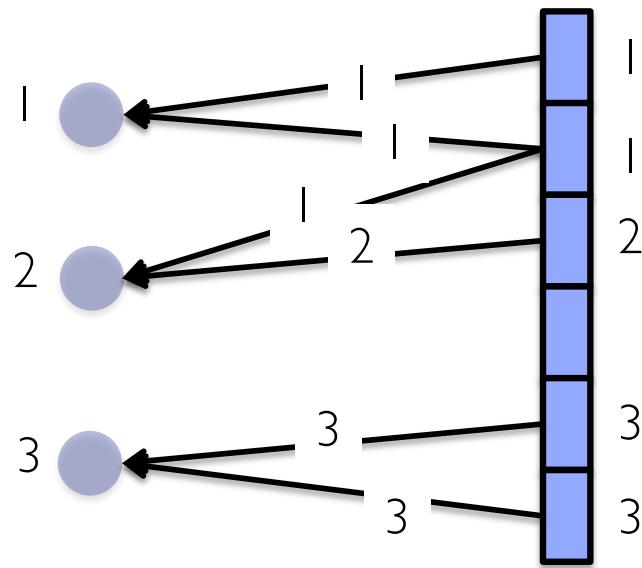
# Fast Connected Components



## Simple Message passing Idea

- Gradients send their IDs
  - Coordinates Compute Min and Send Back
  - Gradients Compute Min and Send Back
- Iterate till you're done

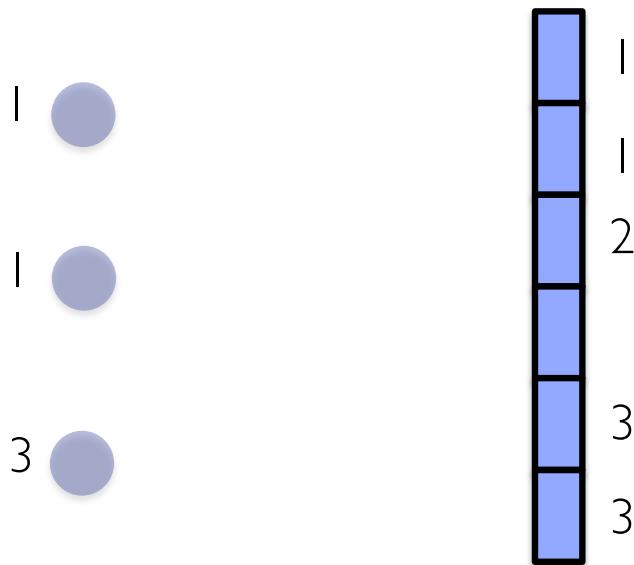
# Fast Connected Components



## Simple Message passing Idea

- Gradients send their IDs
  - Coordinates Compute Min and Send Back
  - Gradients Compute Min and Send Back
- Iterate till you're done

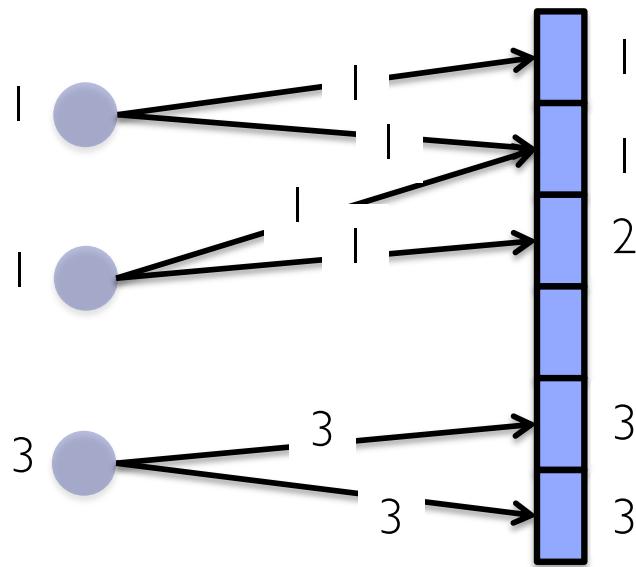
# Fast Connected Components



## Simple Message passing Idea

- Gradients send their IDs
  - Coordinates Compute Min and Send Back
  - Gradients Compute Min and Send Back
- Iterate till you're done

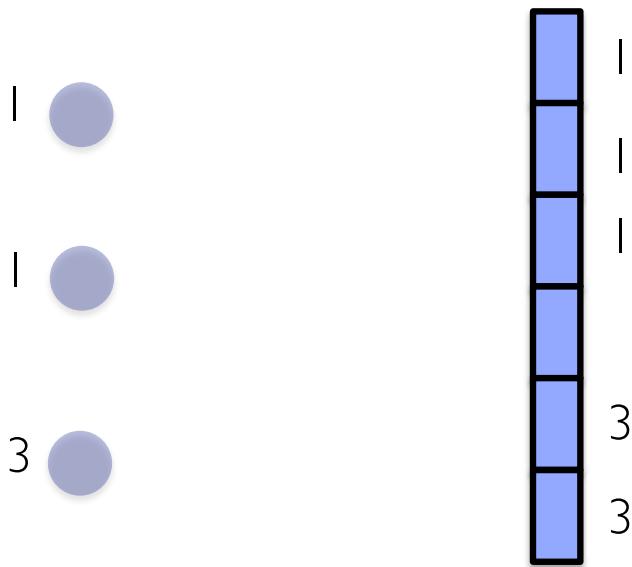
# Fast Connected Components



## Simple Message passing Idea

- Gradients send their IDs
  - Coordinates Compute Min and Send Back
  - Gradients Compute Min and Send Back
- Iterate till you're done

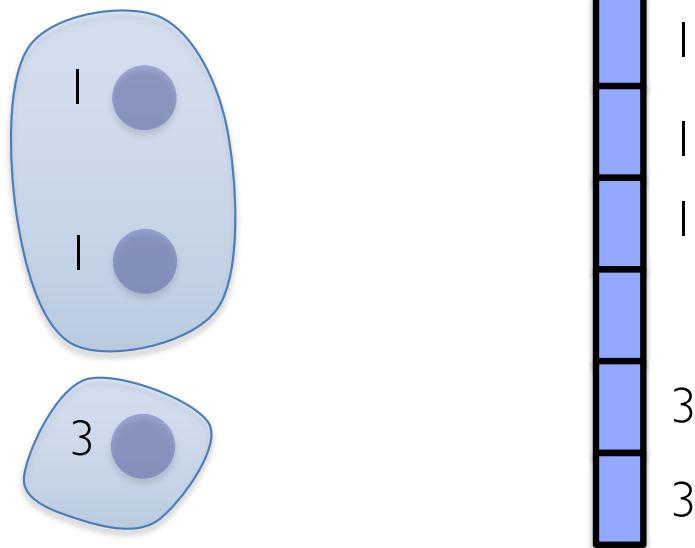
# Fast Connected Components



## Simple Message passing Idea

- Gradients send their IDs
  - Coordinates Compute Min and Send Back
  - Gradients Compute Min and Send Back
- Iterate till you're done

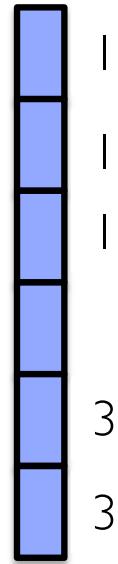
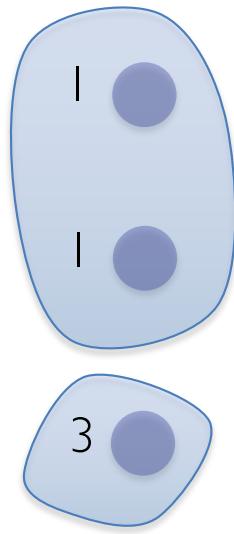
# Fast Connected Components



## Simple Message passing Idea

- Gradients send their IDs
  - Coordinates Compute Min and Send Back
  - Gradients Compute Min and Send Back
- Iterate till you're done

# Fast Connected Components

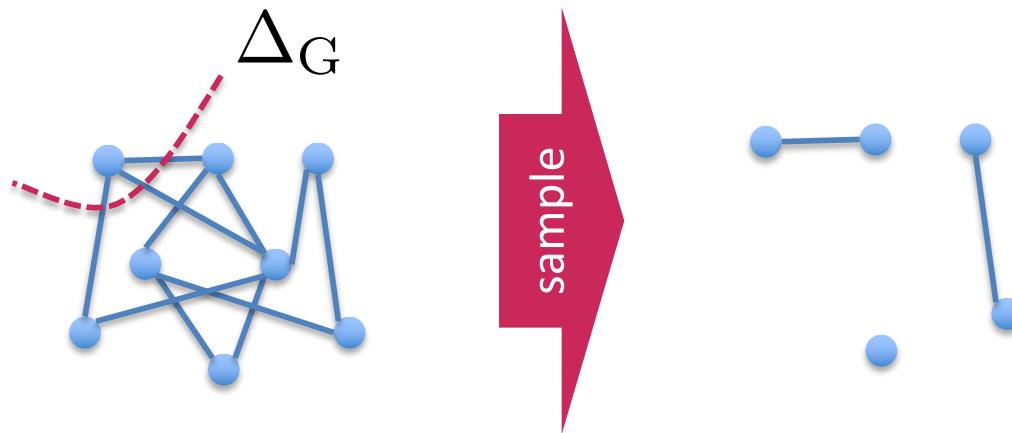


$$\text{Cost} = O\left(\frac{E_u \log^2 n}{P}\right)$$

Same assumptions as main theorem

# Proof of “The Theorem”

# The Theorem



Lemma:

Activate each vertex with probability

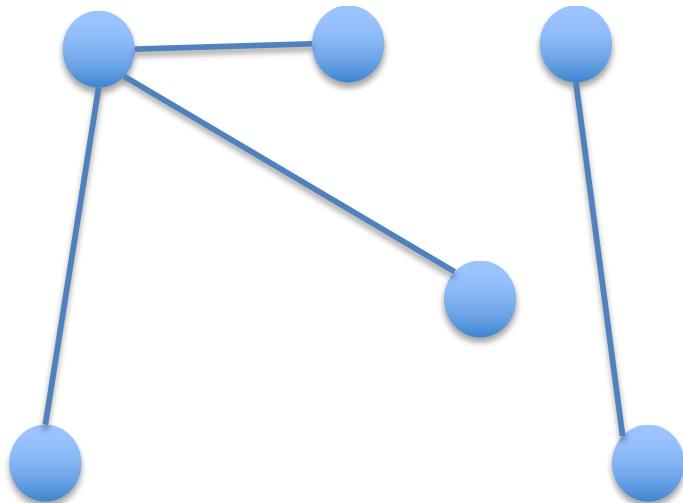
$$p = (1 - \varepsilon) / \Delta$$

Then, the induced subgraph shatters,  
and the largest connected component has size

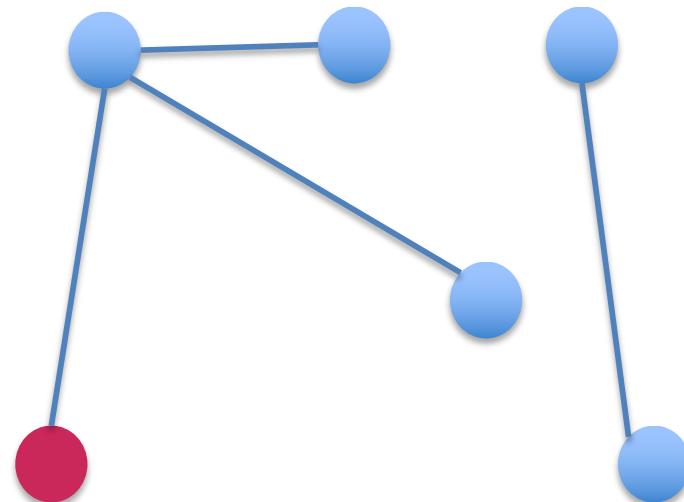
$$\frac{4}{\varepsilon^2} \cdot \log n$$

# DFS 101

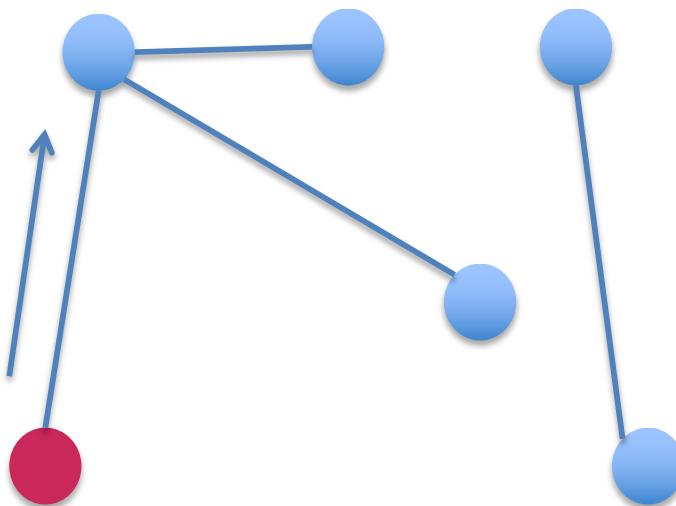
# DFS



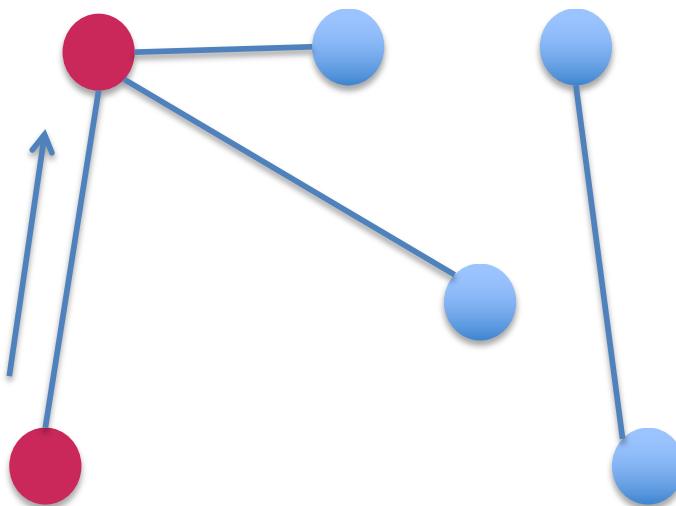
# DFS



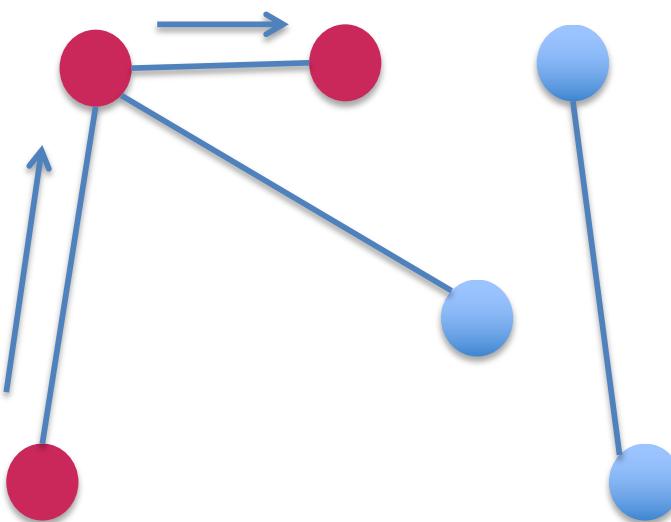
# DFS



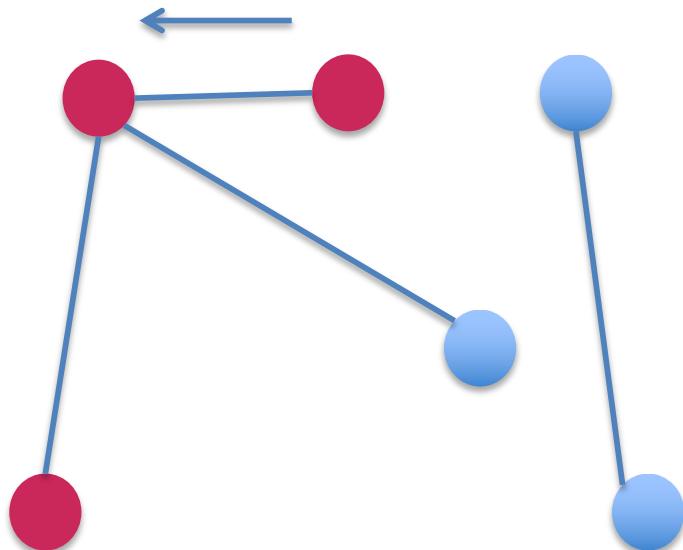
# DFS



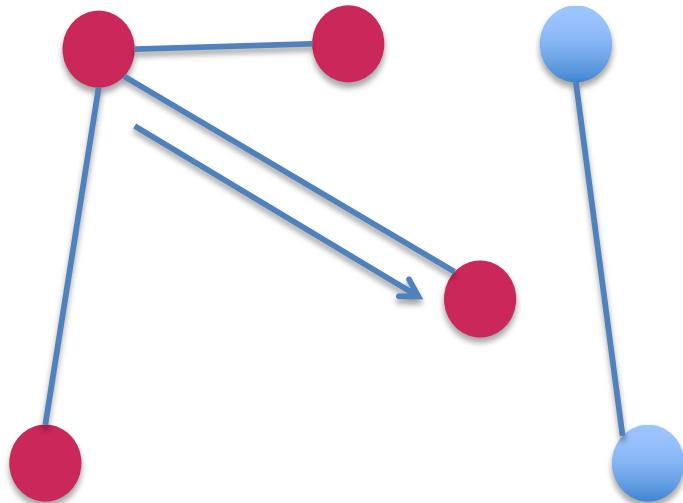
# DFS



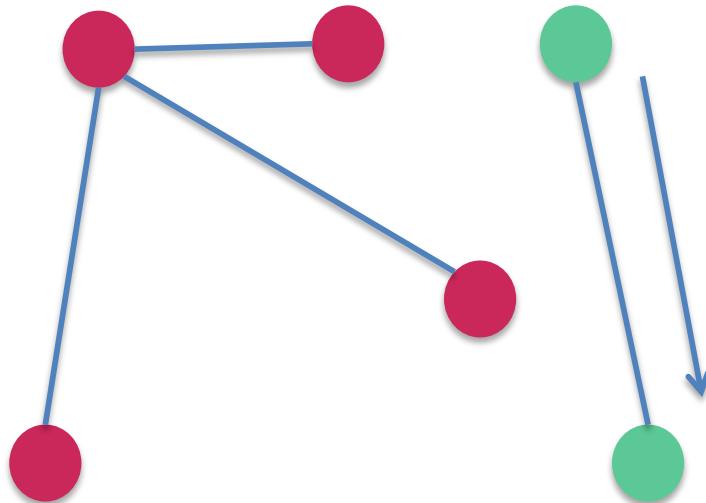
# DFS



# DFS

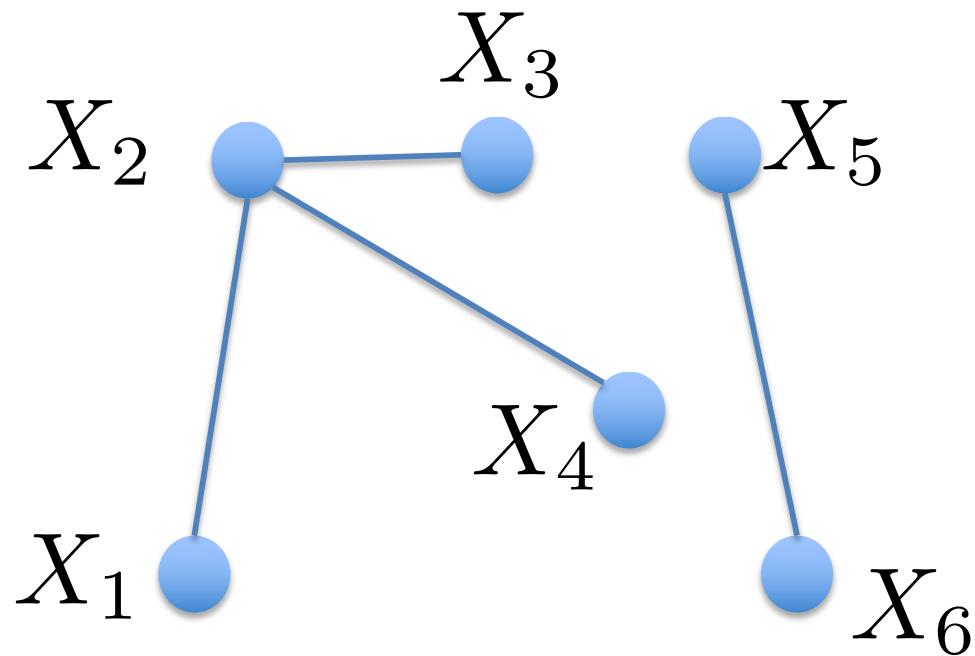


# DFS



# Probabilistic DFS

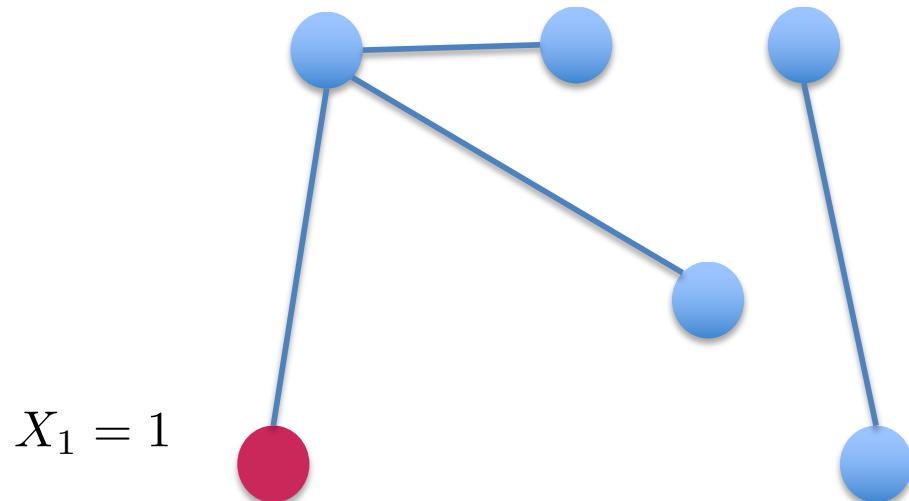
# DFS with random coins



Algorithm:

- Flip a coin for each vertex DFS wants to visit
- If 1 visit, if 0 don't visit and delete with its edges

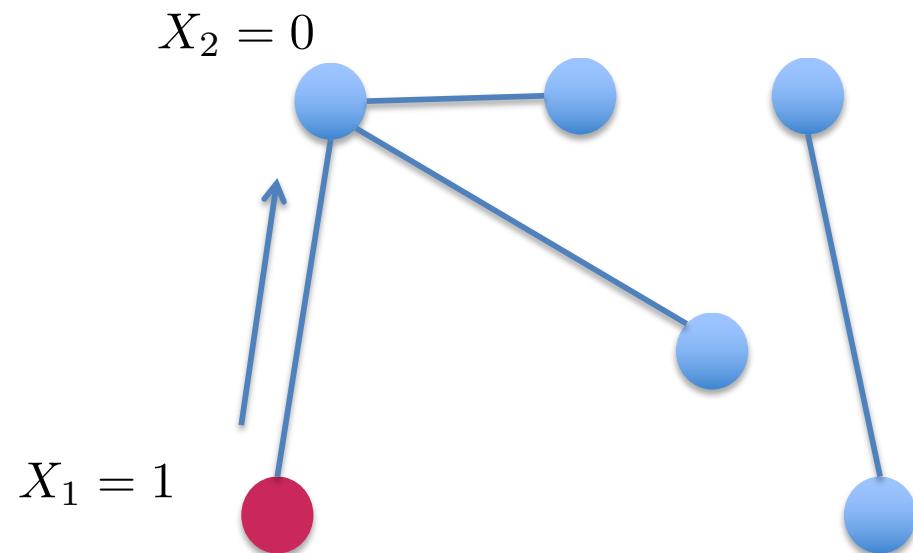
# DFS with random coins



Algorithm:

- Flip a coin for each vertex DFS wants to visit
- If 1 visit, if 0 don't visit and delete with its edges

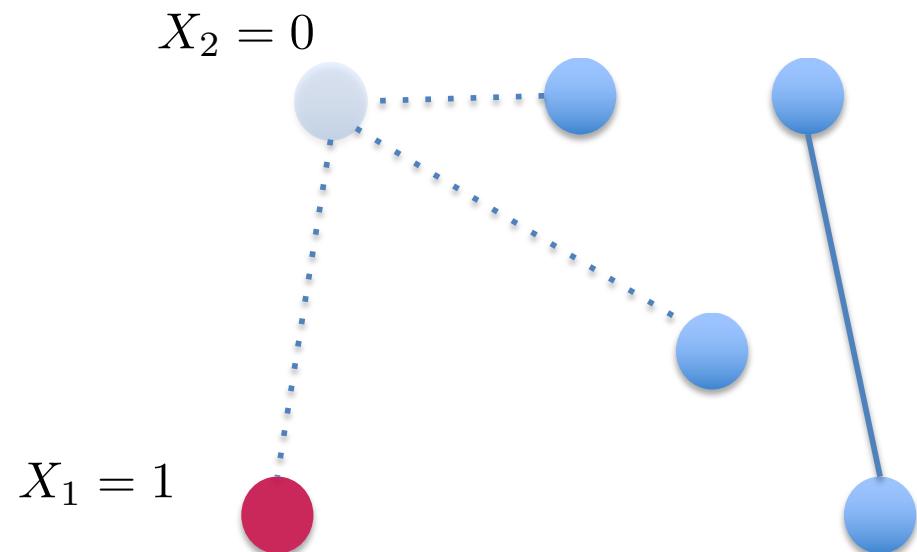
# DFS with random coins



Algorithm:

- Flip a coin for each vertex DFS wants to visit
- If 1 visit, if 0 don't visit and delete with its edges

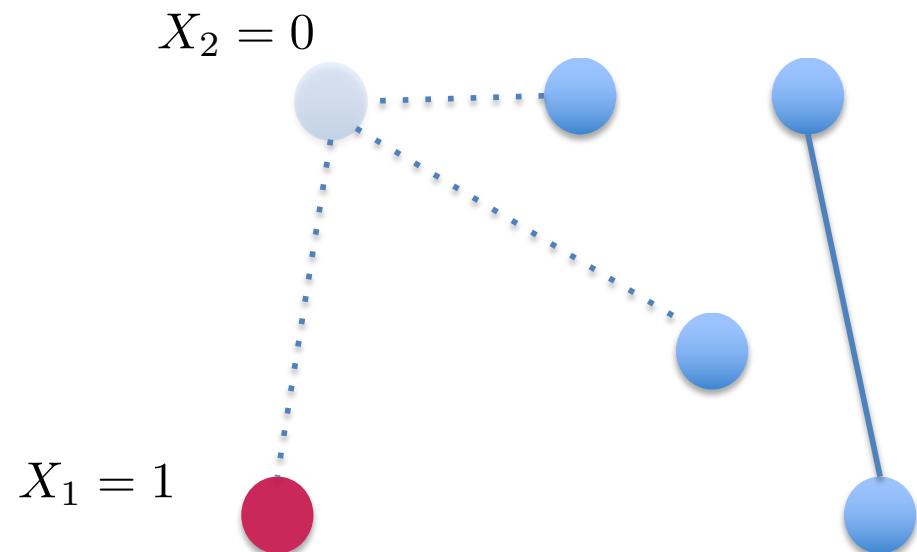
# DFS with random coins



Algorithm:

- Flip a coin for each vertex DFS wants to visit
- If 1 visit, if 0 don't visit and delete with its edges

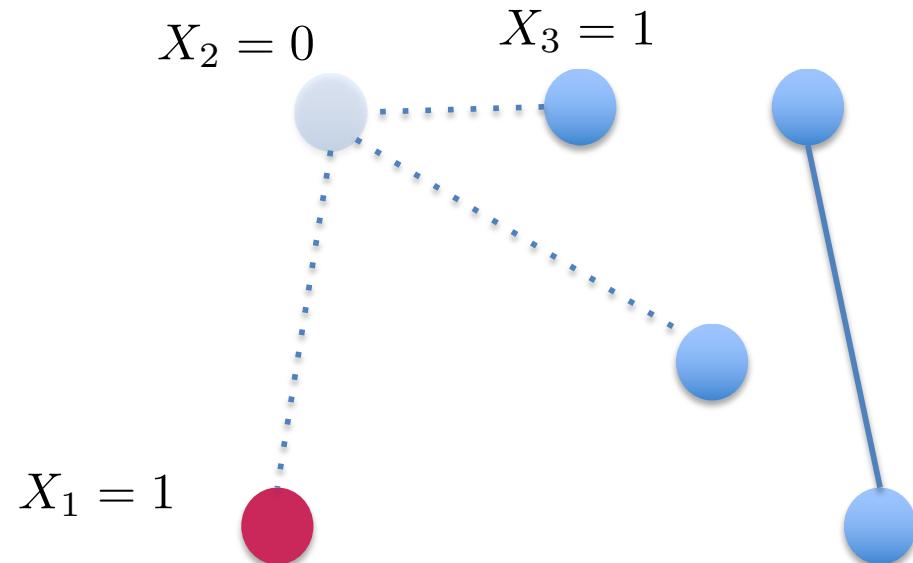
# DFS with random coins



Algorithm:

- Flip a coin for each vertex DFS wants to visit
- If 1 visit, if 0 don't visit and delete with its edges

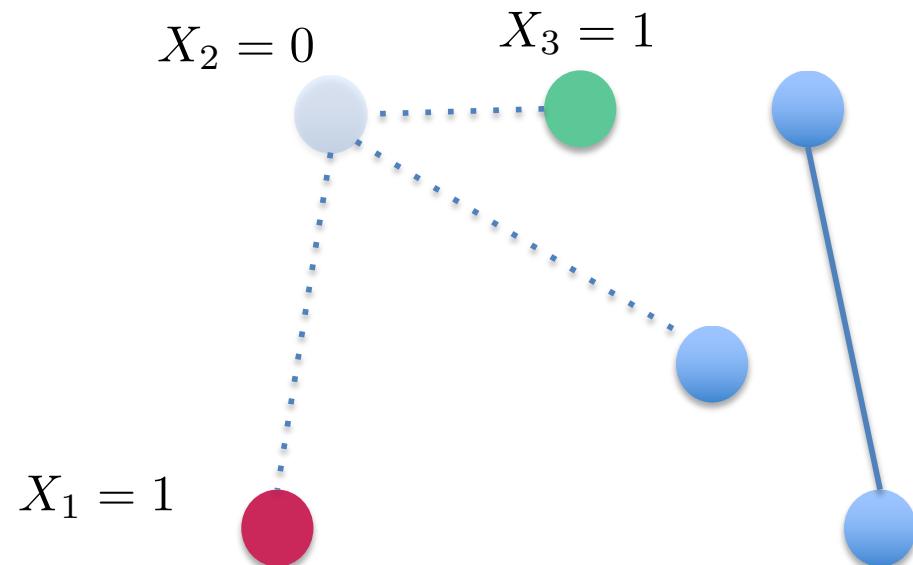
# DFS with random coins



Algorithm:

- Flip a coin for each vertex DFS wants to visit
- If 1 visit, if 0 don't visit and delete with its edges

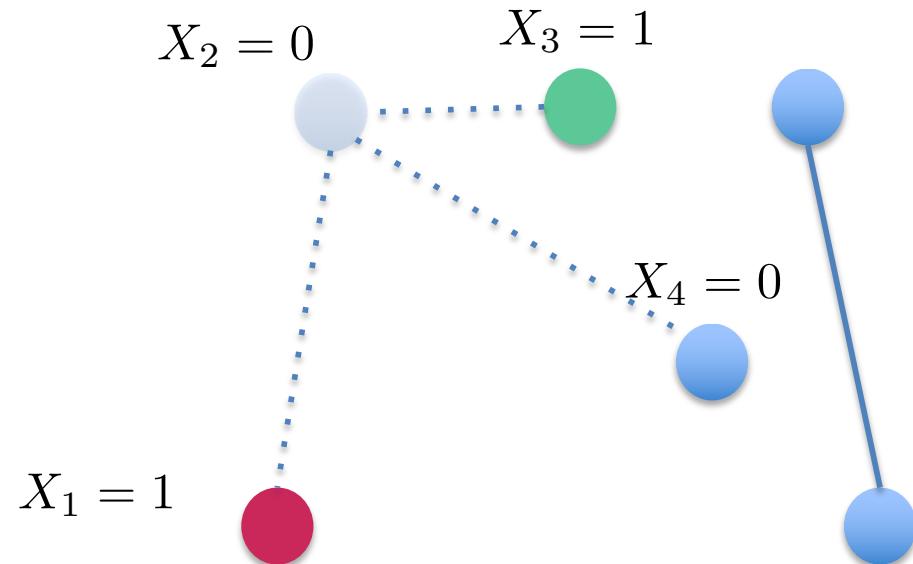
# DFS with random coins



Algorithm:

- Flip a coin for each vertex DFS wants to visit
- If 1 visit, if 0 don't visit and delete with its edges

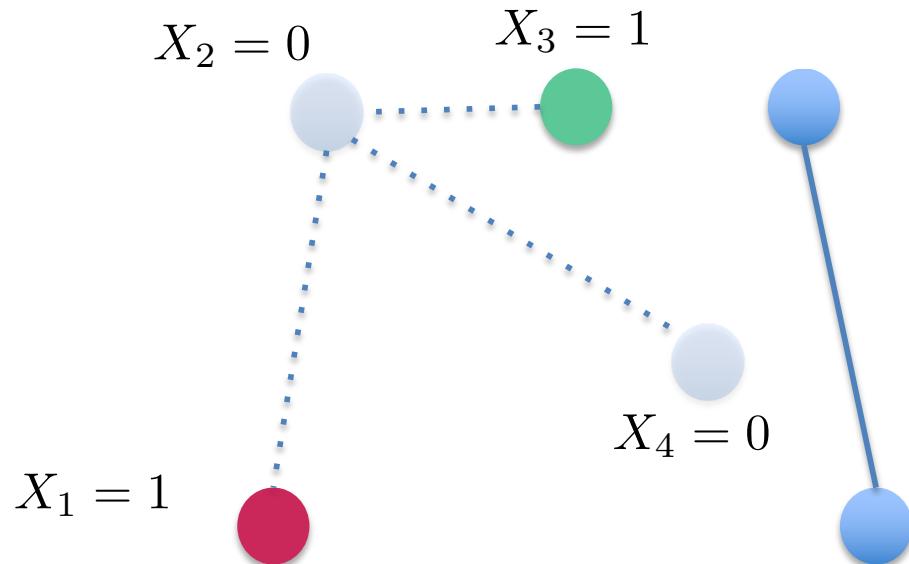
# DFS with random coins



Algorithm:

- Flip a coin for each vertex DFS wants to visit
- If 1 visit, if 0 don't visit and delete with its edges

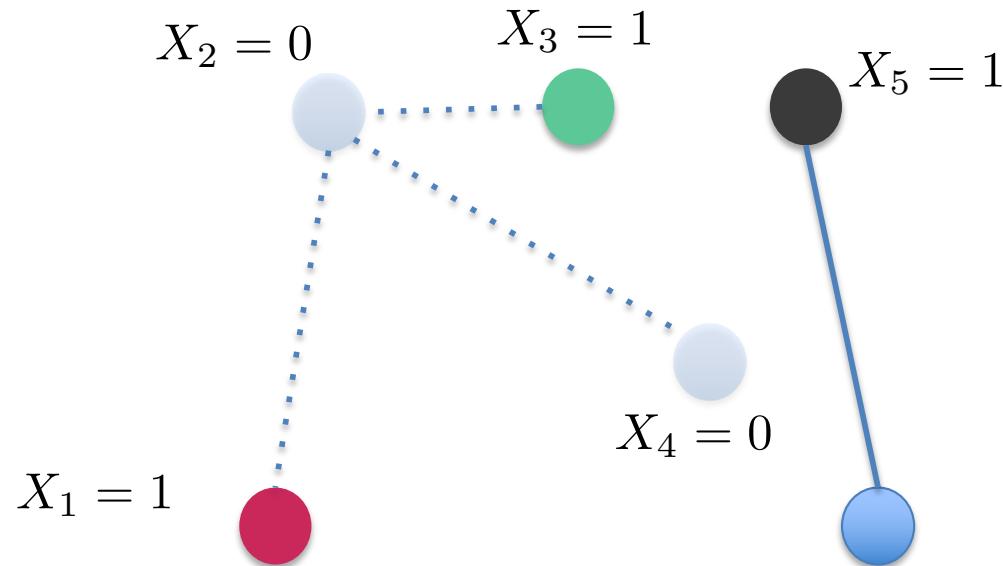
# DFS with random coins



Algorithm:

- Flip a coin for each vertex DFS wants to visit
- If 1 visit, if 0 don't visit and delete with its edges

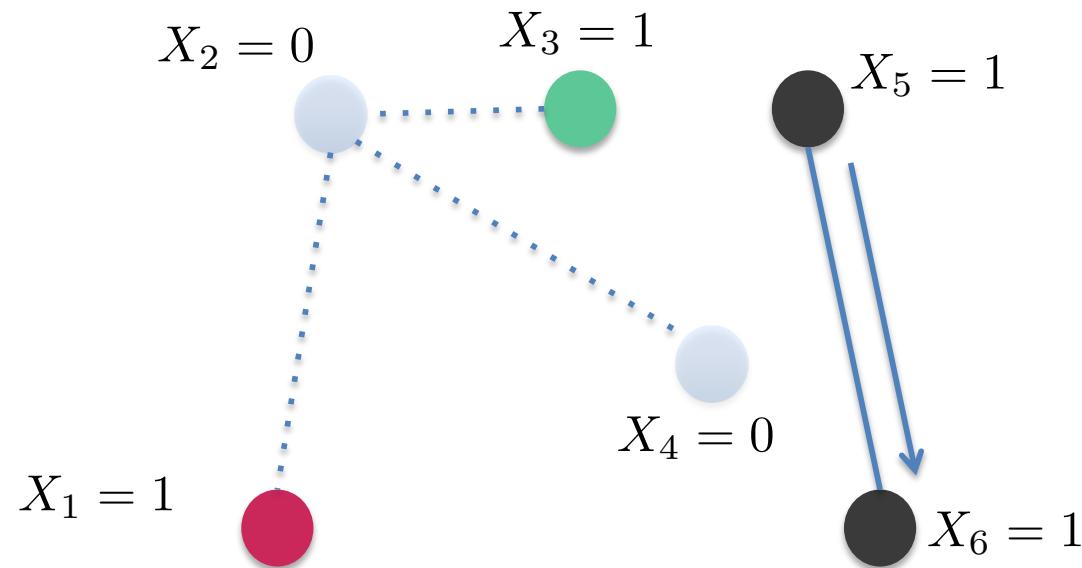
# DFS with random coins



Algorithm:

- Flip a coin for each vertex DFS wants to visit
- If 1 visit, if 0 don't visit and delete with its edges

# DFS with random coins

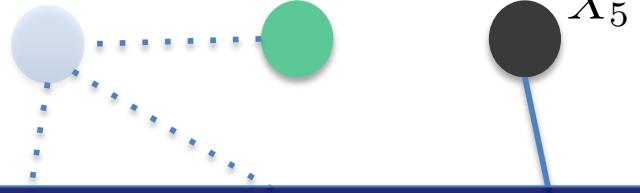


Algorithm:

- Flip a coin for each vertex DFS wants to visit
- If 1 visit, if 0 don't visit and delete with its edges

# DFS with random coins

$$X_2 = 0 \quad X_3 = 1 \quad X_5 = 1$$



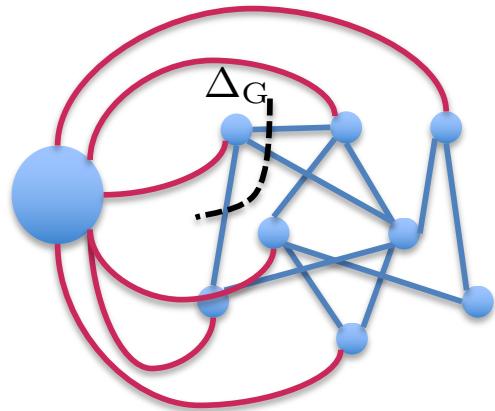
A Little extra trickery to turn this statement to a with or without replacement Theorem.

- Say I have a connected component of size  $k$
- #random coins flipped “associated” to that component  $\leq k * \Delta$
- Since I have a size  $k$  component it means that I had the event “at least  $k$  coins are “ON” in a set of  $k * \Delta$  coins”

$$(n - kd + 1)Pr[B(kd, p) \geq k] < n \cdot e^{-\frac{\epsilon^2}{3}(1-\epsilon)k} < n \cdot e^{-\frac{\epsilon^2(1-\epsilon)}{3} \frac{4}{\epsilon^2} \ln n} :$$

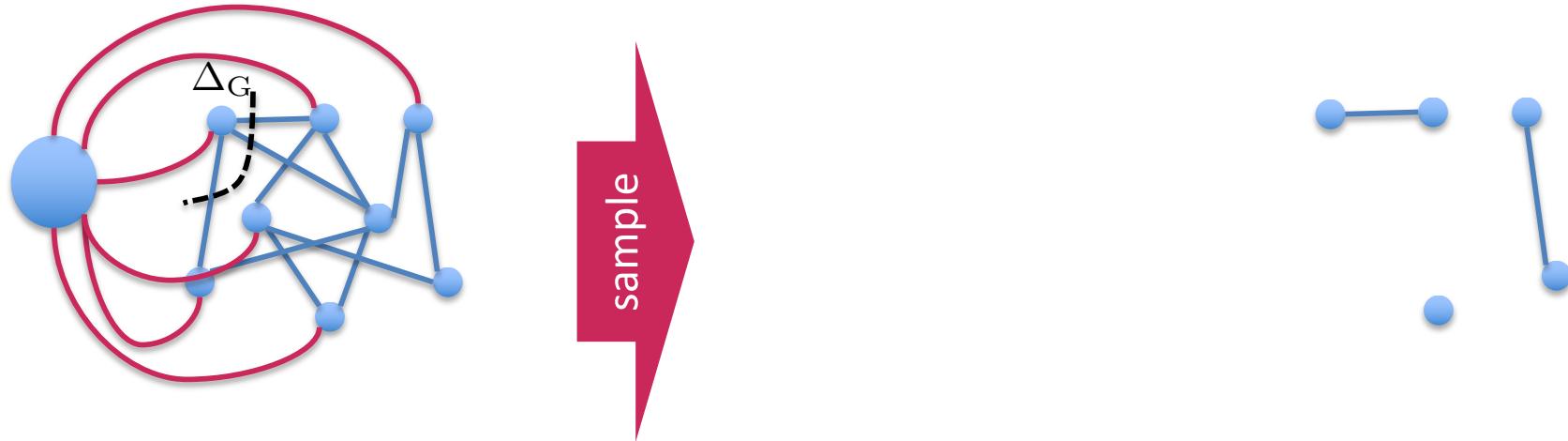
Is Max.Degree really an issue?

# High Degree Vertices (outliers)



Say we have a graph with a low-degree component  
+ some high degree vertices

# High Degree Vertices (outliers)



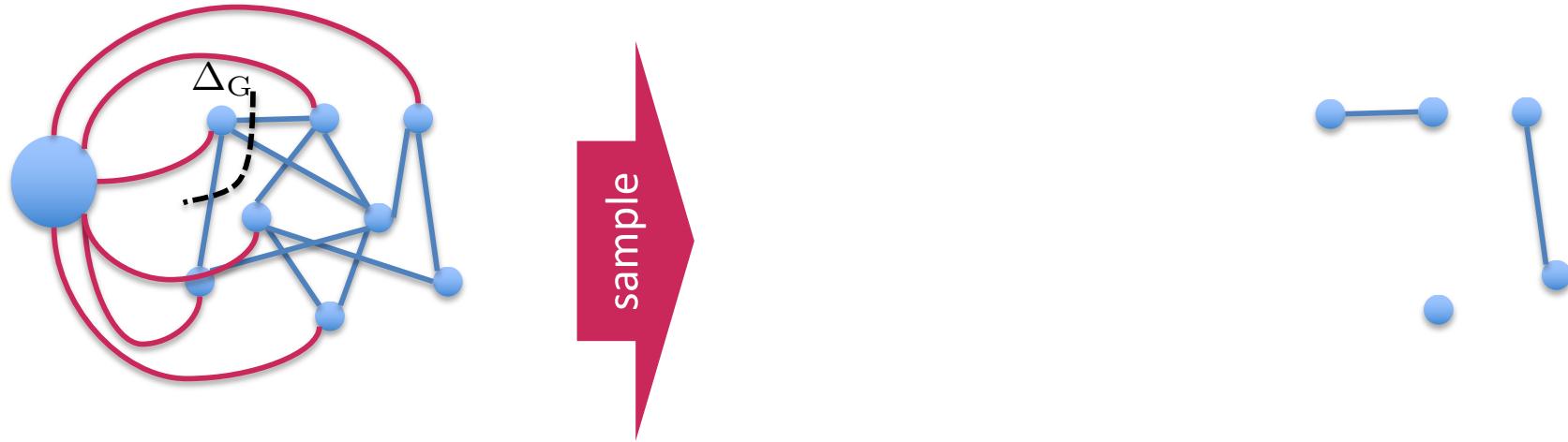
Lemma:

If you sample uniformly less than  $P \leq (1 - \epsilon) \frac{n}{\Delta_G}$  vertices  
Then, the induced subgraph of the low-degree (!) part  
shatters, and the largest connected component has size (whp)

$$\frac{4}{\epsilon^2} \cdot \log n$$

$\Delta_G$  is the max degree in the low-degree subgraph

# High Degree Vertices (outliers)



**Lemma 7.** Let us assume that there are  $O(n^\delta)$  outlier vertices in the original conflict graph  $G$  with degree at most  $\Delta_o$ , and let the remaining vertices have degree (induced on the remaining graph) at most  $\Delta$ . Let the induced update-variable graph on these low degree vertices abide to the same graph assumptions as those of Theorem 4. Moreover, let the batch size be bounded as

$$B \leq \min \left\{ (1 - \epsilon) \frac{n - O(n^\delta)}{\Delta}, \ O \left( \frac{n^{1-\delta}}{P} \right) \right\}.$$

Then, the expected runtime of CYCLADES will be  $O \left( \frac{E_u \cdot \kappa}{P} \cdot \log^2 n \right)$ .

# Experiments

# Experiments

Implementation in C++  
Experiments on Intel Xeon CPU E7-8870 v3  
1 TB RAM

	Dataset	# datapoints	# features	Density (average number of features per datapoint)
SAGA	NH2010	48,838	48,838	4.8026
SVRG	DBLP	5,425,964	5,425,964	3.1880
L2-SGD	MovieLens	~10M	82,250	200
SGD	EN-Wiki	20,207,156	213,272	200

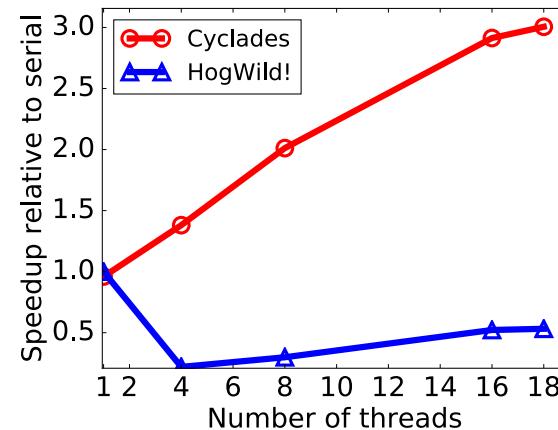
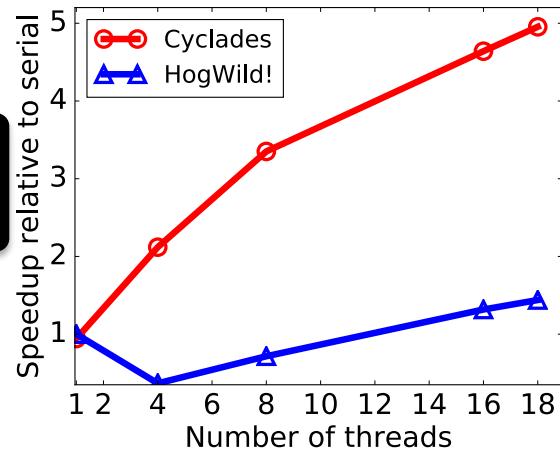
Full asynchronous (Hogwild!)

vs

CYCLADES

# Speedups

SVRG/SAGA

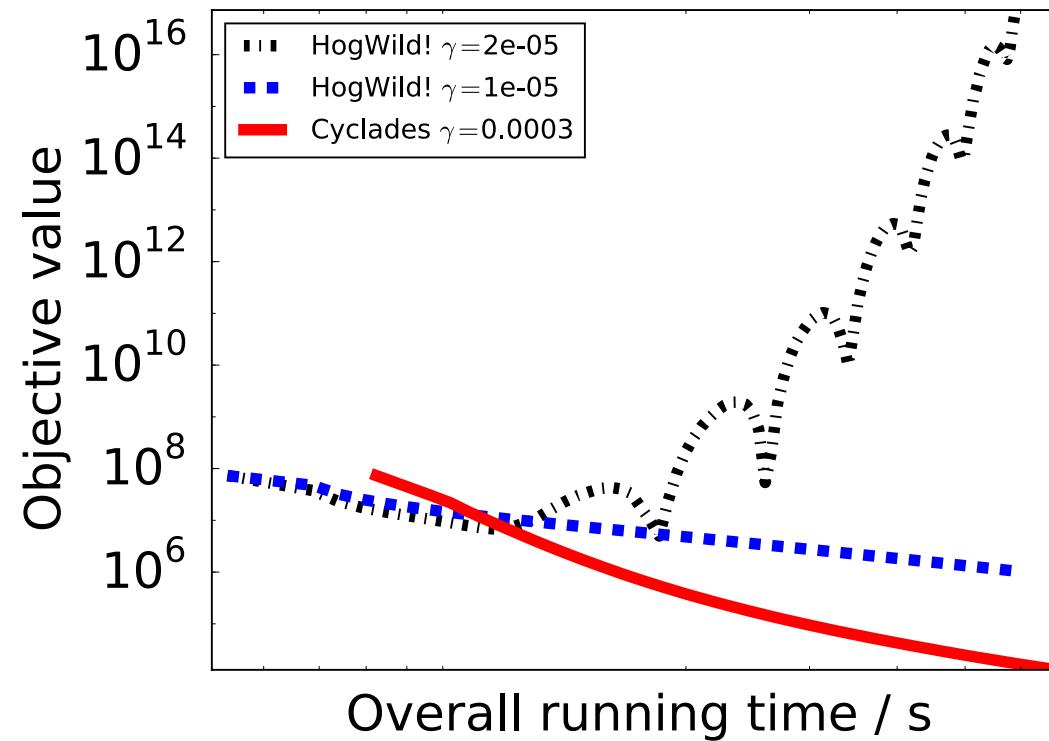


~ 500% gain

(a) Least squares, DBLP, SAGA (b) Graph Eig., NH2010, SVRG

# Convergence

Least Squares SAGA  
16 threads



# Open Problems

Assumptions:

Sparsity is Key

O.P.:

Can we handle Dense Data?

O.P.:

Data sparsification  
for  $f(\langle a, x \rangle)$  problems?

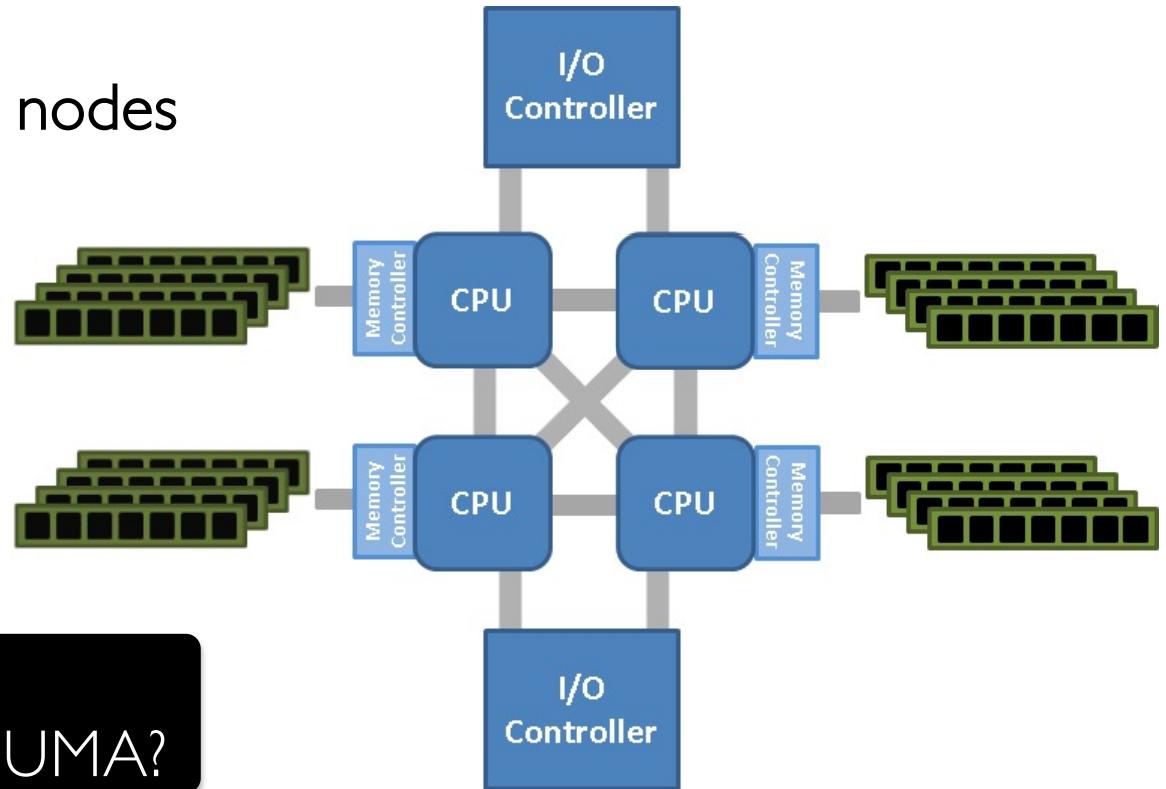
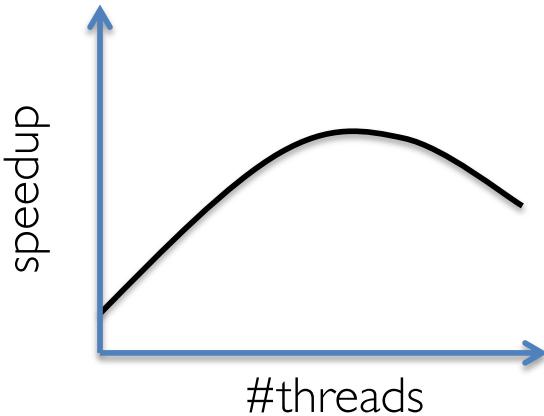
maybe...

We can relax serial equivalence to an “expected” one?

# Open Problems

Asynchronous algorithms great for Shared Memory Systems

- Issues when scaling across nodes



O.P.:

How to provably scale on NUMA?

- Similar Issues for Distributed:

O.P.:

What is the right ML Paradigm  
for Distributed?

# CYCLADES

a framework for Parallel Sparse ML algorithms

- Lock-free + (maximally) Asynchronous
- No Conflicts
- Serializable
- Black-box analysis



# Next Time

- Communication Bottlenecks
- Compressed Gradients
- Quantization

# Reading List

- Krivelevich, M., 2014. The phase transition in site percolation on pseudo-random graphs. arXiv preprint arXiv:1404.5731.
- Pan, X., Lam, M., Tu, S., Papailiopoulos, D., Zhang, C., Jordan, M.I., Ramchandran, K. and Ré, C., 2016. Cyclades: Conflict-free asynchronous machine learning. Advances in Neural Information Processing Systems, 29.
- Pan, X., Papailiopoulos, D., Oymak, S., Recht, B., Ramchandran, K. and Jordan, M.I., 2015. Parallel correlation clustering on big graphs. Advances in Neural Information Processing Systems, 28.