

Big Data Infrastructure

Session 5: MapReduce – Graphs

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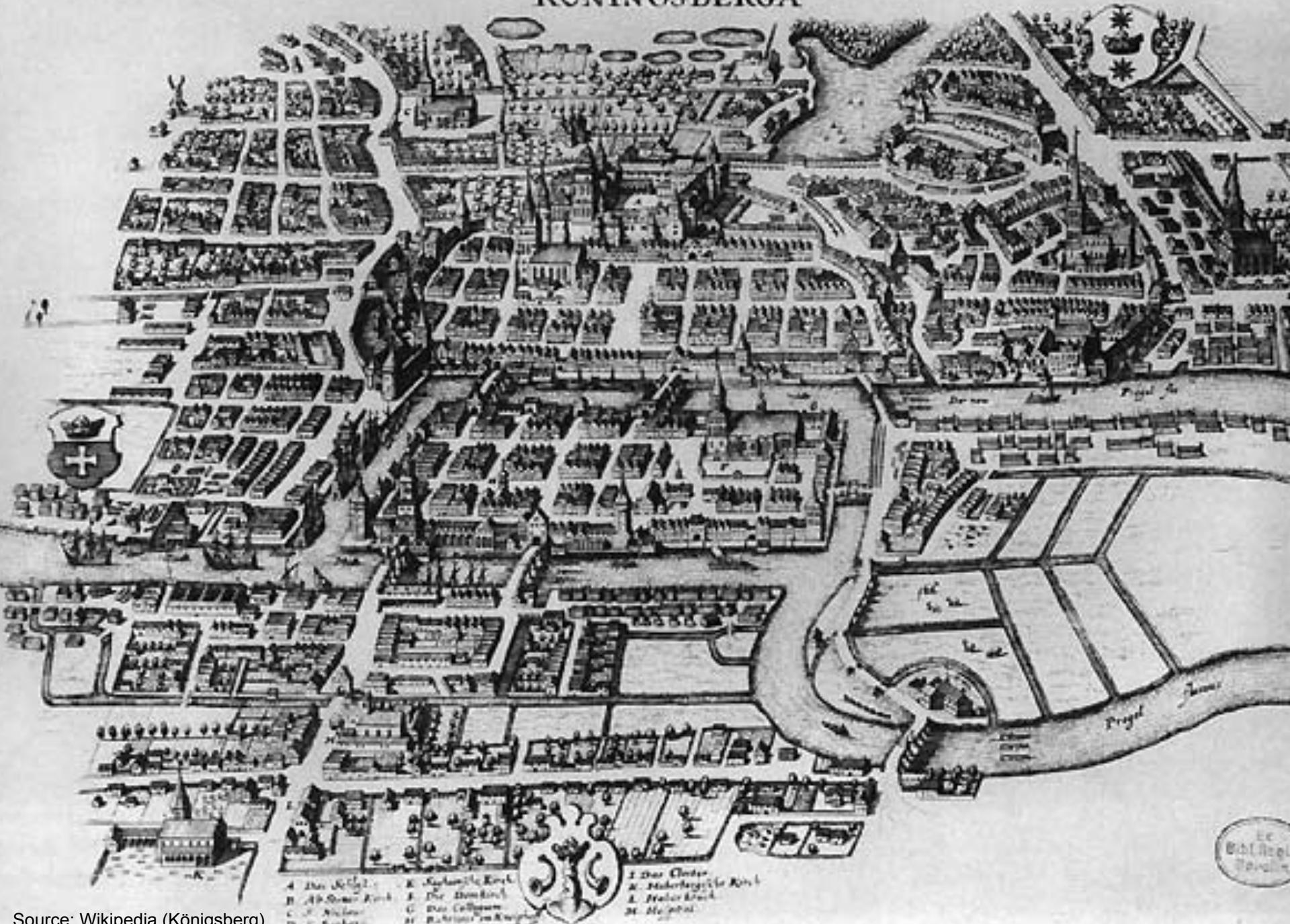
Today's Agenda

- Graph problems and representations
- Parallel breadth-first search
- PageRank
- Beyond PageRank and other graph algorithms
- Optimizing graph algorithms

What's a graph?

- $G = (V, E)$, where
 - V represents the set of vertices (nodes)
 - E represents the set of edges (links)
 - Both vertices and edges may contain additional information
- Different types of graphs:
 - Directed vs. undirected edges
 - Presence or absence of cycles
- Graphs are everywhere:
 - Hyperlink structure of the web
 - Physical structure of computers on the Internet
 - Interstate highway system
 - Social networks

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Some Graph Problems

- Finding shortest paths
 - Routing Internet traffic and UPS trucks
- Finding minimum spanning trees
 - Telco laying down fiber
- Finding Max Flow
 - Airline scheduling
- Identify “special” nodes and communities
 - Breaking up terrorist cells, spread of avian flu
- Bipartite matching
 - Monster.com, Match.com
- And of course... PageRank

Graphs and MapReduce

- A large class of graph algorithms involve:
 - Performing computations at each node: based on node features, edge features, and local link structure
 - Propagating computations: “traversing” the graph
- Key questions:
 - How do you represent graph data in MapReduce?
 - How do you traverse a graph in MapReduce?

Representing Graphs

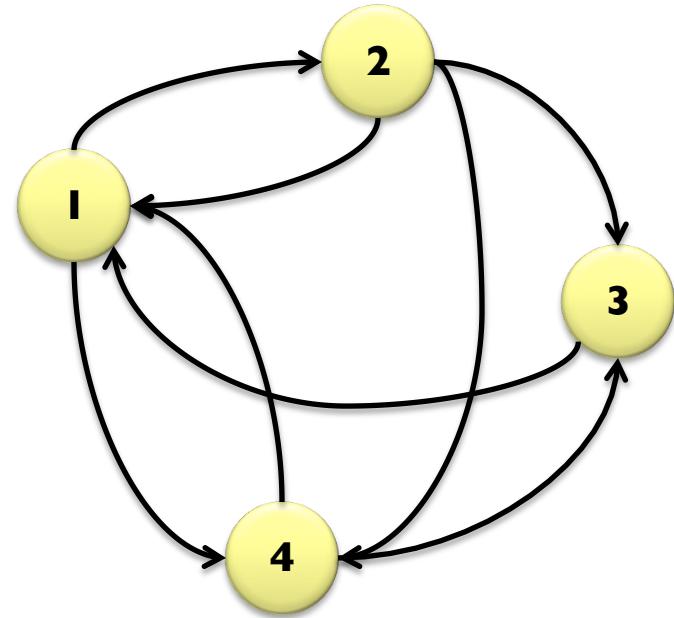
- $G = (V, E)$
- Two common representations
 - Adjacency matrix
 - Adjacency list

Adjacency Matrices

Represent a graph as an $n \times n$ square matrix M

- $n = |V|$
- $M_{ij} = 1$ means a link from node i to j

	1	2	3	4
1	0	1	0	1
2	1	0	1	1
3	1	0	0	0
4	1	0	1	0



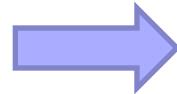
Adjacency Matrices: Critique

- Advantages:
 - Amenable to mathematical manipulation
 - Iteration over rows and columns corresponds to computations on outlinks and inlinks
- Disadvantages:
 - Lots of zeros for sparse matrices
 - Lots of wasted space

Adjacency Lists

Take adjacency matrices... and throw away all the zeros

	1	2	3	4
1	0	1	0	1
2	1	0	1	1
3	1	0	0	0
4	1	0	1	0



- 1: 2, 4
- 2: 1, 3, 4
- 3: 1
- 4: 1, 3

Adjacency Lists: Critique

- Advantages:

- Much more compact representation
- Easy to compute over outlinks

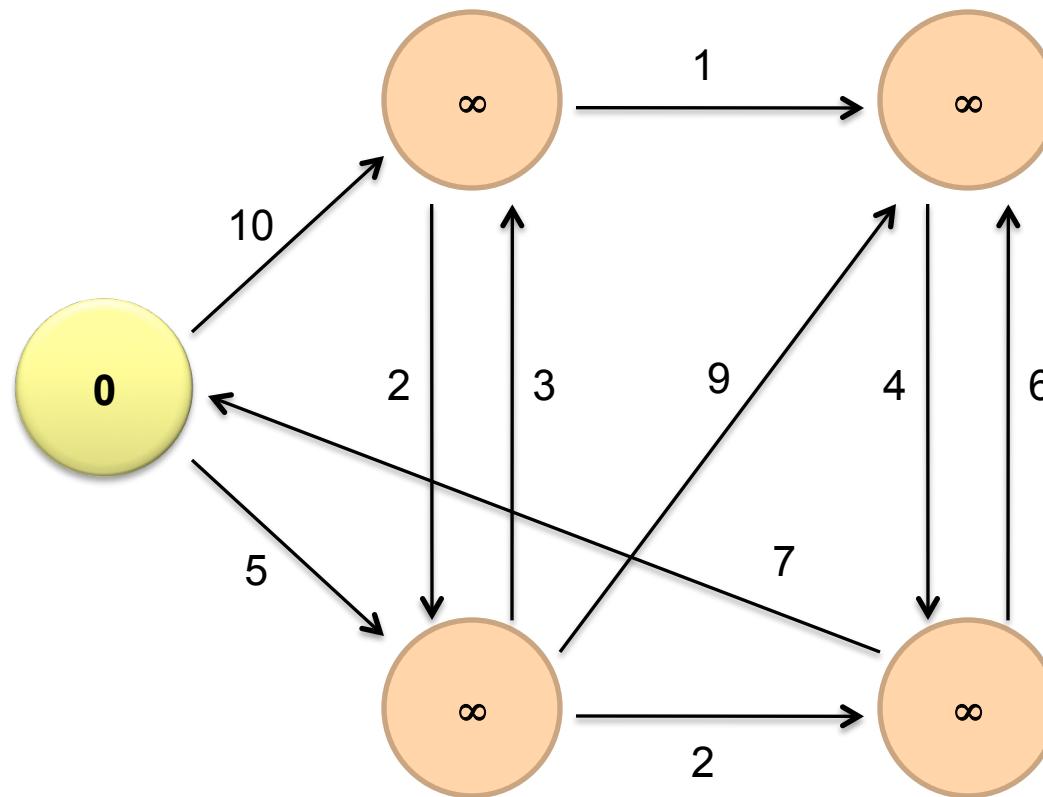
- Disadvantages:

- Much more difficult to compute over inlinks

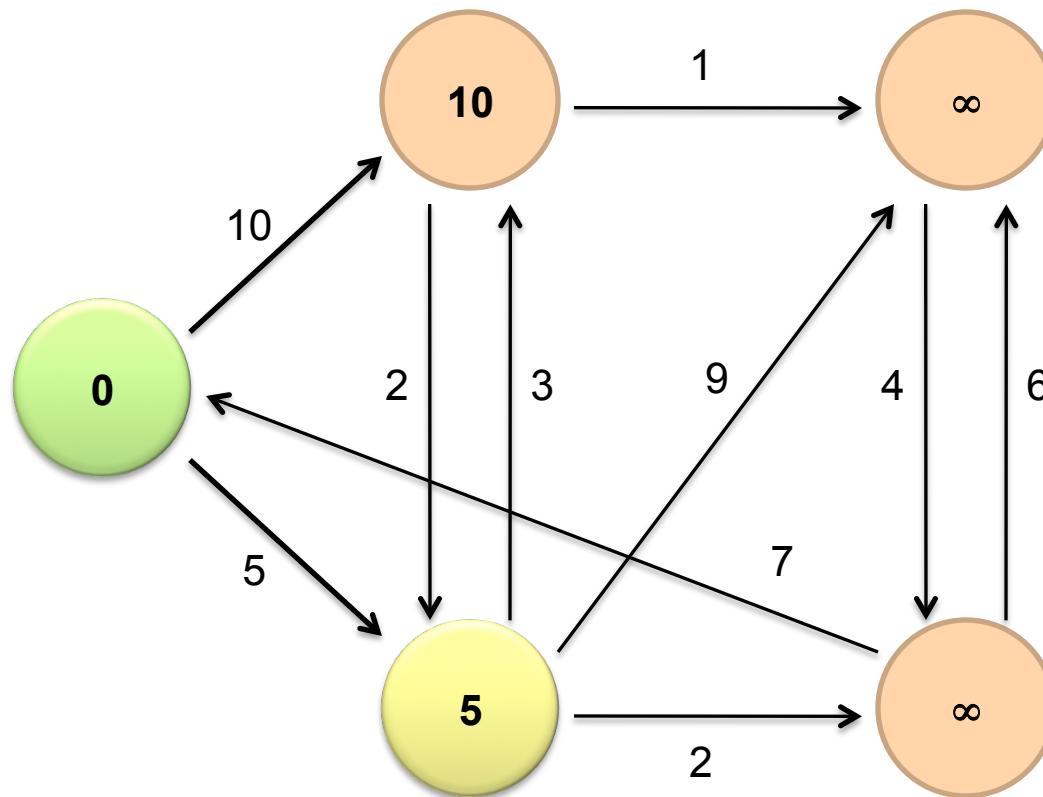
Single-Source Shortest Path

- **Problem:** find shortest path from a source node to one or more target nodes
 - Shortest might also mean lowest weight or cost
- First, a refresher: Dijkstra's Algorithm

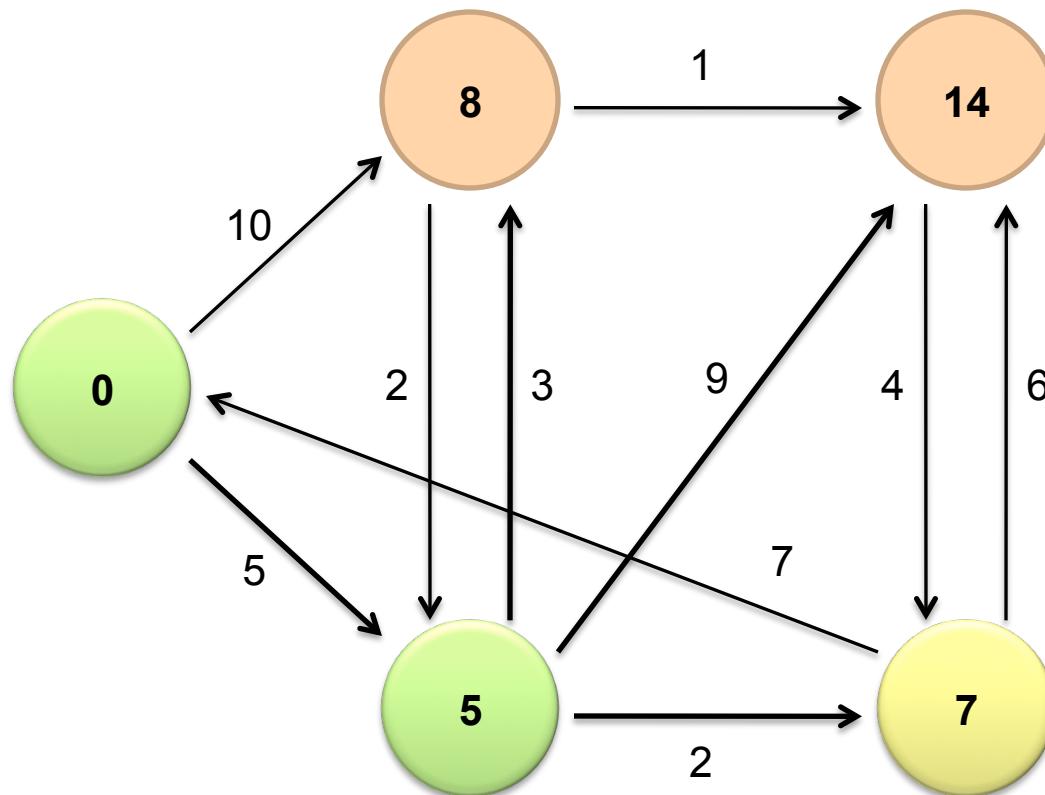
Dijkstra's Algorithm Example



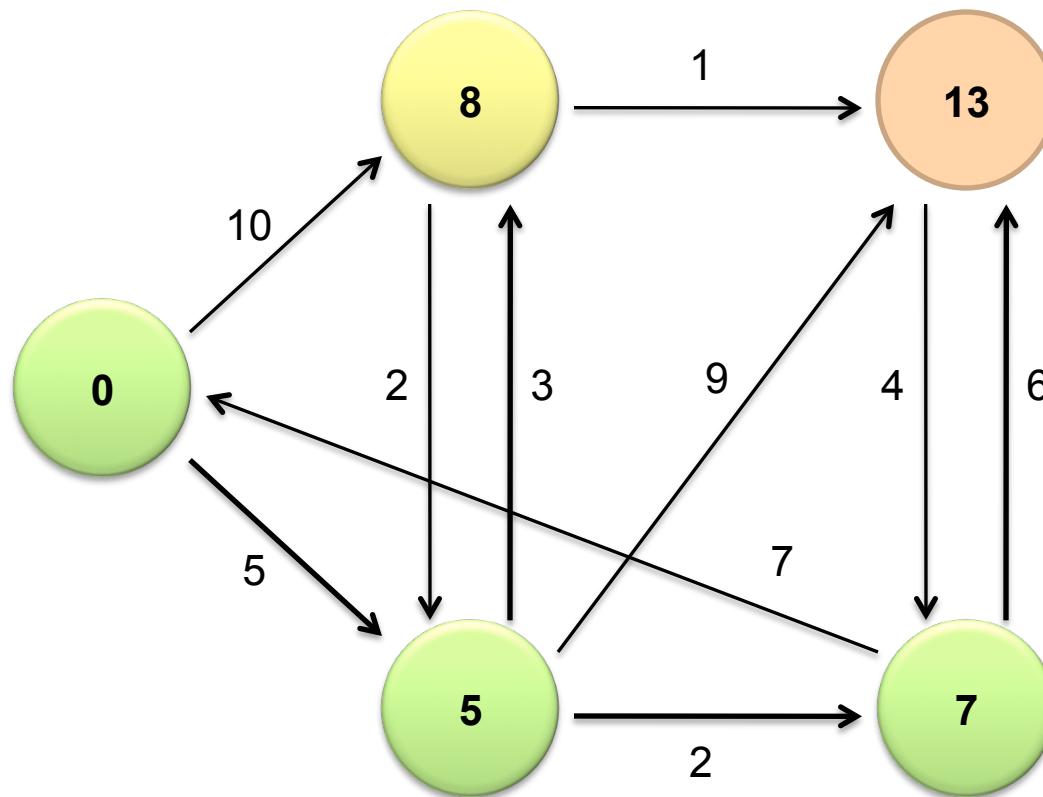
Dijkstra's Algorithm Example



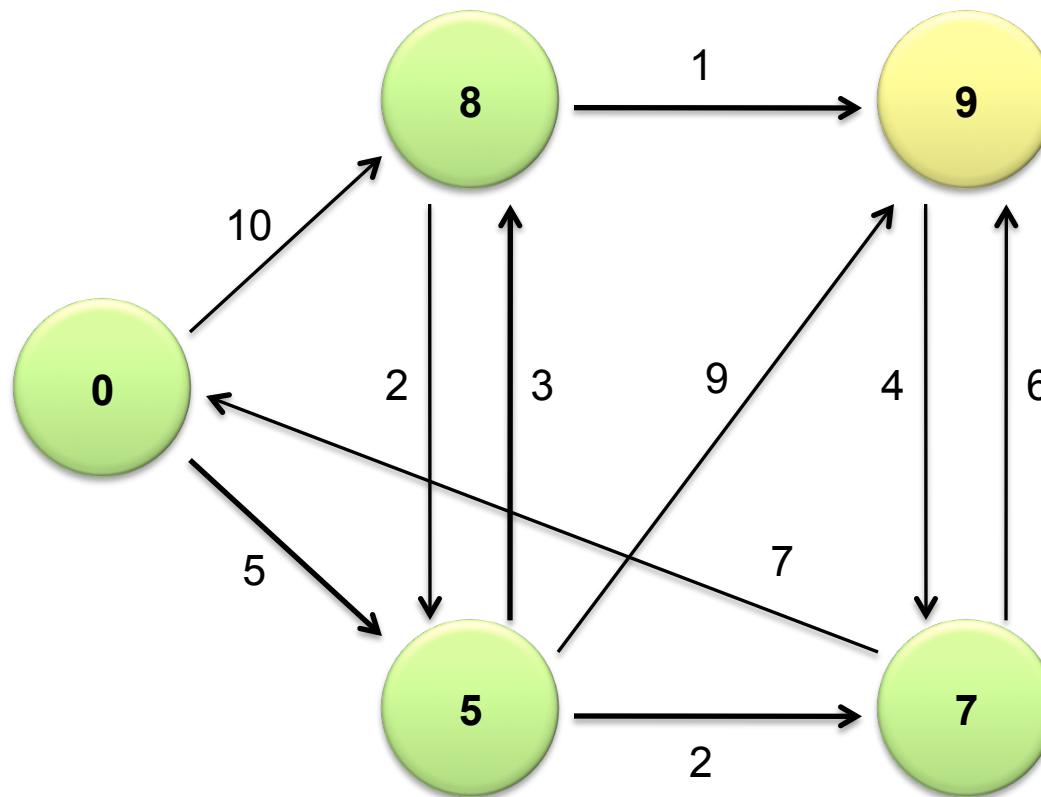
Dijkstra's Algorithm Example



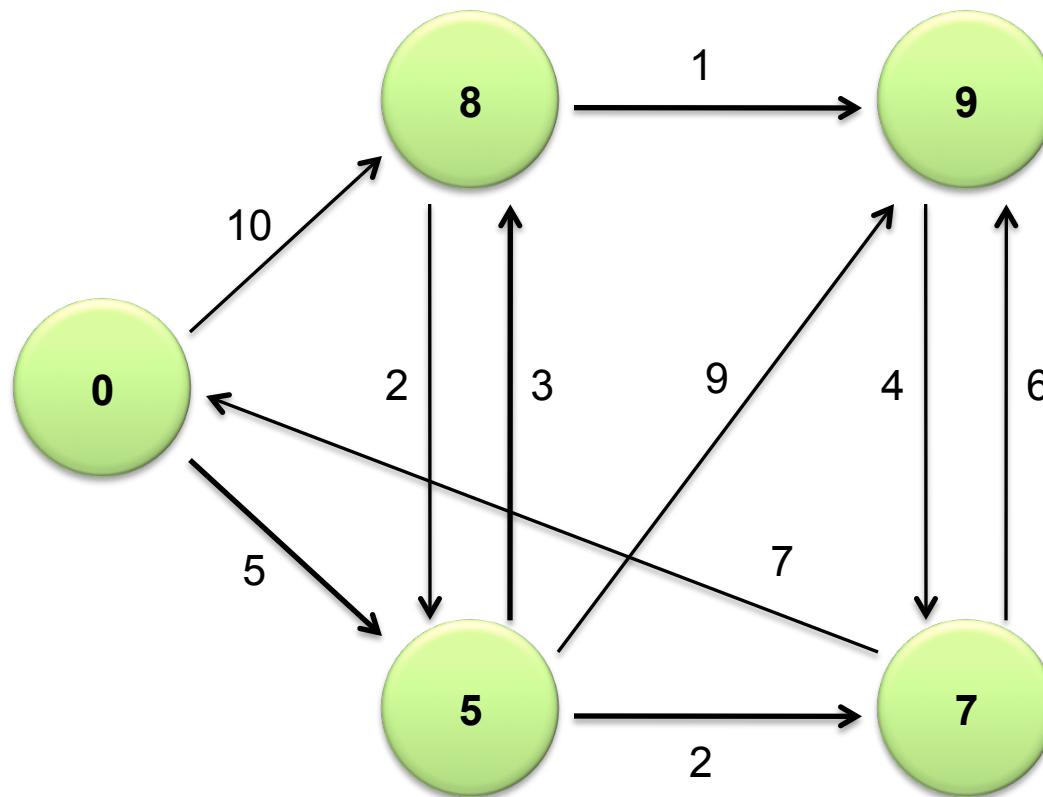
Dijkstra's Algorithm Example



Dijkstra's Algorithm Example



Dijkstra's Algorithm Example

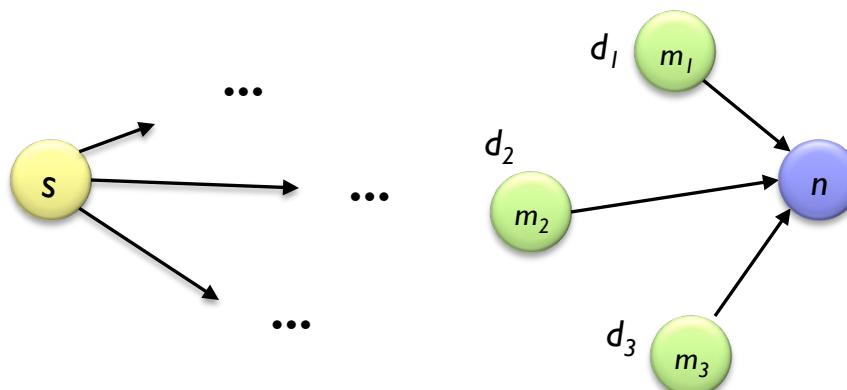


Single-Source Shortest Path

- **Problem:** find shortest path from a source node to one or more target nodes
 - Shortest might also mean lowest weight or cost
- Single processor machine: Dijkstra's Algorithm
- MapReduce: parallel breadth-first search (BFS)

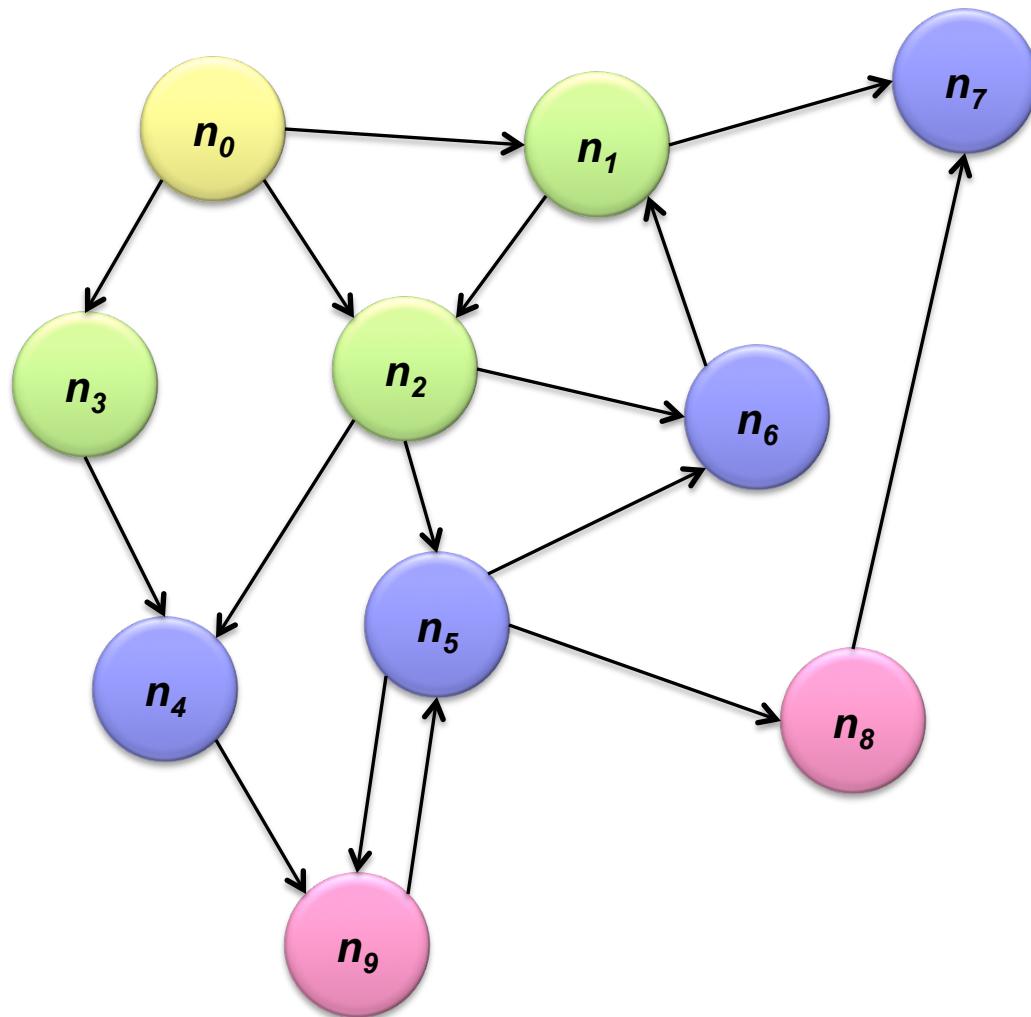
Finding the Shortest Path

- Consider simple case of equal edge weights
- Solution to the problem can be defined inductively
- Here's the intuition:
 - Define: b is reachable from a if b is on adjacency list of a
 $\text{DISTANCETo}(s) = 0$
 - For all nodes p reachable from s ,
 $\text{DISTANCETo}(p) = 1$
 - For all nodes n reachable from some other set of nodes M ,
 $\text{DISTANCETo}(n) = 1 + \min(\text{DISTANCETo}(m), m \in M)$





Visualizing Parallel BFS



From Intuition to Algorithm

- Data representation:
 - Key: node n
 - Value: d (distance from start), adjacency list (nodes reachable from n)
 - Initialization: for all nodes except for start node, $d = \infty$
- Mapper:
 - $\forall m \in$ adjacency list: emit $(m, d + 1)$
 - Remember to also emit distance to yourself
- Sort/Shuffle
 - Groups distances by reachable nodes
- Reducer:
 - Selects minimum distance path for each reachable node
 - Additional bookkeeping needed to keep track of actual path

Multiple Iterations Needed

- Each MapReduce iteration advances the “frontier” by one hop
 - Subsequent iterations include more and more reachable nodes as frontier expands
 - Multiple iterations are needed to explore entire graph
- Preserving graph structure:
 - Problem: Where did the adjacency list go?
 - Solution: mapper emits $(n, \text{adjacency list})$ as well

BFS Pseudo-Code

```
1: class MAPPER
2:   method MAP(nid  $n$ , node  $N$ )
3:      $d \leftarrow N.\text{DISTANCE}$ 
4:     EMIT(nid  $n, N$ )                                 $\triangleright$  Pass along graph structure
5:     for all nodeid  $m \in N.\text{ADJACENCYLIST}$  do
6:       EMIT(nid  $m, d + 1$ )                           $\triangleright$  Emit distances to reachable nodes

1: class REDUCER
2:   method REDUCE(nid  $m, [d_1, d_2, \dots]$ )
3:      $d_{min} \leftarrow \infty$ 
4:      $M \leftarrow \emptyset$ 
5:     for all  $d \in \text{counts } [d_1, d_2, \dots]$  do
6:       if IsNODE( $d$ ) then
7:          $M \leftarrow d$                                  $\triangleright$  Recover graph structure
8:       else if  $d < d_{min}$  then
9:          $d_{min} \leftarrow d$                           $\triangleright$  Look for shorter distance
10:         $M.\text{DISTANCE} \leftarrow d_{min}$              $\triangleright$  Update shortest distance
11:        EMIT(nid  $m, \text{node } M$ )
```

Stopping Criterion

- How many iterations are needed in parallel BFS (equal edge weight case)?
- Convince yourself: when a node is first “discovered”, we’ve found the shortest path
- Now answer the question...
 - Six degrees of separation?
- Practicalities of implementation in MapReduce

Comparison to Dijkstra

- Dijkstra's algorithm is more efficient
 - At each step, only pursues edges from minimum-cost path inside frontier
- MapReduce explores all paths in parallel
 - Lots of “waste”
 - Useful work is only done at the “frontier”
- Why can't we do better using MapReduce?

Single Source: Weighted Edges

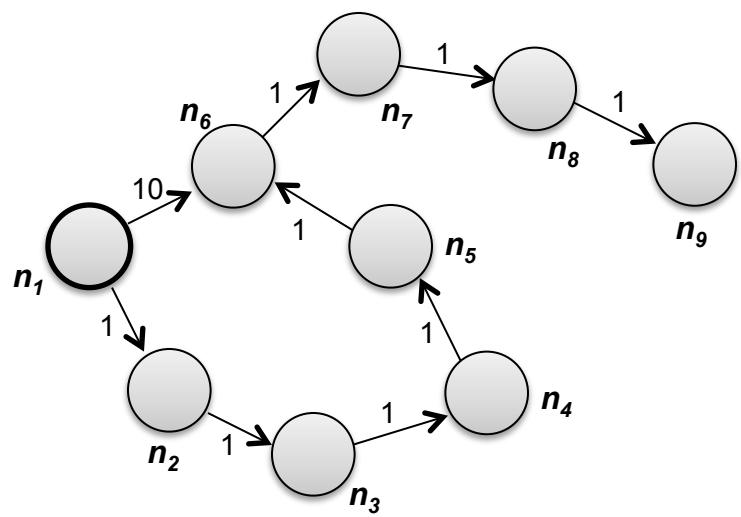
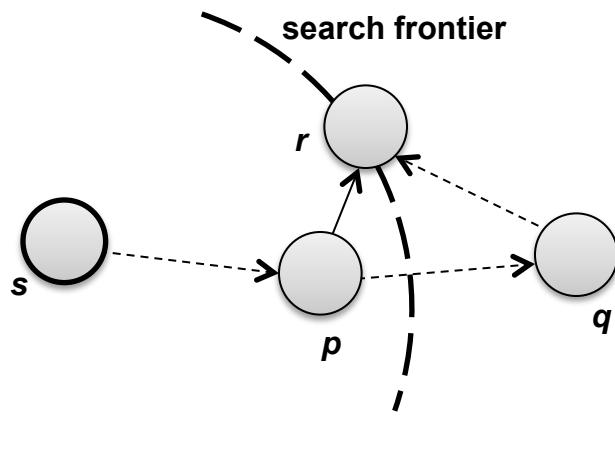
- Now add positive weights to the edges
 - Why can't edge weights be negative?
- Simple change: add weight w for each edge in adjacency list
 - In mapper, emit $(m, d + w_p)$ instead of $(m, d + 1)$ for each node m
- That's it?

Stopping Criterion

- How many iterations are needed in parallel BFS (positive edge weight case)?
- Convince yourself: when a node is first “discovered”, we’ve found the shortest path

Not true!

Additional Complexities



Stopping Criterion

- How many iterations are needed in parallel BFS (positive edge weight case)?
- Practicalities of implementation in MapReduce

Application: Social Search



Social Search

- When searching, how to rank friends named “John”?
 - Assume undirected graphs
 - Rank matches by distance to user
- Naïve implementations:
 - Precompute all-pairs distances
 - Compute distances at query time
- Can we do better?

All-Pairs?

- Floyd-Warshall Algorithm: difficult to MapReduce-ify...
- Multiple-source shortest paths in MapReduce: run multiple parallel BFS *simultaneously*
 - Assume source nodes $\{s_0, s_1, \dots s_n\}$
 - Instead of emitting a single distance, emit an array of distances, with respect to each source
 - Reducer selects minimum for each element in array
- Does this scale?

Landmark Approach (aka sketches)

- Select n seeds $\{s_0, s_1, \dots s_n\}$
- Compute distances from seeds to every node:

$$A = [2, 1, 1]$$

$$B = [1, 1, 2]$$

$$C = [4, 3, 1]$$

$$D = [1, 2, 4]$$

- What can we conclude about distances?
- Insight: landmarks bound the maximum path length
- Lots of details:
 - How to more tightly bound distances
 - How to select landmarks (random isn't the best...)
- Use multi-source parallel BFS implementation in MapReduce!



<pause/>

Graphs and MapReduce

- A large class of graph algorithms involve:
 - Performing computations at each node: based on node features, edge features, and local link structure
 - Propagating computations: “traversing” the graph
- Generic recipe:
 - Represent graphs as adjacency lists
 - Perform local computations in mapper
 - Pass along partial results via outlinks, keyed by destination node
 - Perform aggregation in reducer on inlinks to a node
 - Iterate until convergence: controlled by external “driver”
 - Don’t forget to pass the graph structure between iterations

Random Walks Over the Web

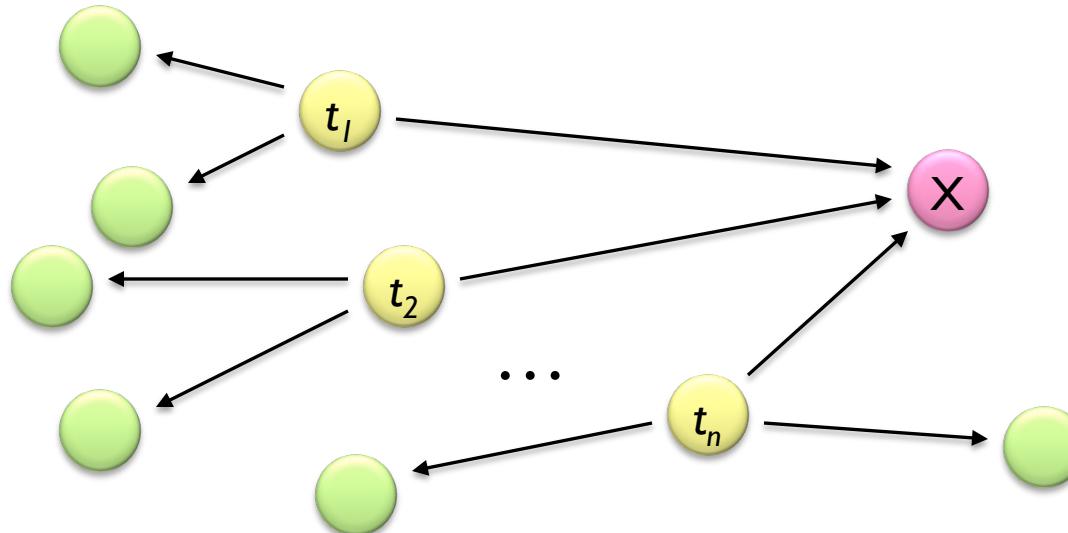
- Random surfer model:
 - User starts at a random Web page
 - User randomly clicks on links, surfing from page to page
- PageRank
 - Characterizes the amount of time spent on any given page
 - Mathematically, a probability distribution over pages
- PageRank captures notions of page importance
 - Correspondence to human intuition?
 - One of thousands of features used in web search (query-independent)

PageRank: Defined

Given page x with inlinks t_1, \dots, t_n , where

- $C(t)$ is the out-degree of t
- α is probability of random jump
- N is the total number of nodes in the graph

$$PR(x) = \alpha \left(\frac{1}{N} \right) + (1 - \alpha) \sum_{i=1}^n \frac{PR(t_i)}{C(t_i)}$$



Computing PageRank

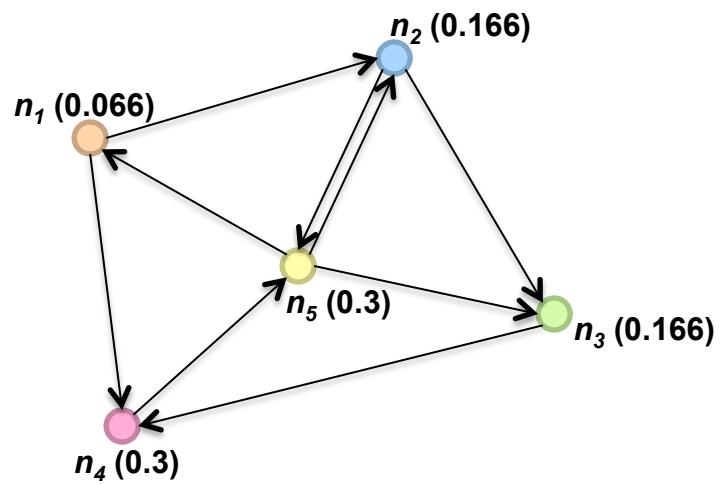
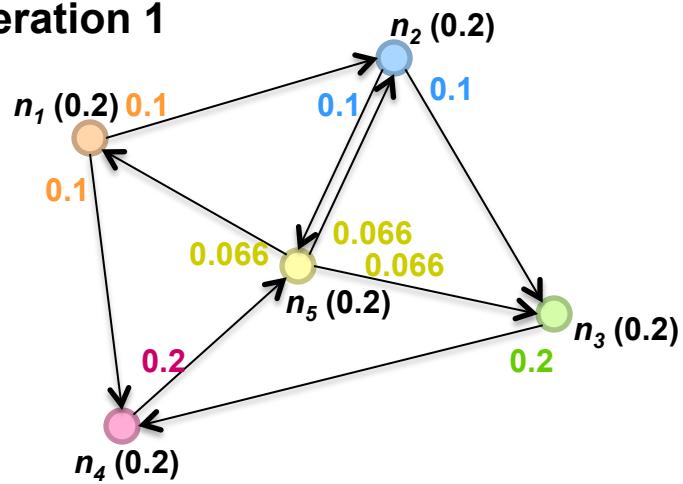
- Properties of PageRank
 - Can be computed iteratively
 - Effects at each iteration are local
- Sketch of algorithm:
 - Start with seed PR_i values
 - Each page distributes PR_i “credit” to all pages it links to
 - Each target page adds up “credit” from multiple in-bound links to compute PR_{i+1}
 - Iterate until values converge

Simplified PageRank

- First, tackle the simple case:
 - No random jump factor
 - No dangling nodes
- Then, factor in these complexities...
 - Why do we need the random jump?
 - Where do dangling nodes come from?

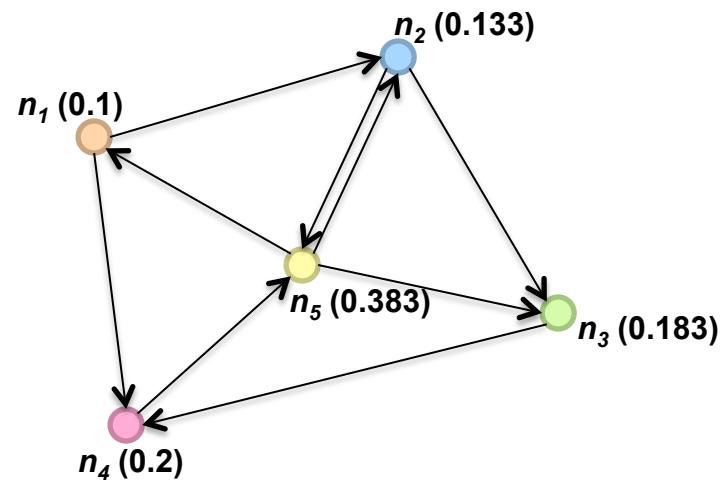
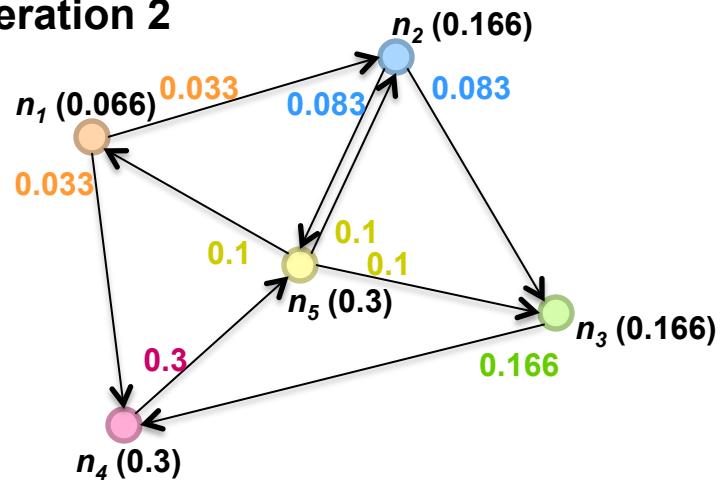
Sample PageRank Iteration (I)

Iteration 1

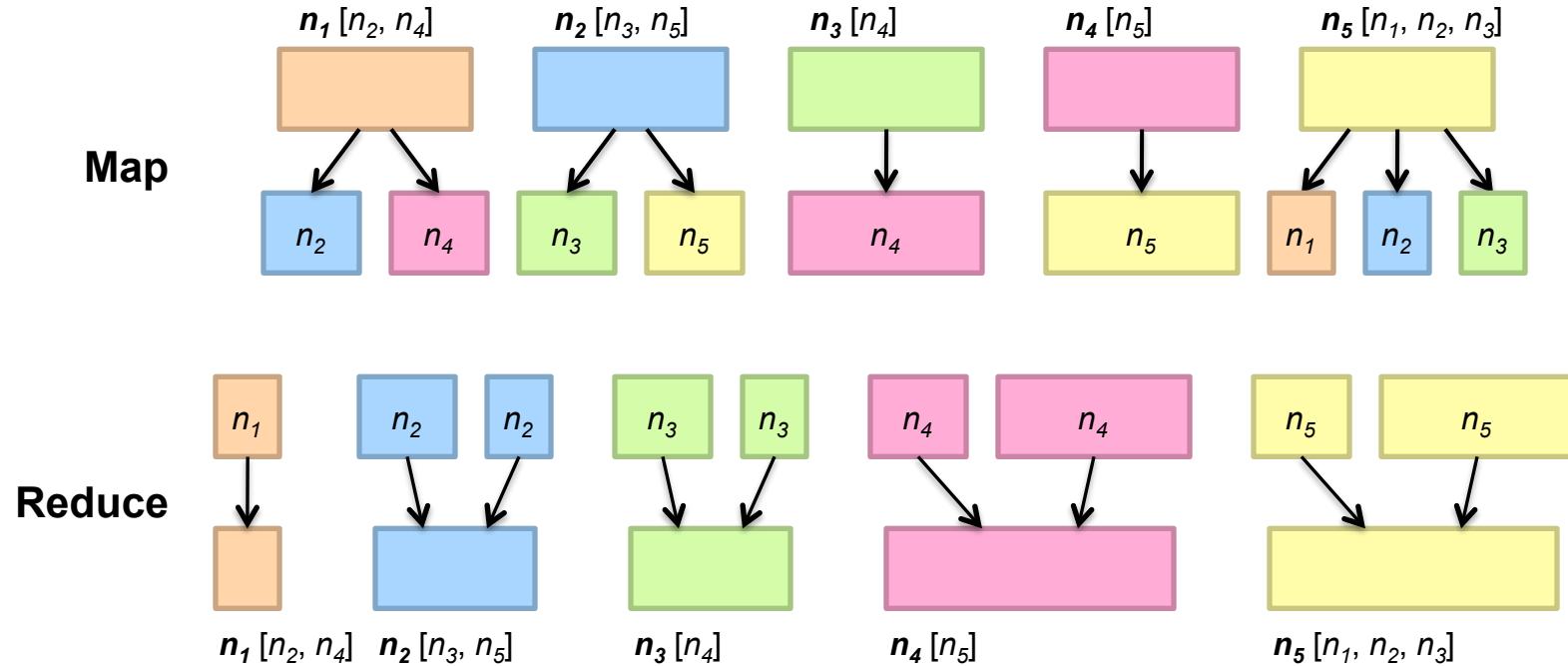


Sample PageRank Iteration (2)

Iteration 2



PageRank in MapReduce



PageRank Pseudo-Code

```
1: class MAPPER
2:   method MAP(nid  $n$ , node  $N$ )
3:      $p \leftarrow N.\text{PAGERANK}/|N.\text{ADJACENCYLIST}|$ 
4:     EMIT(nid  $n, N$ )                                ▷ Pass along graph structure
5:     for all nodeid  $m \in N.\text{ADJACENCYLIST}$  do
6:       EMIT(nid  $m, p$ )                            ▷ Pass PageRank mass to neighbors
7:
8: class REDUCER
9:   method REDUCE(nid  $m, [p_1, p_2, \dots]$ )
10:     $M \leftarrow \emptyset$ 
11:    for all  $p \in \text{counts} [p_1, p_2, \dots]$  do
12:      if IsNODE( $p$ ) then
13:         $M \leftarrow p$                                 ▷ Recover graph structure
14:      else
15:         $s \leftarrow s + p$                           ▷ Sums incoming PageRank contributions
16:     $M.\text{PAGERANK} \leftarrow s$ 
17:    EMIT(nid  $m, \text{node } M$ )
```

PageRank vs. BFS

PageRank

BFS

Map

PR/N

$d + I$

Reduce

sum

min

Complete PageRank

- Two additional complexities
 - What is the proper treatment of dangling nodes?
 - How do we factor in the random jump factor?
- Solution:
 - Second pass to redistribute “missing PageRank mass” and account for random jumps

$$p' = \alpha \left(\frac{1}{N} \right) + (1 - \alpha) \left(\frac{m}{N} + p \right)$$

- p is PageRank value from before, p' is updated PageRank value
 - N is the number of nodes in the graph
 - m is the missing PageRank mass
- Additional optimization: make it a single pass!

PageRank Convergence

- Alternative convergence criteria
 - Iterate until PageRank values don't change
 - Iterate until PageRank rankings don't change
 - Fixed number of iterations
- Convergence for web graphs?
 - Not a straightforward question
- Watch out for link spam:
 - Link farms
 - Spider traps
 - ...

Beyond PageRank

- Variations of PageRank
 - Weighted edges
 - Personalized PageRank
- Variants on graph random walks
 - Hubs and authorities (HITS)
 - SALSA

Applications

- Static prior for web ranking
- Identification of “special nodes” in a network
- Link recommendation
- Additional feature in any machine learning problem

Other Classes of Graph Algorithms

- Subgraph pattern matching
- Computing simple graph statistics
 - Degree vertex distributions
- Computing more complex graph statistics
 - Clustering coefficients
 - Counting triangles

General Issues for Graph Algorithms

- Sparse vs. dense graphs
- Graph topologies



MapReduce Sucks

- Java verbosity
- Hadoop task startup time
- Stragglers
- Needless graph shuffling
- Checkpointing at each iteration

Iterative Algorithms

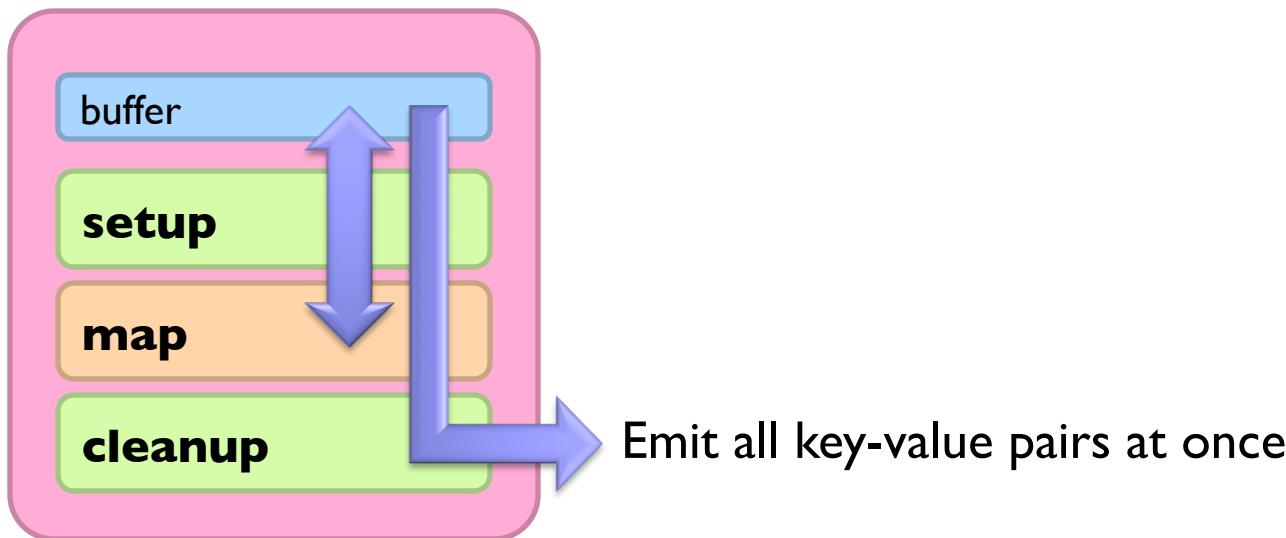


MapReduce sucks at iterative algorithms

- Alternative programming models (later)
- Easy fixes (now)

In-Mapper Combining

- Use combiners
 - Perform local aggregation on map output
 - Downside: intermediate data is still materialized
- Better: in-mapper combining
 - Preserve state across multiple map calls, aggregate messages in buffer, emit buffer contents at end
 - Downside: requires memory management



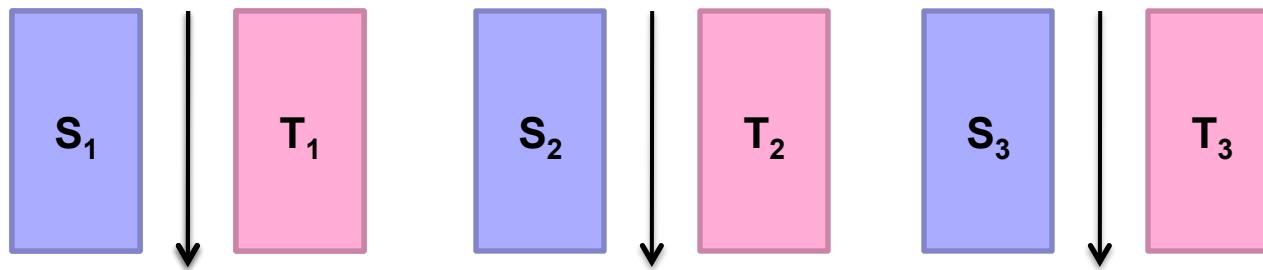
Better Partitioning

- Default: hash partitioning
 - Randomly assign nodes to partitions
- Observation: many graphs exhibit local structure
 - E.g., communities in social networks
 - Better partitioning creates more opportunities for local aggregation
- Unfortunately, partitioning is **hard!**
 - Sometimes, chick-and-egg...
 - But cheap heuristics sometimes available
 - For webgraphs: range partition on domain-sorted URLs

Schimmy Design Pattern

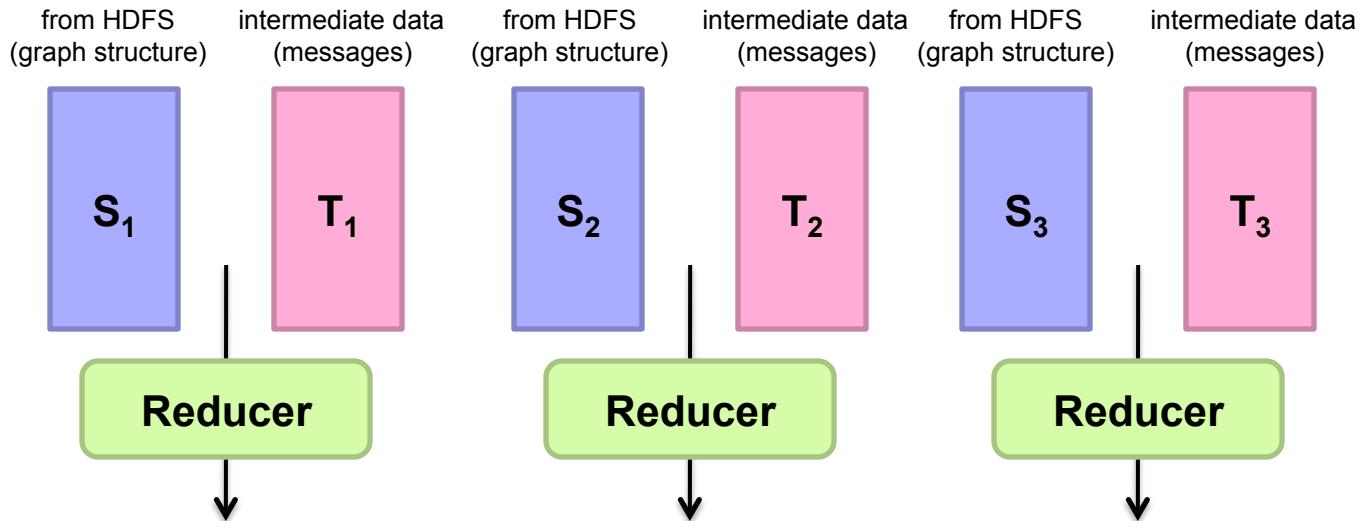
- Basic implementation contains two dataflows:
 - Messages (actual computations)
 - Graph structure (“bookkeeping”)
- Schimmy: separate the two dataflows, shuffle only the messages
 - Basic idea: merge join between graph structure and messages

both relationships consistently partitioned and sorted by join key



Do the Schimmy!

- Schimmy = reduce side parallel merge join between graph structure and messages
 - Consistent partitioning between input and intermediate data
 - Mappers emit only messages (actual computation)
 - Reducers read graph structure directly from HDFS



Experiments

- Cluster setup:

- 10 workers, each 2 cores (3.2 GHz Xeon), 4GB RAM, 367 GB disk
- Hadoop 0.20.0 on RHELS 5.3

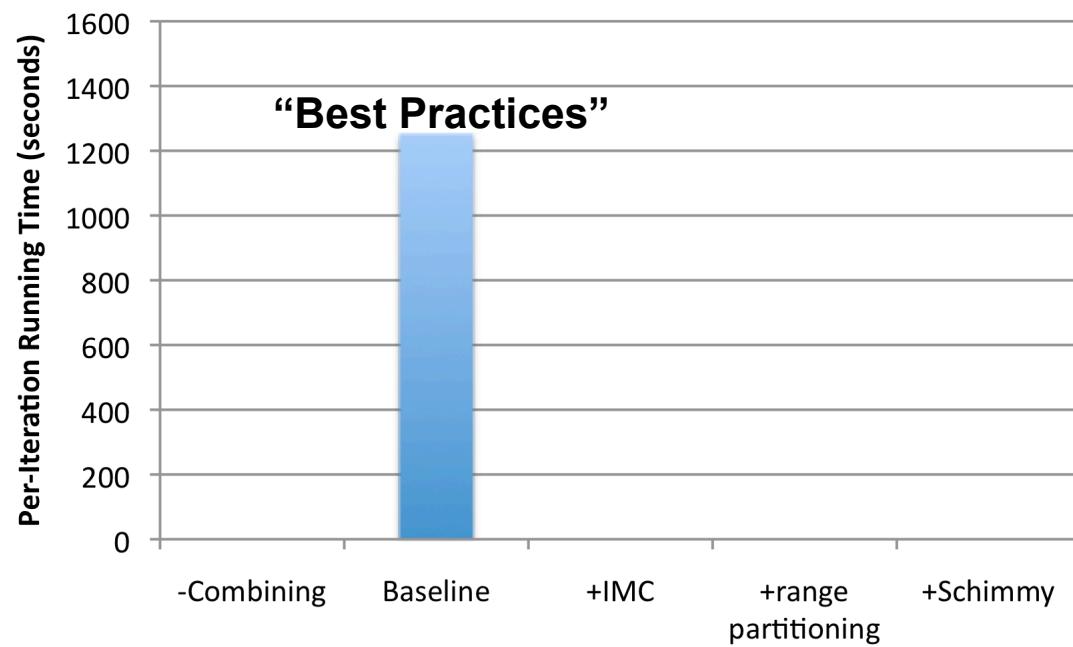
- Dataset:

- First English segment of ClueWeb09 collection
- 50.2m web pages (1.53 TB uncompressed, 247 GB compressed)
- Extracted webgraph: 1.4 billion links, 7.0 GB
- Dataset arranged in crawl order

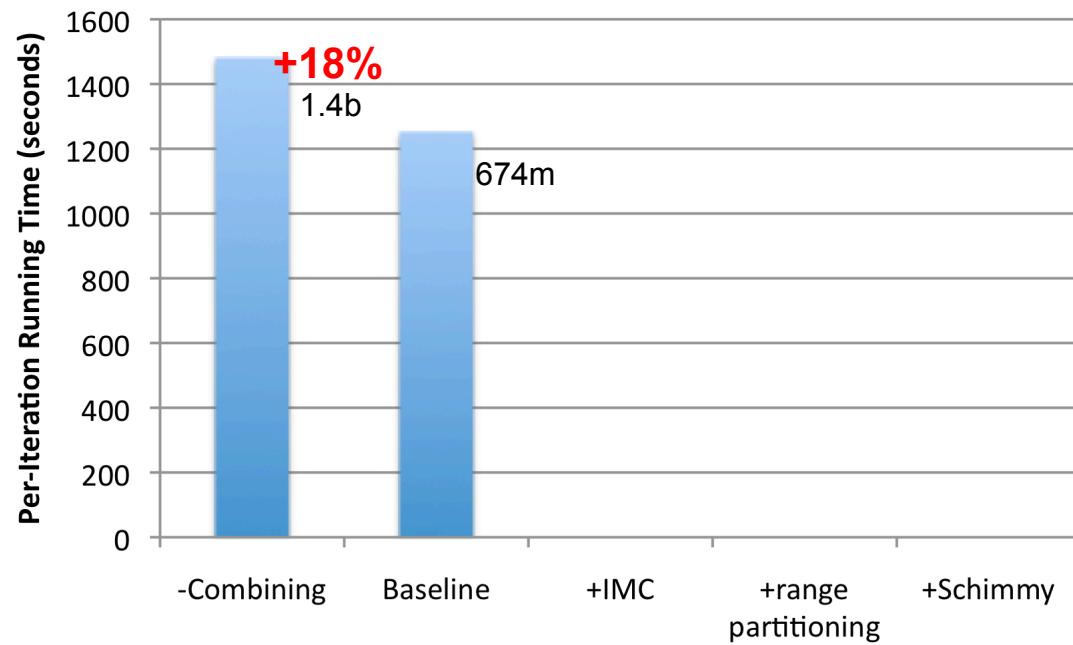
- Setup:

- Measured per-iteration running time (5 iterations)
- 100 partitions

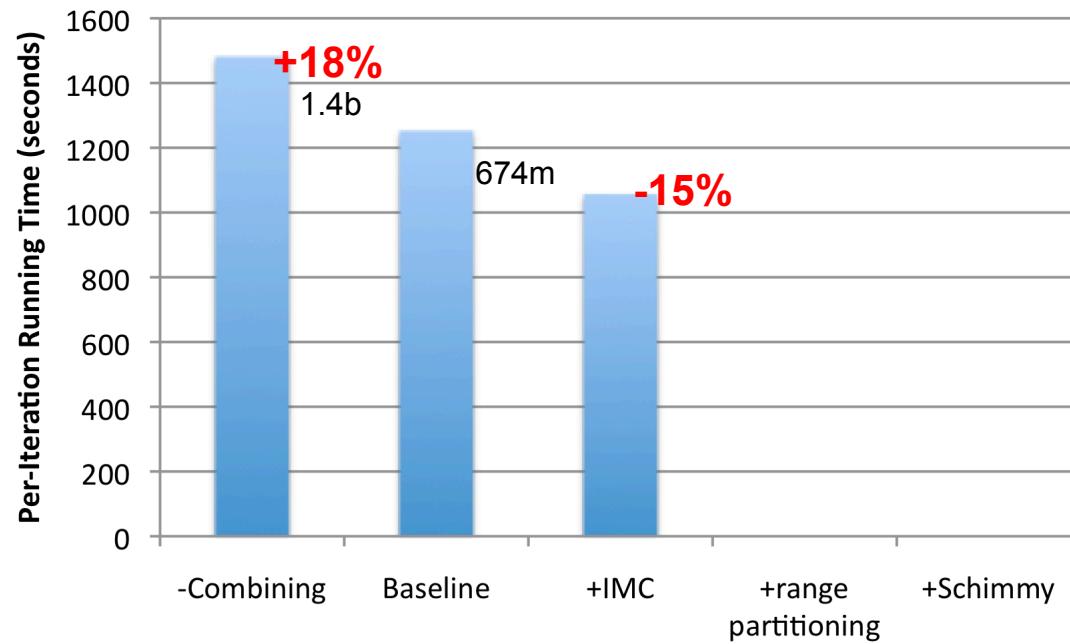
Results



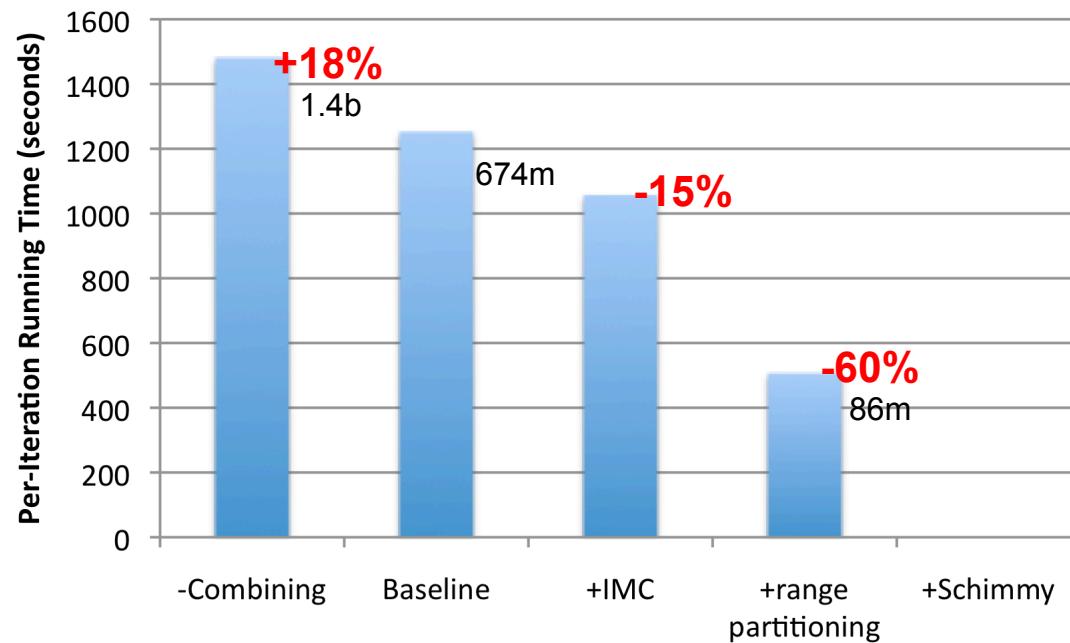
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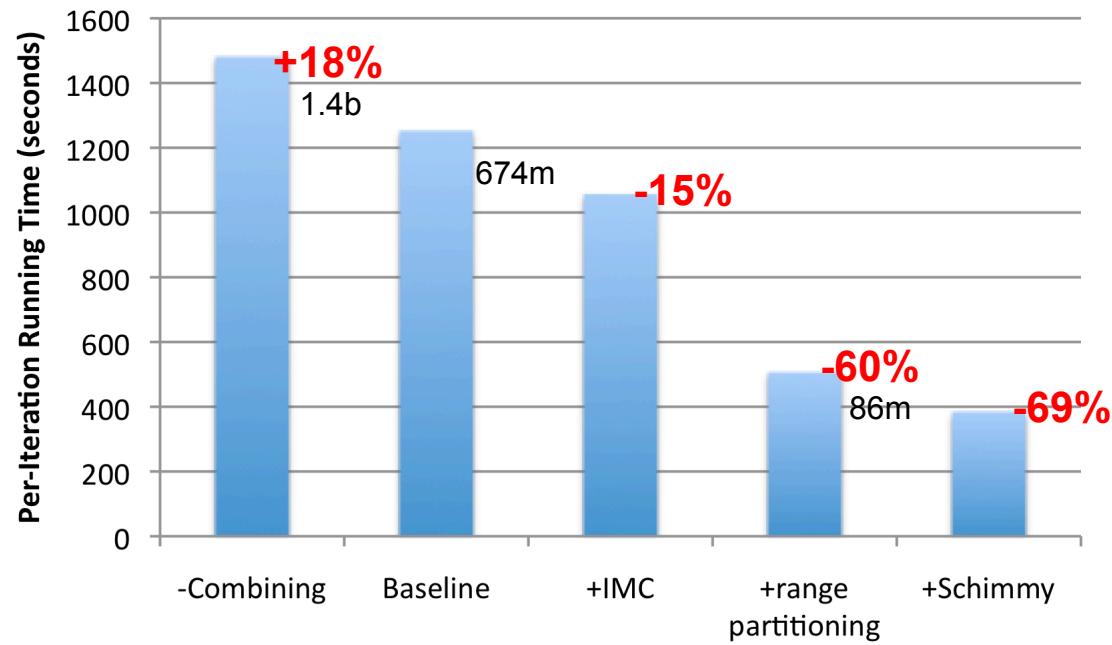
Results



Results



Results



MapReduce sucks at iterative algorithms

- Alternative programming models (later)
- Easy fixes (now)

Later, the “hammer” argument...

Today's Agenda

- Graph problems and representations
- Parallel breadth-first search
- PageRank
- Beyond PageRank and other graph algorithms
- Optimizing graph algorithms

A photograph of a traditional Japanese rock garden. In the foreground, a gravel path is raked into fine, parallel lines. Several large, dark, irregular stones are scattered across the garden. A small, shallow pond is visible in the middle ground, surrounded by more stones and low-lying green plants. In the background, there are more trees and shrubs, and the wooden buildings of a residence are visible behind the garden wall.

Questions?