

Large-Scale Graph Processing

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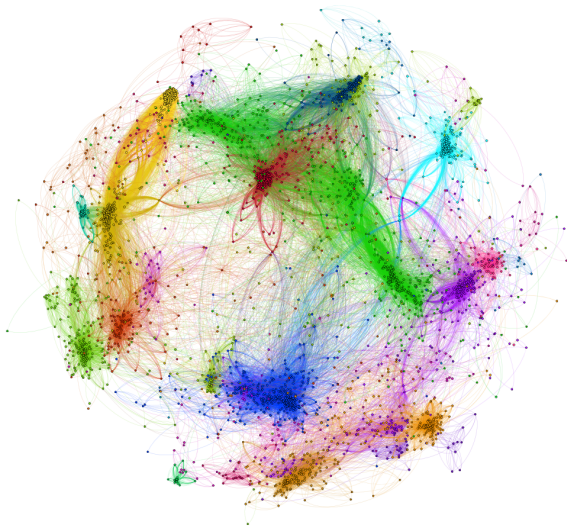
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May 13-15, 2014





- ▶ **Graphs** provide a **flexible abstraction** for describing relationships between **discrete objects**.
- ▶ Many problems can be **modeled by graphs** and solved with appropriate **graph algorithms**.

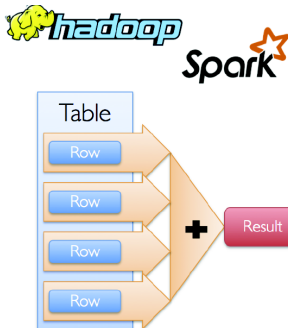
Large Graph



- ▶ Large graphs need large-scale processing.
- ▶ A large graph either cannot fit into memory of single computer or it fits with huge cost.

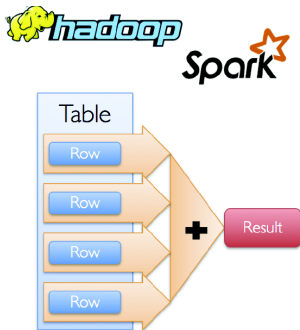
Question

Can we use platforms like [MapReduce](#) or [Spark](#), which are based on **data-parallel** model, for large-scale graph proceeding?



Data-Parallel Model for Large-Scale Graph Processing

- ▶ The platforms that have worked well for developing **parallel applications** are not necessarily effective for **large-scale graph** problems.
- ▶ Why?



Graph Algorithms Characteristics (1/2)

► Unstructured problems

- Difficult to extract parallelism based on partitioning of the data: the irregular structure of graphs.
- Limited scalability: unbalanced computational loads resulting from poorly partitioned data.

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► Data-driven computations

- Difficult to express parallelism based on partitioning of computation: the structure of computations in the algorithm is not known a priori.
- The computations are dictated by nodes and links of the graph.

Graph Algorithms Characteristics (2/2)

- ▶ Poor data locality
 - The computations and data access patterns do not have much locality: the irregular structure of graphs.

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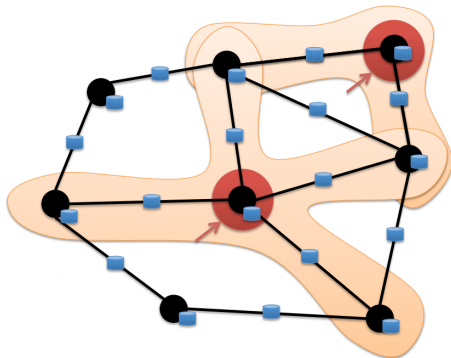
► High data access to computation ratio

- Graph algorithms are often based on exploring the structure of a graph to perform computations on the graph data.
- Runtime can be dominated by waiting memory fetches: low locality.

Graph-Parallel Processing

Proposed Solution

Graph-Parallel Processing



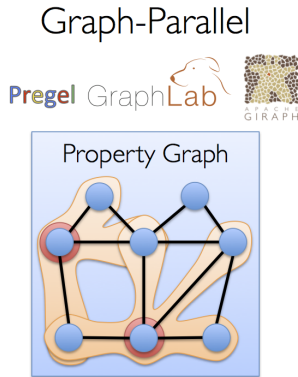
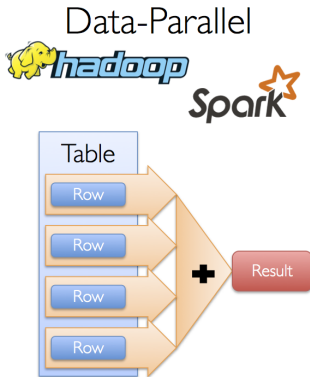
- Computation typically depends on the **neighbors**.

Graph-Parallel Processing

- ▶ Restricts the types of computation.
- ▶ New techniques to partition and distribute graphs.
- ▶ Exploit graph structure.
- ▶ Executes graph algorithms orders-of-magnitude faster than more general data-parallel systems.

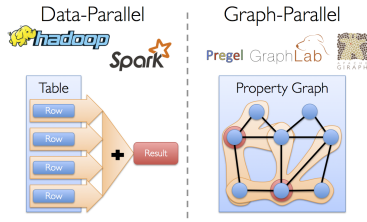


Data-Parallel vs. Graph-Parallel Computation



Data-Parallel vs. Graph-Parallel Computation

- ▶ **Data-parallel** computation
 - **Record-centric** view of data.
 - **Parallelism**: processing **independent** data on separate resources.
- ▶ **Graph-parallel** computation
 - **Vertex-centric** view of graphs.
 - **Parallelism**: partitioning graph (**dependent**) data across processing resources, and **resolving dependencies** (**along edges**) through **iterative** computation and communication.

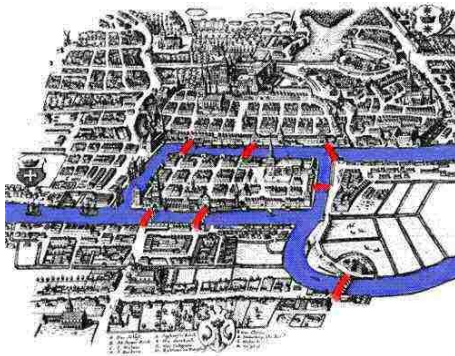


- ▶ Pregel
- ▶ GraphLab
- ▶ PowerGraph
- ▶ GraphX



Seven Bridges of Königsberg

- ▶ Finding a walk through the city that would cross each bridge once and only once.
- ▶ Euler proved that the problem has no solution.



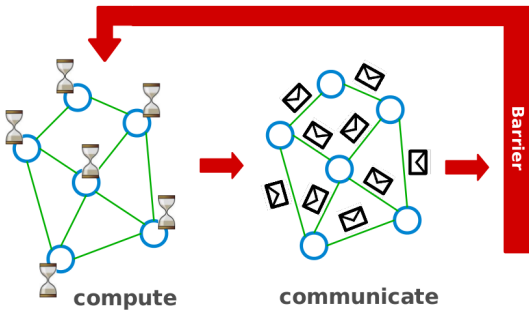
Map of Königsberg in Euler's time, highlighting the river Pregel and the bridges.

- ▶ Large-scale **graph-parallel** processing platform developed at Google.
- ▶ Inspired by **bulk synchronous parallel (BSP)** model.

Bulk Synchronous Parallel (1/2)

- ▶ It is a **parallel programming model**.
- ▶ The model consists of:
 - A set of **processor-memory** pairs.
 - A **communications network** that delivers messages in a **point-to-point** manner.
 - A mechanism for the efficient **barrier synchronization** for all or a subset of the processes.
 - There are **no special** combining, replicating, or broadcasting facilities.

Bulk Synchronous Parallel (2/2)



All vertices update in parallel (at the same time).

Vertex-Centric Programs

- ▶ Think like a vertex.
- ▶ Each vertex computes **individually** its value: in **parallel**
- ▶ Each vertex can see its **local** context, and updates its value accordingly.

- ▶ A **directed graph** that stores the program **state**, e.g., the current value.

Execution Model (1/3)

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 - **reads** messages sent to it in superstep **S-1**.
 - **sends** messages to other vertices: receiving at superstep **S+1**.
 - **modifies** its state.

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- ▶ Vertices communicate directly with one another by **sending messages**.

Execution Model (2/3)

- Superstep 0: all vertices are in the active state.

Execution Model (2/3)

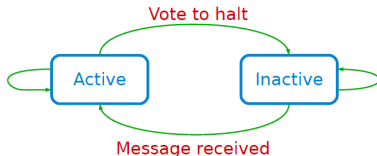
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Execution Model (2/3)

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- ▶ A vertex **deactivates** itself by voting to **halt**: no further work to do.
- ▶ A halted vertex can be active if it **receives a message**.
- ▶ The whole algorithm terminates when:
 - All vertices are **simultaneously inactive**.
 - There are **no messages in transit**.



Execution Model (3/3)

- **Aggregation**: a mechanism for **global** communication, monitoring, and data.

Execution Model (3/3)

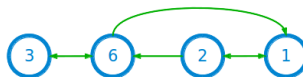
- ▶ **Aggregation**: a mechanism for **global** communication, monitoring, and data.
- ▶ Runs after each **superstep**.
- ▶ Each **vertex** can provide a value to an aggregator in superstep **S** .
- ▶ The system **combines** those values and the resulting value is made available to all vertices in superstep **$S + 1$** .

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- ▶ Each **vertex** can provide a value to an aggregator in superstep S .
- ▶ The system **combines** those values and the resulting value is made available to all vertices in superstep $S + 1$.
- ▶ A number of **predefined aggregators**, e.g., **min**, **max**, **sum**.
- ▶ Aggregation operators should be **commutative** and **associative**.

Example: Max Value (1/4)

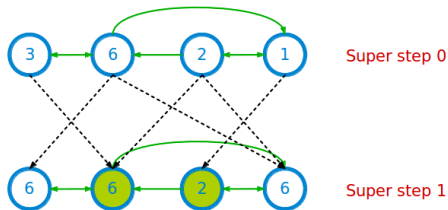
```
i_val := val  
  
for each message m  
  if m > val then val := m  
  
if i_val == val then  
  vote_to_halt  
else  
  for each neighbor v  
    send_message(v, val)
```



Super step 0

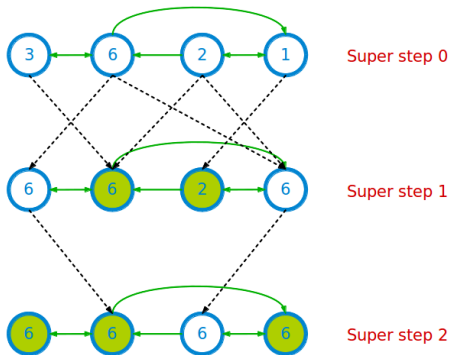
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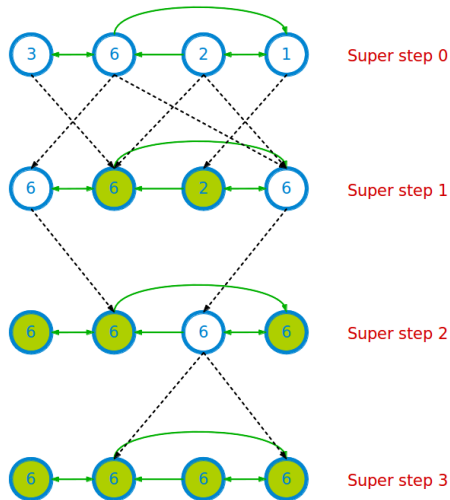
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Example: Max Value (4/4)

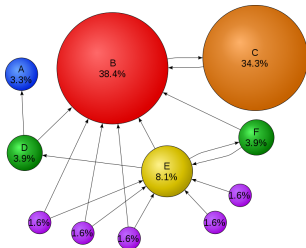
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Example: PageRank

- Update ranks in **parallel**.
- **Iterate** until convergence.

$$R[i] = 0.15 + \sum_{j \in Nbrs(i)} w_{ji} R[j]$$



Example: PageRank

```
Pregel_PageRank(i, messages):  
    // receive all the messages  
    total = 0  
    foreach(msg in messages):  
        total = total + msg  
  
    // update the rank of this vertex  
    R[i] = 0.15 + total  
  
    // send new messages to neighbors  
    foreach(j in out_neighbors[i]):  
        sendmsg(R[i] * wij) to vertex j
```

$$R[i] = 0.15 + \sum_{j \in Nbrs(i)} w_{ji} R[j]$$

Partitioning the Graph

- ▶ The pregel library divides a graph into a number of **partitions**.
- ▶ Each consisting of a set of **vertices** and all of those vertices' **outgoing edges**.
- ▶ Vertices are assigned to partitions based on their **vertex-ID** (e.g., `hash(ID)`).

Implementation (1/4)

- ▶ Master-worker model.
- ▶ User programs are copied on machines.
- ▶ One copy becomes the master.

Implementation (2/4)

- ▶ The **master** is responsible for
 - **Coordinating** workers activity.
 - Determining the **number of partitions**.
- ▶ Each **worker** is responsible for
 - Maintaining the **state** of its partitions.
 - Executing the user's **Compute()** method on its vertices.
 - Managing **messages** to and from other workers.

Implementation (3/4)

- ▶ The **master** assigns one or more **partitions** to each **worker**.

Implementation (3/4)

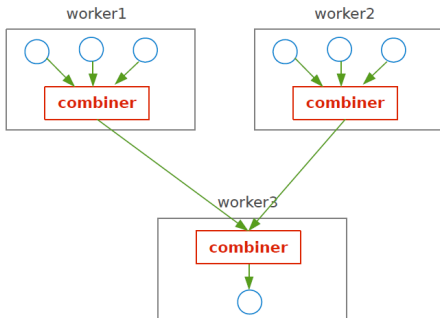
- ▶ The **master** assigns one or more **partitions** to each **worker**.
- ▶ The **master** assigns a portion of **user input** to each **worker**.
 - Set of records containing an arbitrary **number of vertices and edges**.
 - If a worker loads a vertex that **belongs to that worker's partitions**, the appropriate data structures are immediately updated.
 - Otherwise the worker enqueues a message to the **remote peer** that owns the vertex.

Implementation (4/4)

- ▶ After the **input has finished loading**, all vertices are marked as **active**.
- ▶ The master instructs each worker to perform a **superstep**.
- ▶ After the computation **halts**, the master may instruct each worker to save its portion of the graph.

Combiner

- ▶ Sending a message between workers incurs some **overhead**: use **combiner**.
- ▶ This can be reduced in some cases: sometimes vertices only care about a **summary value** for the messages it is sent (e.g., **min**, **max**, **sum**, **avg**).



Fault Tolerance (1/2)

- ▶ Fault tolerance is achieved through **checkpointing**.
- ▶ At **start of each superstep**, master tells workers to **save** their state:
 - Vertex values, edge values, incoming messages
 - Saved to persistent storage
- ▶ Master saves **aggregator values** (if any).
- ▶ This is **not** necessarily done at every superstep: **costly**

Fault Tolerance (2/2)

- ▶ When master **detects** one or more **worker failures**:
 - All workers revert to last **checkpoint**.
 - Continue **from there**.
 - That is a lot of **repeated work**.
 - At least it is better than redoing the whole job.

Pregel Summary

- ▶ Bulk Synchronous Parallel model
- ▶ Vertex-centric
- ▶ Superstep: sequence of iterations
- ▶ Master-worker model
- ▶ Communication: message passing

Pregel Limitations

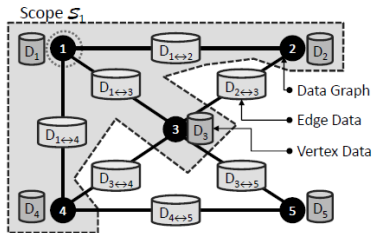
- ▶ **Inefficient** if different regions of the graph converge at **different speed**.
- ▶ Can suffer if one **task** is **more expensive** than the others.
- ▶ Runtime of each phase is determined by the **slowest** machine.



- ▶ A **directed graph** that stores the program **state**, called **data graph**.

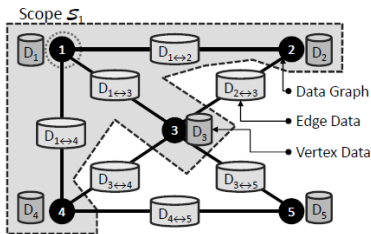
Vertex Scope

- The **scope** of vertex v is the data stored in **vertex v** , in all **adjacent vertices** and **adjacent edges**.



Execution Model (1/4)

- Rather than adopting a **message passing** as in Pregel, GraphLab allows the user defined function of a vertex to **read** and **modify** any of the data in its **scope**.



Execution Model (2/4)

- ▶ **Update** function: user-defined function similar to **Compute** in Pregel.
- ▶ Can **read** and **modify** the data within the **scope** of a vertex.
- ▶ **Schedules** the future execution of other update functions.

Execution Model (3/4)

Input: Data Graph $G = (V, E, D)$

Input: Initial task set $\mathcal{T} = \{(f, v_1), (g, v_2), \dots\}$

while \mathcal{T} is not Empty **do**

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1  |    $(f, v) \leftarrow \text{RemoveNext}(\mathcal{T})$   
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   |
```

Output: Modified Data Graph $G = (V, E, D')$

- ▶ After executing an update function (f, g, \dots) the **modified scope** data in \mathcal{S}_v is **written back** to the data graph.
- ▶ Each **task** in the set of tasks \mathcal{T} , is a tuple (f, v) consisting of an **update function** f and a vertex v .

Execution Model (4/4)

- ▶ **Sync** function: similar to **aggregate** in Pregel.
- ▶ Maintains **global aggregates**.
- ▶ Performs periodically in the **background**.

Example: PageRank

```
GraphLab_PageRank(i)
  // compute sum over neighbors
  total = 0
  foreach(j in in_neighbors(i)):
    total = total + R[j] * wji

  // update the PageRank
  R[i] = 0.15 + total

  // trigger neighbors to run again
  foreach(j in out_neighbors(i)):
    signal vertex-program on j
```

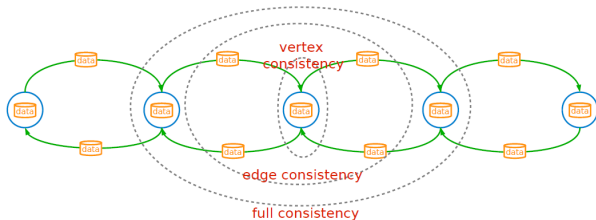
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Data Consistency (1/3)

- ▶ Overlapped scopes: **race-condition** in simultaneous execution of two update functions.

Data Consistency (1/3)

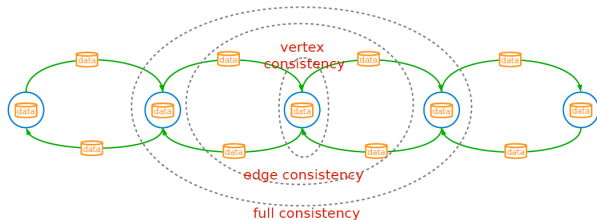
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- **Full** consistency: during the execution $f(v)$, no other function reads or modifies data within the v scope.

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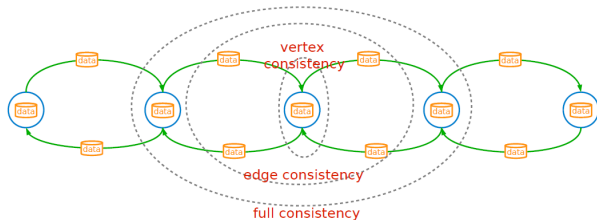
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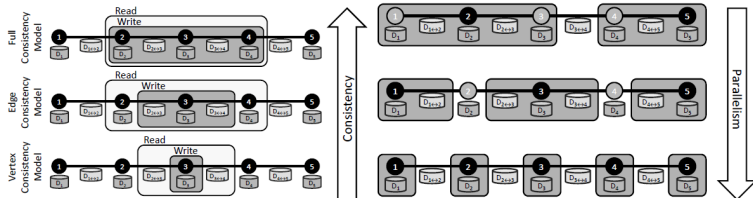
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- ▶ **Vertex** consistency: during the execution $f(v)$, no other function will be applied to v .

Data Consistency (2/3)



Consistency vs. Parallelism

[Low, Y., GraphLab: A Distributed Abstraction for Large Scale Machine Learning (Doctoral dissertation, University of California), 2013.]

- ▶ Proving the **correctness** of a parallel algorithm: **sequential consistency**

Data Consistency (3/3)

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Data Consistency (3/3)

- ▶ Proving the **correctness** of a parallel algorithm: **sequential consistency**
- ▶ **Sequential consistency**: if for every **parallel** execution, there exists a **sequential** execution of update functions that produces an **equivalent** result.
- ▶ A simple method to **achieve serializability** is to ensure that the **scopes** of concurrently executing update functions **do not overlap**.
 - The **full consistency** model is used.
 - The **edge consistency** model is used and update functions do not modify data in adjacent vertices.
 - The **vertex consistency** model is used and update functions only access local vertex data.

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- ▶ Distributed implementation

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Tasks Schedulers (1/2)

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- In what **order** should the **tasks** (**vertex-update function pairs**) be called?
 - A collection of base schedules, e.g., round-robin, and synchronous.
 - **Set scheduler**: enables users to compose custom update schedules.

Tasks Schedulers (2/2)

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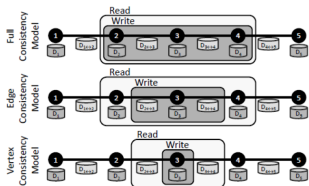
- **FIFO**: only permits task **creation** but do **not permit task reordering**.
- **Prioritized**: **permits task reordering** at the cost of increased overhead.

Consistency

- ▶ Implemented in C++ using PThreads for parallelism.
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- ▶ Implemented in C++ using PThreads for parallelism.
- ▶ Consistency: **read-write lock**
- ▶ **Vertex consistency**
 - Central vertex (**write-lock**)
- ▶ **Edge consistency**
 - Central vertex (**write-lock**)
 - Adjacent vertices (**read-locks**)
- ▶ **Full consistency**
 - Central vertex (**write-locks**)
 - Adjacent vertices (**write-locks**)
- ▶ **Deadlocks** are avoided by acquiring **locks sequentially** following a **canonical order**.



- ▶ Shared memory implementation
- ▶ Distributed implementation

Distributed Implementation

▶ Graph partitioning

- How to efficiently load, partition and distribute the data graph across machines?

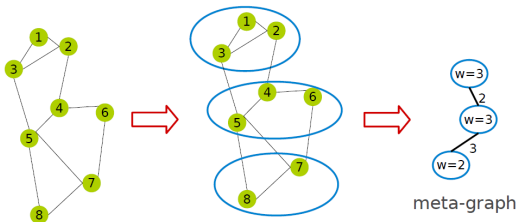
▶ Consistency

- How to achieve consistency in the distributed setting?

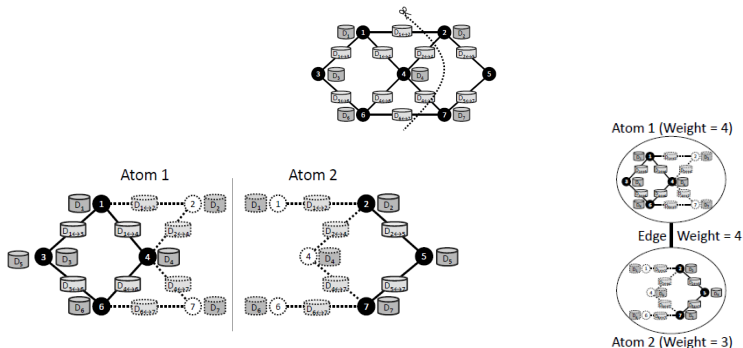
▶ Fault tolerance

Graph Partitioning - Phase 1 (1/2)

- ▶ **Two-phase** partitioning.
- ▶ Partitioning the data graph into k parts, called **atom**: $k \gg$ number of machines.
- ▶ **meta-graph**: the graph of atoms (one vertex for each atom).
- ▶ **Atom weight**: the amount of data it stores.
- ▶ **Edge weight**: the number of edges crossing the atoms.



Graph Partitioning - Phase 1 (2/2)



- ▶ Each **atom** is stored as a separate file on a distributed storage system, e.g., HDFS.
- ▶ Each atom file is a simple **binary** that stores **interior** and the **ghosts** of the partition information.
- ▶ **Ghost**: set of **vertices and edges adjacent** to the partition boundary.

Graph Partitioning - Phase 2

- ▶ Meta-graph is very **small**.
- ▶ A **fast balanced partition** of the **meta-graph** over the physical machines.
- ▶ Assigning graph atoms to machines.

- ▶ To achieve a **serializable parallel execution** of a set of **dependent tasks**.
- ▶ Chromatic Engine
- ▶ Distributed Locking Engine

- ▶ Construct a **vertex coloring**: assigns a color to each vertex such that **no adjacent** vertices share the same color.

Consistency - Chromatic Engine

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- ▶ **Full consistency**: no vertex shares the same color as any of its distance **two** neighbors.
- ▶ **Vertex consistency**: assigning all vertices the **same** color.

Consistency - Distributed Locking Engine

- ▶ Associating a **readers-writer** lock with each vertex.
- ▶ **Vertex consistency**
 - Central vertex (**write-lock**)
- ▶ **Edge consistency**
 - Central vertex (**write-lock**), Adjacent vertices (**read-locks**)
- ▶ **Full consistency**
 - Central vertex (**write-locks**), Adjacent vertices (**write-locks**)
- ▶ **Deadlocks** are avoided by acquiring **locks sequentially** following a **canonical order**.

Fault Tolerance - Synchronous

- ▶ The systems **periodically** signals all computation activity to **halt**.
- ▶ Then synchronizes all caches (ghosts) and **saves to disk** all data which has been modified since the last snapshot.
- ▶ **Simple**, but eliminates the systems advantage of **asynchronous** computation.

Fault Tolerance - Asynchronous

- ▶ Based on the [Chandy-Lamport](#) algorithm.
- ▶ The [snapshot](#) function is implemented as an [update function](#) in vertices.
- ▶ The Snapshot update takes [priority](#) over all other update functions.
- ▶ [Edge Consistency](#) is used on all update functions.

```
if v was already snapshotted then
```

```
    Quit
```

```
Save  $D_v$  // Save current vertex
```

```
// Save all edges connected to un-snapshotted vertices
```

```
foreach  $u \in N[v]$  do // Loop over neighbors
```

```
    if u was not snapshotted then
```

```
        Save  $D_{u \rightarrow v}$  if edge  $u \rightarrow v$  exists
```

```
        Save  $D_{v \rightarrow u}$  if edge  $v \rightarrow u$  exists
```

```
        Reschedule u for a Snapshot Update
```

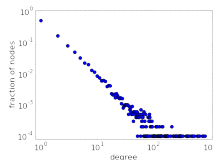
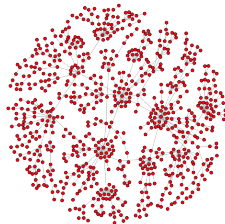
```
Mark v as snapshotted
```

- ▶ **Asynchronous** model
- ▶ Vertex-centric
- ▶ Communication: distributed **shared memory**
- ▶ Three **consistency** levels: full, edge-level, and vertex-level

- ▶ Poor performance on **Natural** graphs.

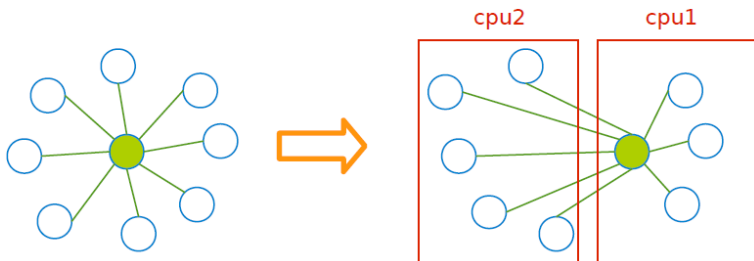
Natural Graphs

- ▶ Graphs derived from **natural phenomena**.
- ▶ Skewed **Power-Law** degree distribution.



Natural Graphs Challenges

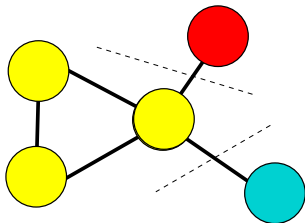
- ▶ Traditional graph-partitioning algorithms (**edge-cut** algorithms) perform **poorly** on Power-Law Graphs.
- ▶ Challenges of **high-degree** vertices.



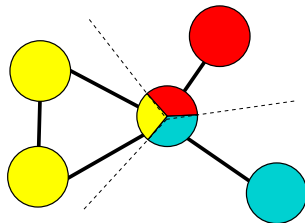
Vertex-Cut Partitioning

Proposed Solution

Vertex-Cut Partitioning

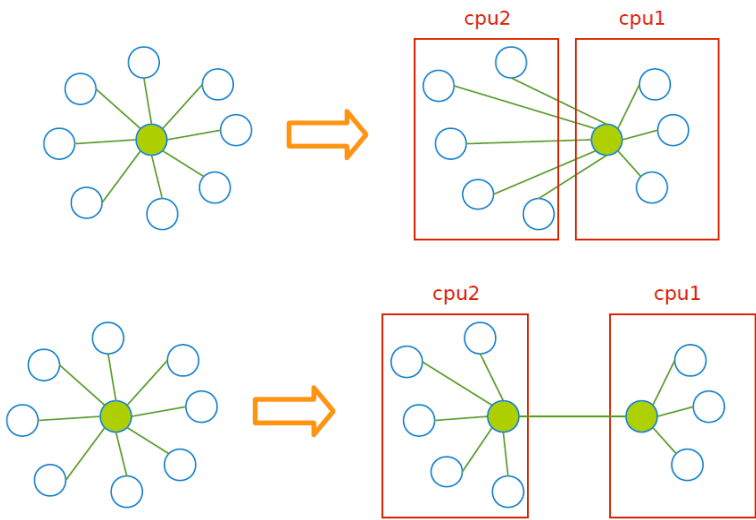


Edge-cut

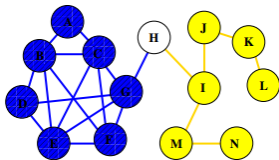


Vertex-cut

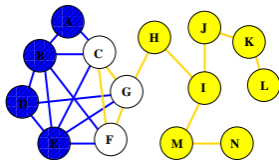
Edge-cut vs. Vertex-cut Partitioning



Edge-cut vs. Vertex-cut Partitioning



Edge-cut



Vertex-cut

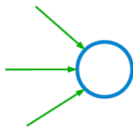
PowerGraph

- ▶ Vertex-cut partitioning of graphs.
- ▶ Factorizes the GraphLab update function into the Gather, Apply and Scatter phases (GAS).

Gather-Apply-Scatter Programming Model

► Gather

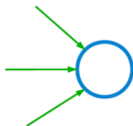
- **Accumulate** information about neighborhood through a generalized **sum**.



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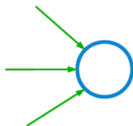
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- Apply the accumulated value to center vertex.

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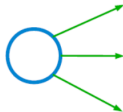


► Apply

- **Apply** the accumulated value to center vertex.

► Scatter

- **Update** adjacent edges and vertices.



- ▶ A **directed graph** that stores the program **state**, called **data graph**.

Execution Model (1/2)

- ▶ **Vertex-centric** programming: implementing the **GASVertexProgram** interface (**vertex-program** for short).
- ▶ Similar to **Compute** in **Pregel**, and **update** function in **GraphLab**.

```
interface GASVertexProgram(u) {  
  // Run on gather_nbrs(u)  
  gather( $D_u$ ,  $D_{u-v}$ ,  $D_v$ )  $\rightarrow$   $Accum$   
  sum( $Accum$  left,  $Accum$  right)  $\rightarrow$   $Accum$   
  apply( $D_u$ ,  $Accum$ )  $\rightarrow D_u^{new}$   
  // Run on scatter_nbrs(u)  
  scatter( $D_u^{new}$ ,  $D_{u-v}$ ,  $D_v$ )  $\rightarrow$  ( $D_{u-v}^{new}$ ,  $Accum$ )  
}
```


Execution Model (2/2)

Input: Center vertex u

if *Cache Disabled* **then**

 // Basic Gather-Apply-Scatter Model

foreach neighbor v in $\text{gather_nbrs}(u)$ **do**

$a_u \leftarrow \text{sum}(a_u, \text{gather}(D_u, D_{u-v}, D_v))$

$D_u \leftarrow \text{apply}(D_u, a_u)$

foreach neighbor v in $\text{scatter_nbrs}(u)$ **do**

$(D_{u-v}) \leftarrow \text{scatter}(D_u, D_{u-v}, D_v)$

else if *Cache Enabled* **then**

 // Faster GAS Model with Delta Caching

if *cached accumulator a_u is empty* **then**

foreach neighbor v in $\text{gather_nbrs}(u)$ **do**

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if a_v and Δa are not Empty **then** $a_v \leftarrow \text{sum}(a_v, \Delta a)$

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Example: PageRank

```
PowerGraph_PageRank(i):  
  Gather(j -> i):  
    return wji * R[j]  
  
  sum(a, b):  
    return a + b  
  
  // total: Gather and sum  
  Apply(i, total):  
    R[i] = 0.15 + total  
  
  Scatter(i -> j):  
    if R[i] changed then activate(j)
```

$$R[i] = 0.15 + \sum_{j \in Nbrs(i)} w_{ji} R[j]$$

Scheduling (1/5)

Input: Data Graph $G = (V, E, D)$

Input: Initial task set $\mathcal{T} = \{(f, v_1), (g, v_2), \dots\}$

while \mathcal{T} is not Empty **do**

```
1    $(f, v) \leftarrow \text{RemoveNext}(\mathcal{T})$   
2    $(\mathcal{T}', \mathcal{S}_v) \leftarrow f(v, \mathcal{S}_v)$   
3    $\mathcal{T} \leftarrow \mathcal{T} \cup \mathcal{T}'$ 
```

Output: Modified Data Graph $G = (V, E, D')$

- PowerGraph inherits the **dynamic scheduling** of **GraphLab**.

Scheduling (2/5)

- Initially all vertices are active.

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- ▶ The order of executing activated vertices is up to the PowerGraph execution engine.
- ▶ Once a vertex-program completes the scatter phase it becomes inactive until it is reactivated.
- ▶ Vertices can activate themselves and neighboring vertices.

- ▶ PowerGraph can execute both **synchronously** and **asynchronously**.
 - Bulk synchronous execution
 - Asynchronous execution

Scheduling - Bulk Synchronous Execution (4/5)

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- ▶ **Super-step**: a complete series of GAS minor-steps.

Scheduling - Bulk Synchronous Execution (4/5)

- ▶ Similar to Pregel.
- ▶ **Minor-step**: executing the **gather, apply, and scatter in order**.
 - Runs **synchronously** on all **active** vertices with a **barrier** at the end.
- ▶ **Super-step**: a complete series of GAS minor-steps.
- ▶ Changes made to the vertex/edge data are committed at the **end** of each **minor-step** and are visible in the **subsequent minor-steps**.

Scheduling - Asynchronous Execution (5/5)

- ▶ Changes made to the vertex/edge data during the **apply and scatter** functions are **immediately** committed to the graph.
 - **Visible** to subsequent computation on neighboring vertices.

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 - **GraphLab** implements **Dijkstras** solution, where forks are acquired **sequentially** according to a total ordering.
 - **PowerGraph** implements **Chandy-Misra** solution, which acquires all forks **simultaneously**.

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- ▶ The gather operation is invoked on all neighbors: wasting computation cycles
- ▶ Maintaining a cache of the accumulator a_v from the previous gather phase for each vertex.
- ▶ The scatter can return an additional Δa , which is added to the cached accumulator a_v .

Delta Caching (2/2)

Input: Center vertex u

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 // Basic Gather-Apply-Scatter Model

foreach neighbor v in $\text{gather_nbrs}(u)$ **do**

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~~**else if** *Cache Enabled* **then**~~

~~// Faster GAS Model with Delta Caching~~

~~**if** *cached accumulator* a_u *is empty* **then**~~

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Example: PageRank (Delta-Caching)

```
PowerGraph_PageRank(i):  
  Gather(j -> i):  
    return wji * R[j]  
  
sum(a, b):  
  return a + b  
  
// total: Gather and sum  
Apply(i, total):  
  new = 0.15 + total  
  R[i].delta = new - R[i]  
  R[i] = new  
  
Scatter(i -> j):  
  if R[i] changed then activate(j)  
  return wij * R[i].delta
```

$$R[i] = 0.15 + \sum_{j \in \text{Nbrs}(i)} w_{ji} R[j]$$

Graph Partitioning

- ▶ Vertex-cut partitioning.
- ▶ Evenly assign edges to machines.
 - Minimize machines spanned by each vertex.
- ▶ Two proposed solutions:
 - Random edge placement.
 - Greedy edge placement.

- ▶ Randomly assign edges to machines.
- ▶ Completely parallel and easy to distribute.
- ▶ High replication factor.

Greedy Vertex-Cuts (1/2)

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- ▶ **Case 4**: If neither vertex has been assigned, then assign the edge to the least loaded machine.

Greedy Vertex-Cuts (2/2)

- ▶ **Coordinated** edge placement:
 - Requires coordination to place each edge
 - **Slower**, but **higher** quality cuts
- ▶ **Oblivious** edge placement:
 - Approx. greedy objective without coordination
 - **Faster**, but **lower** quality cuts

PowerGraph Summary

- ▶ **Gather-Apply-Scatter** programming model
- ▶ **Synchronous** and **Asynchronous** models
- ▶ **Vertex-cut** graph partitioning

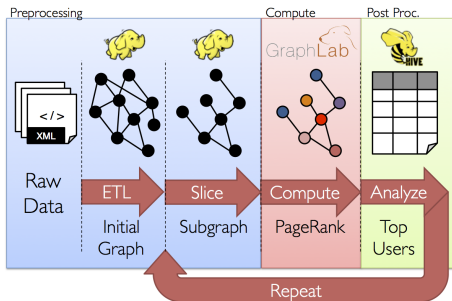
► Any limitations?

Data-Parallel vs. Graph-Parallel Computation

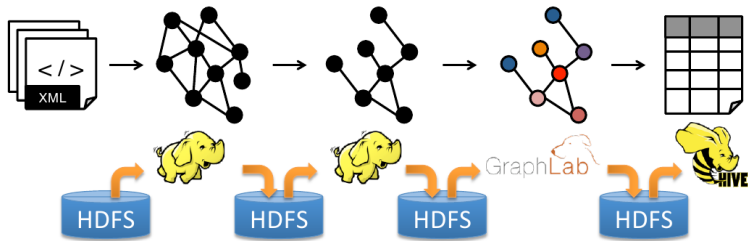
- ▶ **Graph-parallel** computation: **restricting** the types of computation to achieve **performance**.

Data-Parallel vs. Graph-Parallel Computation

- ▶ **Graph-parallel** computation: **restricting** the types of computation to achieve **performance**.
- ▶ **But**, the same restrictions make it **difficult** and **inefficient** to express many stages in a typical graph-analytics **pipeline**.

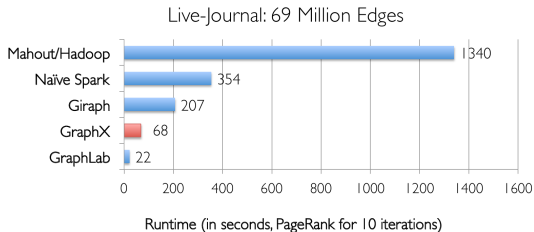


Data-Parallel and Graph-Parallel Pipeline

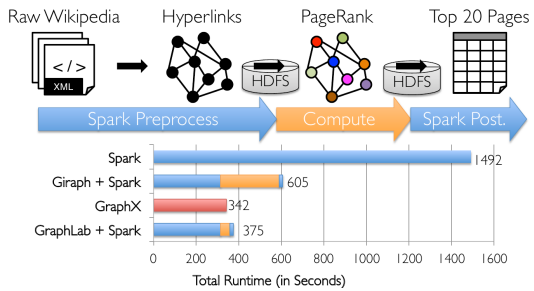
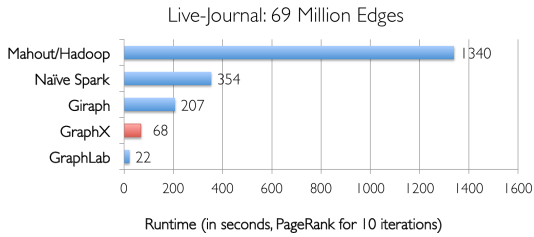


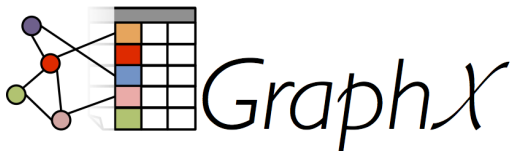
- ▶ **Moving** between **table** and **graph** views of the **same physical data**.
- ▶ **Inefficient**: extensive **data movement** and **duplication** across the network and file system.

GraphX vs. Data-Parallel/Graph-Parallel Systems



GraphX vs. Data-Parallel/Graph-Parallel Systems

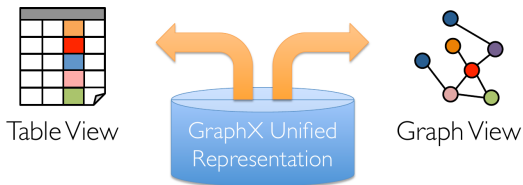




- ▶ New API that blurs the distinction between Tables and Graphs.
- ▶ New system that unifies Data-Parallel and Graph-Parallel systems.
- ▶ It is implemented on top of Spark.

Unifying Data-Parallel and Graph-Parallel Analytics

- ▶ **Tables** and **Graphs** are **composable views** of the same physical data.
- ▶ Each view has its **own operators** that **exploit the semantics** of the view to achieve **efficient** execution.

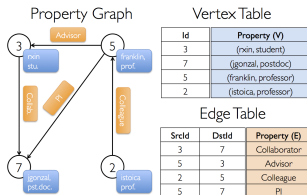


Data Model

► **Property Graph**: represented using **two** Spark **RDDs**:

- **Edge collection**: VertexRDD
- **Vertex collection**: EdgeRDD

```
// VD: the type of the vertex attribute
// ED: the type of the edge attribute
class Graph[VD, ED] {
  val vertices: VertexRDD[VD]
  val edges: EdgeRDD[ED]
}
```



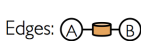
Primitive Data Types

```
// Vertex collection
class VertexRDD[VD] extends RDD[(VertexId, VD)]

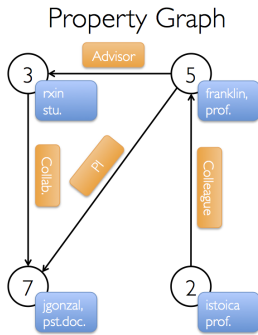
// Edge collection
class EdgeRDD[ED] extends RDD[Edge[ED]]
case class Edge[ED](srcId: VertexId = 0, dstId: VertexId = 0,
                    attr: ED = null.asInstanceOf[ED])

// Edge Triple
class EdgeTriplet[VD, ED] extends Edge[ED]
```

- **EdgeTriplet** represents an **edge** along with the **vertex attributes** of its **neighboring** vertices.



Example (1/3)



Vertex Table

Id	Property (V)
3	(rxin, student)
7	(jgonzal, postdoc)
5	(franklin, professor)
2	(istoica, professor)

Edge Table

SrcId	DstId	Property (E)
3	7	Collaborator
5	3	Advisor
2	5	Colleague
5	7	PI

Example (2/3)

```
val sc: SparkContext

// Create an RDD for the vertices
val users: VertexRDD[(String, String)] = sc.parallelize(
    Array((3L, ("rxin", "student")), (7L, ("jgonzal", "postdoc")),
        (5L, ("franklin", "prof")), (2L, ("istoica", "prof"))))

// Create an RDD for edges
val relationships: EdgeRDD[String] = sc.parallelize(
    Array(Edge(3L, 7L, "collab"),      Edge(5L, 3L, "advisor"),
        Edge(2L, 5L, "colleague"), Edge(5L, 7L, "pi")))

// Define a default user in case there are relationship with missing user
val defaultUser = ("John Doe", "Missing")

// Build the initial Graph
val userGraph: Graph[(String, String), String] =
    Graph(users, relationships, defaultUser)
```

Example (3/3)

```
// Constructed from above
val userGraph: Graph[(String, String), String]

// Count all users which are postdocs
userGraph.vertices.filter((id, (name, pos)) => pos == "postdoc").count

// Count all the edges where src > dst
userGraph.edges.filter(e => e.srcId > e.dstId).count

// Use the triplets view to create an RDD of facts
val facts: RDD[String] = graph.triplets.map(triplet =>
    triplet.srcAttr._1 + " is the " +
    triplet.attr + " of " + triplet.dstAttr._1)

// Remove missing vertices as well as the edges to connected to them
val validGraph = graph.subgraph(vpred = (id, attr) => attr._2 != "Missing")

facts.collect.foreach(println(_))
```

Property Operators (1/2)

```
class Graph[VD, ED] {  
  def mapVertices[VD2](map: (VertexId, VD) => VD2): Graph[VD2, ED]  
  
  def mapEdges[ED2](map: Edge[ED] => ED2): Graph[VD, ED2]  
  
  def mapTriplets[ED2](map: EdgeTriplet[VD, ED] => ED2): Graph[VD, ED2]  
}
```

- ▶ They yield **new graphs** with the vertex or edge properties modified by the map function.
- ▶ The graph **structure** is **unaffected**.

Property Operators (2/2)

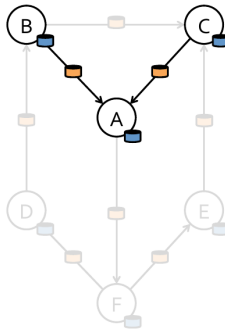
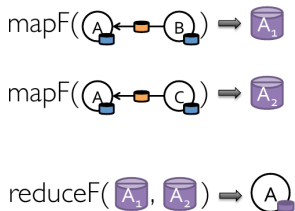
```
val newGraph = graph.mapVertices((id, attr) => mapUdf(id, attr))
```

```
val newVertices = graph.vertices.map((id, attr) => (id, mapUdf(id, attr)))  
val newGraph = Graph(newVertices, graph.edges)
```

- ▶ Both are logically equivalent, but the second one **does not preserve** the structural indices and would not benefit from the GraphX system **optimizations**.

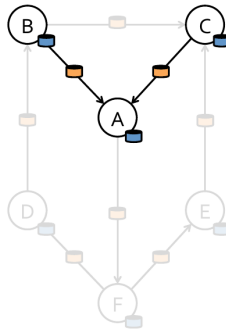
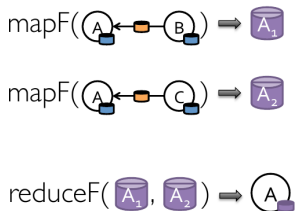
Map Reduce Triplets

- ▶ Map-Reduce for each vertex



Map Reduce Triplets

- Map-Reduce for each vertex



```
// what is the age of the oldest follower for each user?  
val oldestFollowerAge = graph.mapReduceTriplets(  
  e => (e.dstAttr, e.srcAttr), // Map  
  (a, b) => max(a, b) // Reduce  
) .vertices
```

Structural Operators

```
class Graph[VD, ED] {  
  // returns a new graph with all the edge directions reversed  
  def reverse: Graph[VD, ED]  
  
  // returns the graph containing only the vertices and edges that satisfy  
  // the vertex predicate  
  def subgraph(epred: EdgeTriplet[VD,ED] => Boolean,  
              vpred: (VertexId, VD) => Boolean): Graph[VD, ED]  
  
  // a subgraph by returning a graph that contains the vertices and edges  
  // that are also found in the input graph  
  def mask[VD2, ED2](other: Graph[VD2, ED2]): Graph[VD, ED]  
}
```

Structural Operators Example

```
// Build the initial Graph  
val graph = Graph(users, relationships, defaultUser)  
  
// Run Connected Components  
val ccGraph = graph.connectedComponents()  
  
// Remove missing vertices as well as the edges to connected to them  
val validGraph = graph.subgraph(vpred = (id, attr) => attr._2 != "Missing")  
  
// Restrict the answer to the valid subgraph  
val validCCGraph = ccGraph.mask(validGraph)
```

Join Operators

- To join data from external collections (RDDs) with graphs.

```
class Graph[VD, ED] {  
  // joins the vertices with the input RDD and returns a new graph  
  // by applying the map function to the result of the joined vertices  
  def joinVertices[U](table: RDD[(VertexId, U)])  
    (map: (VertexId, VD, U) => VD): Graph[VD, ED]  
  
  // similarly to joinVertices, but the map function is applied to  
  // all vertices and can change the vertex property type  
  def outerJoinVertices[U, VD2](table: RDD[(VertexId, U)])  
    (map: (VertexId, VD, Option[U]) => VD2): Graph[VD2, ED]  
}
```

Graph Builders

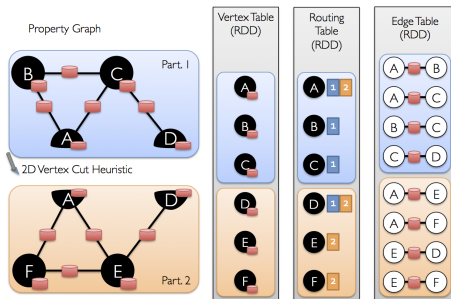
```
// load a graph from a list of edges on disk
object GraphLoader {
  def edgeListFile(
    sc: SparkContext,
    path: String,
    canonicalOrientation: Boolean = false,
    minEdgePartitions: Int = 1)
    : Graph[Int, Int]
}

// graph file
# This is a comment
2 1
4 1
1 2
```

- ▶ GraphX is implemented on top of Spark
- ▶ In-memory caching
- ▶ Lineage-based fault tolerance
- ▶ Programmable partitioning

Distributed Graph Representation (1/2)

- ▶ Representing graphs using **two RDDs**: **edge-collection** and **vertex-collection**
- ▶ **Vertex-cut** partitioning (like **PowerGraph**)



Distributed Graph Representation (2/2)

- ▶ Each vertex partition contains a **bitmask** and **routing table**.
- ▶ **Routing table**: a **logical map** from a vertex id to the set of edge partitions that contains adjacent edges.
- ▶ **Bitmask**: enables the set intersection and filtering.
 - Vertices bitmasks are updated after each operation (e.g., mapReduceTriplets).
 - Vertices hidden by the bitmask **do not** participate in the graph operations.

Summary

▶ Pregel

- Synchronous model: super-step
- Message passing

▶ GraphLab

- Asynchronous model: distributed shared-memory
- Edge-cut partitioning

▶ PowerGraph

- GAS programming model
- Vertex-cut partitioning

▶ GraphX

- Unifying data-parallel and graph-parallel analytics
- Vertex-cut partitioning

Questions?

Acknowledgements

Some pictures were derived from the Spark web site
(<http://spark.apache.org/>).