



MCSN - N. Tonellotto - Distributed Enabling Platforms

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- Written locally
- Transferred from mappers to reducers over network
- •Issue
 - Performance bottleneck
- Solution
 - Use combiners
 - Use In-Mapper Combining







1:	class MAPPER
2:	method MAP(docid a , doc d)
3:	for all term $t \in \text{doc } d$ do
4:	EMIT(term t , count 1)
1:	class Reducer
2:	method REDUCE(term t , counts $[c_1, c_2,]$)
3:	$sum \leftarrow 0$
4:	for all count $c \in$ counts $[c_1, c_2, \ldots]$ do
5:	$sum \leftarrow sum + c$
6:	Еміт(term <i>t</i> , count <i>sum</i>)

- How many intermediate keys per mapper?
- How can we improve this?
- Is it a "real" improvement?







- 1: class MAPPER
- 2: **method** MAP(docid a, doc d) 3: $H \leftarrow$ new AssociativeArray
- 4: **for all** term $t \in \operatorname{doc} d$ **do**
- 5: $H\{t\} \leftarrow H\{t\} + 1$
- 6: **for all** term $t \in H$ **do**
- 7: EMIT(term t, count $H{t}$)

- Custom local aggregator
- Coding overhead
- Is it a "real" improvement?





- 1: class Mapper
- 2: method Initialize
- 3: $H \leftarrow \text{new AssociativeArray}$
- 4: **method** MAP(docid a, doc d)
- 5: **for all** term $t \in \operatorname{doc} d$ **do**
- 6: $H\{t\} \leftarrow H\{t\} + 1$
- 7: method CLOSE
- 8: for all term $t \in H$ do
- 9: $E_{MIT}(term t, count H{t})$

- Custom local aggregator
- Coding overhead
- Is it a "real" improvement?







Advantages:

- Complete local aggregation control (how and when)
- Guaranteed to execute
- Direct efficiency control on intermediate data creation
- Avoid unnecessary objects creation and destruction (before combiners)

Disadvantages:

- Breaks the functional programming background (state)
- Potential ordering-dependent bugs
- Memory scalability bottleneck (solved by memory foot-printing and flushing)







- Common problem:
 - Given an input of size N, generate an output matrix of size N x

• Example: word co-occurrence matrix

- Given a document collection, emit the bigram frequencies









1:	class MAPPER
2:	method MAP(docid a , doc d)
3:	for all term $w \in \operatorname{doc} d$ do
4:	for all term $u \in \text{Neighbors}(w)$ do
5:	Еміт(pair (w, u) , count 1)
1:	class Reducer
2:	method REDUCE(pair p , counts $[c_1, c_2,]$)
3:	$s \leftarrow 0$
4:	for all count $c \in$ counts $[c_1, c_2, \ldots]$ do
5:	$s \leftarrow s + c$
6:	Еміт(pair p , count s)

▷ Emit count for each co-occurrence

▷ Sum co-occurrence counts

- We must use custom key type
- Intermediate overhead? Bottlenecks?
- Can we use the reducer as a combiner?
- Keys space?





6:



- 2: **method** MAP(docid a, doc d)
- 3: **for all** term $w \in \operatorname{doc} d$ **do**
- 4: $H \leftarrow \text{new AssociativeArray}$
- 5: for all term $u \in NEIGHBORS(w)$ do
 - $H\{u\} \leftarrow H\{u\} + 1$
- 7: EMIT(Term w, Stripe H)
- 1: class Reducer
- 2: method REDUCE(term w, stripes $[H_1, H_2, H_3, \ldots]$)
- 3: $H_f \leftarrow \text{new AssociativeArray}$
- 4: for all stripe $H \in$ stripes $[H_1, H_2, H_3, \ldots]$ do
- 5: $\operatorname{Sum}(H_f, H)$
- 6: EMIT(term w, stripe H_f)
- We must use custom key and value types
- Intermediate overhead? Bottlenecks?
- Can we use the reducer as a combiner?
- Keys space?



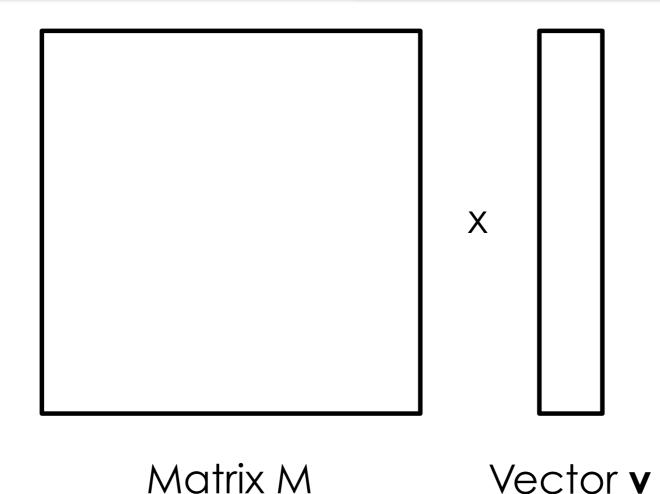
 \triangleright Tally words co-occurring with w

 \triangleright Element-wise sum

"Stripes"







• The matrix does not fit in memory

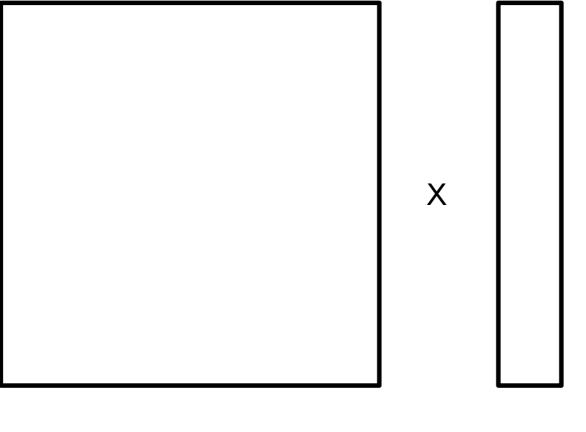
- 1 case: vector v fits in memory
- 2 case: vector **v** does not fit in memory





Vector fits in memory





Matrix M

Vector **v**

• Map

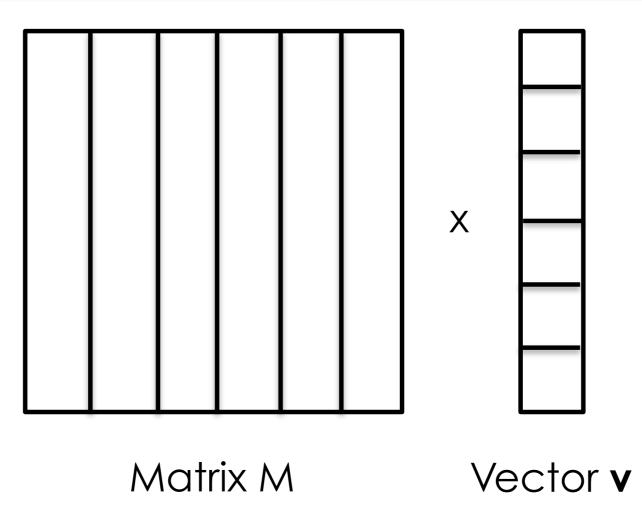
- input = (*, chunk of matrix M)
- vector v read from memory
- output = (i, mijvj)
- Reduce
 - sum up all the values for the given key i





Vector does not in memory



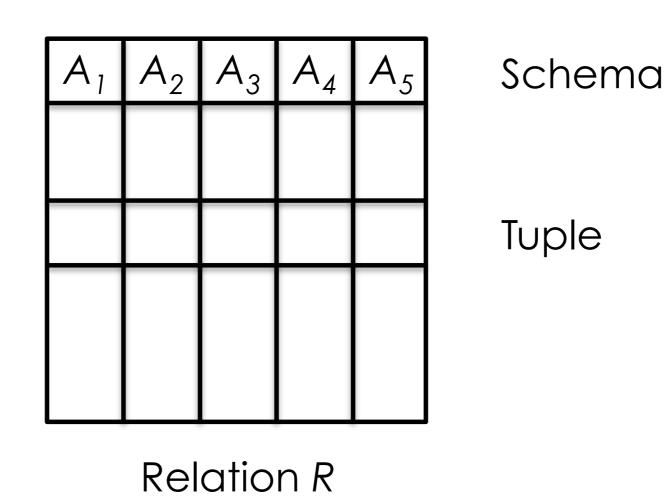


- Divide the vector in equal-sized subvectors that can fit in memory
- According to that, divide the matrix in stripes
- Stripe i and subvector i are independent from other stripes/subvectors
- Use the previous algorithm for each stripe/subvector pair









- •SELECTION: Select from R tuples satisfying condition C
- PROJECTION: For each tuple in R, select only certain attributes
- •UNION, INTERSECTION, DIFFERENCE: Set operations on two relations with same schema
- •NATURAL JOIN
- GROUPING and AGGREGATION







- MAP: Each tuple t, if condition C is satisfied, is outputted as a (t, t) pair
- REDUCE: Identity

MAP: For each tuple t, create a new tuple t' containing only projected attributes. Outpu is (t', t') pair
REDUCE: Coalesce input (t', [t' t' t' t']) in output (t',t')







- •MAP: Each tuple t is outputted as a (t, t) pair
- REDUCE: For each key t, there will be 1 or 2 values t.
 Coalesce them in a single output (t,t)
- •MAP: Each tuple t is outputted as a (t, t) pair
- REDUCE: For each key t, there will be 1 or 2 values t. If 2

values, coalesce them in a single output (t,t), else ignore

- MAP: For each tuple t in R, produce (t, "R"). For each tuple t in S, produce (t, "S").
- REDUCE: For each key t, there will be 1 or 2 values t. If 1 value, and being "R", output (t,t), else ignore







- We have two relations R(A,B) and S(B,C). Find tuples that agree on B components
- MAP: For each tuple (a,b) from R, produce (b,("R",a)).
 For each tuple (b,c) from S, produce (b,("S",c)).
 REDUCE: For each key b, there will a list of values of the form ("R",a) or ("S",c). Construct all pairs and output them with b.







- We have the relation R(A,B,C) and we **group-by** A and **aggregate** on B.
- MAP: For each tuple (a,b,c) from R, output (a,b).
 Each key a represents a group.
- REDUCE: Apply the aggregation operator to the list of b values associate with group a, producing x. Output (a,x).







•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
- Graph algorithms typically 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?







Adjacency Matrix

- Represent a graph as an n x n square matrix M
- n = IVI
- $M_{ij} = 1$ means a link from node i to j
- 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 List

- Take adjacency matrices...
- and throw away all the zeros
- Advantages:
 - Much more compact representation
 - Easy to compute over outlinks
- Disadvantages:
 - Much more difficult to compute over inlinks







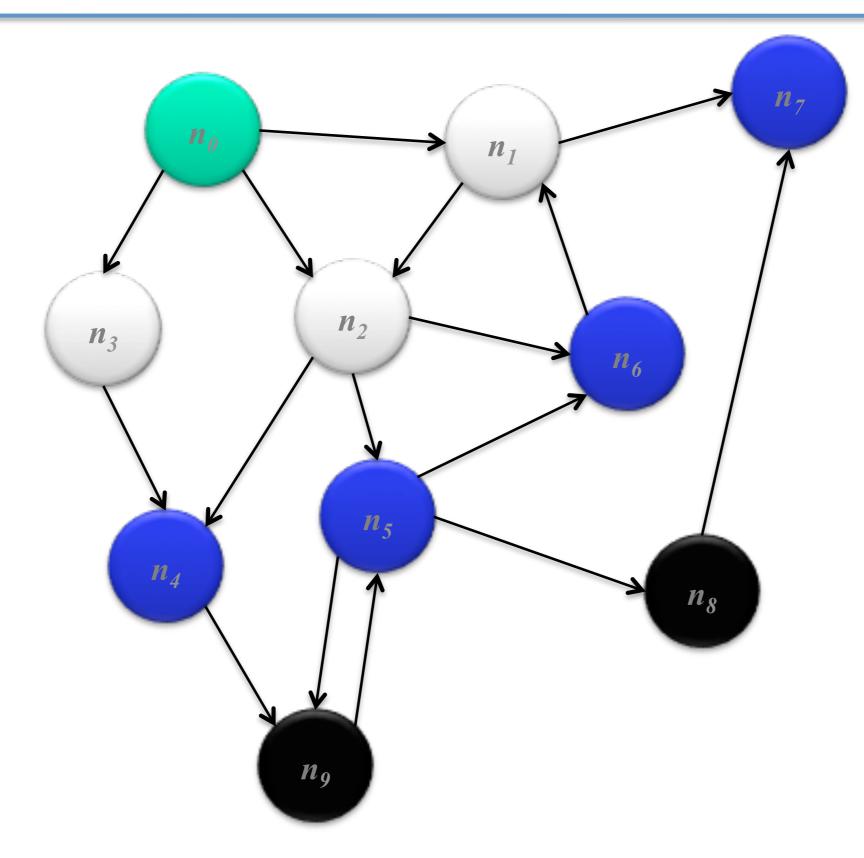
- Consider simple case of equal edge weights
- Solution to the problem can be defined inductively
- Here's the intuition:
 - Define: b is reachable from a if b is on adjacency list of a
 DISTANCETO(s) = 0
 - For all nodes p reachable from s, DISTANCETO(p) = 1
 - For all nodes n reachable from some other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





Shortest Path









Algorithm



• Data representation:

- Key: node n
- Value: d (distance from start), adjacency list (list of nodes reachable from n)
- Initialization: for all nodes except for start node, d = infinity
- Mapper:
 - m Selects minimum distance path for each reachable node
 - Additional bookkeeping needed to keep track of actual path
 - adjacency list: emit (m, d + 1)
- 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







- Each MapReduce iteration advances the "known 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, adjacency list) as well









1:	class MAPPER	
2:	method MAP(nid n , node N)	
3:	$d \leftarrow N.\text{Distance}$	
4:	Eміт(nid n, N)	
5:	for all nodeid $m \in N$. AdjacencyList do	
6:	Еміт(nid $m, d + 1$)	D
1:	class Reducer	
2:	method REDUCE(nid m , $[d_1, d_2, \ldots]$)	
3:	$d_{min} \leftarrow \infty$	
4:	$M \leftarrow \emptyset$	
5:	for all $d \in \text{counts} [d_1, d_2, \ldots]$ do	
6:	if $IsNode(d)$ then	
7:	$M \leftarrow d$	
8:	else if $d < d_{min}$ then	
9:	$d_{min} \leftarrow d$	
10:	$M.Distance \leftarrow d_{min}$	
11:	EмIT(nid m , node M)	

▷ Pass along graph structure

▷ Emit distances to reachable nodes

Recover graph structureLook for shorter distance

▷ Update shortest distance







- Graph algorithms typically 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