

*W4111 – Introduction to Databases  
Section 002, Spring 2024  
Lecture 9  
Module II (2), NoSQL (2)*



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*Lecture 9*  
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We will start in a couple of minutes.

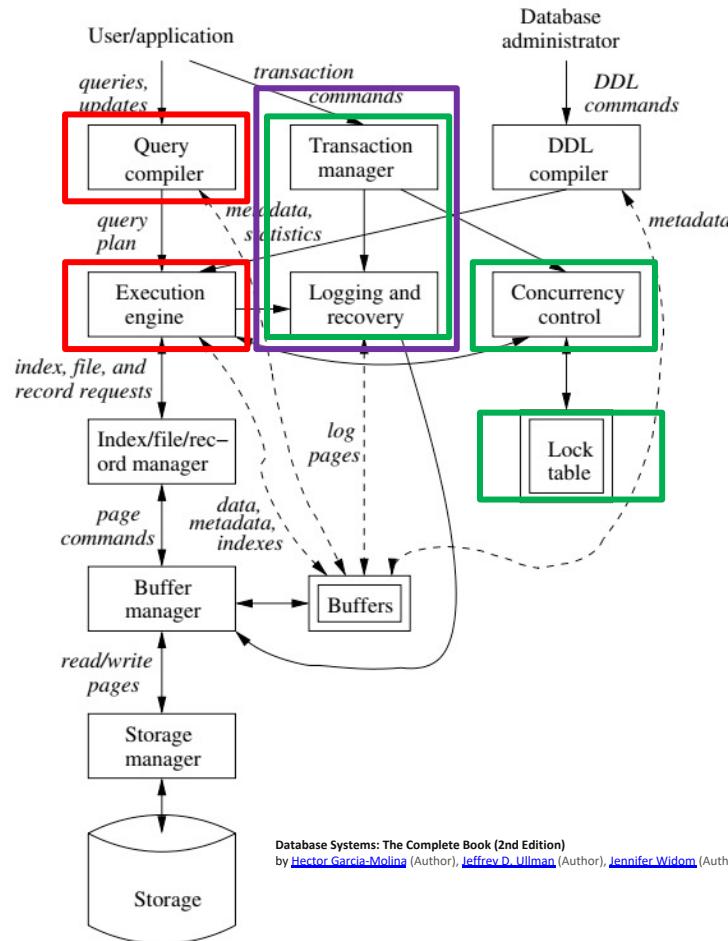
# *Module II, Part 3*

# *Reminder*

# *Query Processing*

# Transactions

- Find things quickly.
- Load/Save quickly.
- Transactions
- Durability



# Query Compilation

## Preview of Query Compilation

Database Systems: The Complete Book (2nd Edition) 2nd Edition  
by [Hector Garcia-Molina](#) (Author), [Jeffrey D. Ullman](#) (Author), [Jennifer Widom](#) (Author)

To set the context for query execution, we offer a very brief outline of the content of the next chapter. Query compilation is divided into the three major steps shown in Fig. 15.2.

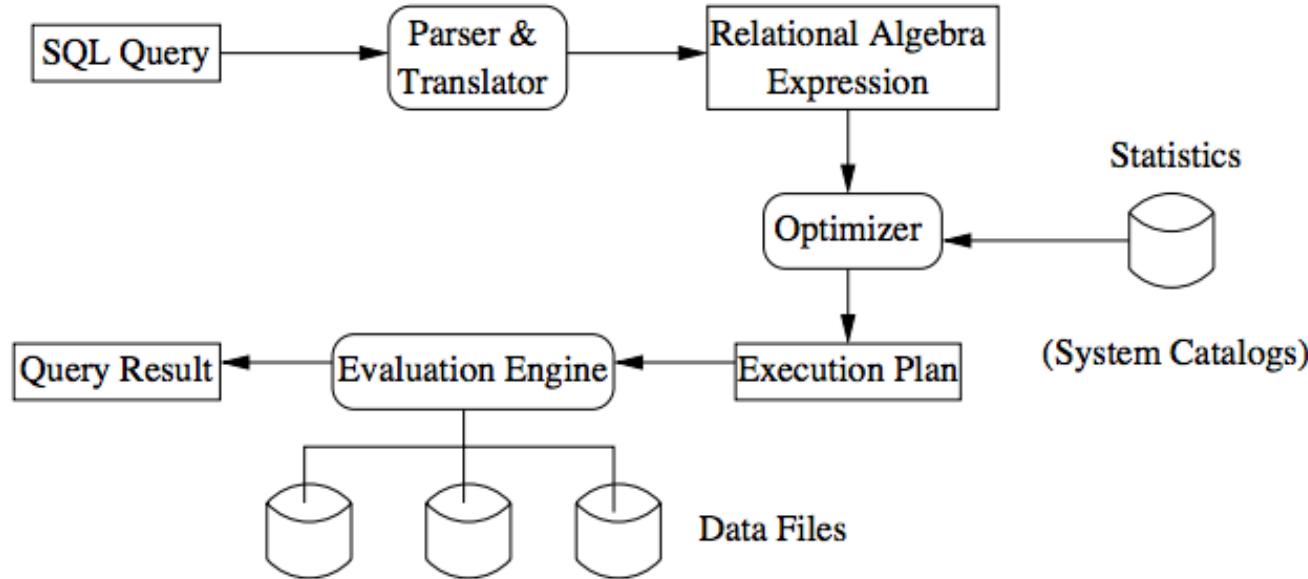
- a) *Parsing.* A *parse tree* for the query is constructed.
- b) *Query Rewrite.* The parse tree is converted to an initial query plan, which is usually an algebraic representation of the query. This initial plan is then transformed into an equivalent plan that is expected to require less time to execute.
- c) *Physical Plan Generation.* The abstract query plan from (b), often called a *logical query plan*, is turned into a *physical query plan* by selecting algorithms to implement each of the operators of the logical plan, and by selecting an order of execution for these operators. The physical plan, like the result of parsing and the logical plan, is represented by an expression tree. The physical plan also includes details such as how the queried relations are accessed, and when and if a relation should be sorted.

# Parsing and Execution

- Parser/Translator
  - Verifies syntax correctness and generates a *parse tree*.
  - Converts to *logical plan tree* that defines how to execute the query.
    - Tree nodes are *operator(tables, parameters)*
    - Edges are the flow of data “up the tree” from node to node.
- Optimizer
  - Modifies the logical plan to define an improved execution.
  - Query rewrite/transformation.
  - Determines *how* to choose among multiple implementations of operators.
- Engine
  - Executes the plan
  - May modify the plan to *optimize* execution, e.g. using indexes.

# Query Processing Overview

## Basic Steps in Processing an SQL Query





# Chapter 15: Query Processing

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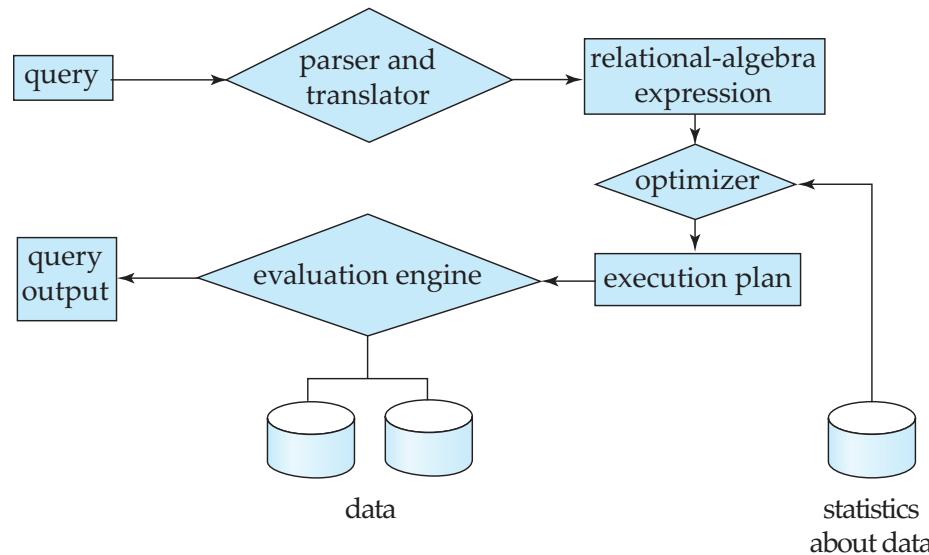
# Chapter 15: Query Processing

- Overview
- Measures of Query Cost
- Selection Operation
- Sorting
- Join Operation
- Other Operations
- Evaluation of Expressions



# Basic Steps in Query Processing

1. Parsing and translation
2. Optimization
3. Evaluation





# Basic Steps in Query Processing (Cont.)

- Parsing and translation
  - translate the query into its internal form. This is then translated into relational algebra.
  - Parser checks syntax, verifies relations
- Evaluation
  - The query-execution engine takes a query-evaluation plan, executes that plan, and returns the answers to the query.



# Basic Steps in Query Processing: Optimization

- A relational algebra expression may have many equivalent expressions
  - E.g.,  $\sigma_{\text{salary} < 75000}(\Pi_{\text{salary}}(\text{instructor}))$  is equivalent to  $\Pi_{\text{salary}}(\sigma_{\text{salary} < 75000}(\text{instructor}))$
- Each relational algebra operation can be evaluated using one of several different algorithms
  - Correspondingly, a relational-algebra expression can be evaluated in many ways.
- Annotated expression specifying detailed evaluation strategy is called an **evaluation-plan**. E.g.,:
  - Use an index on *salary* to find instructors with salary < 75000,
  - Or perform complete relation scan and discard instructors with salary  $\geq 75000$



# Basic Steps: Optimization (Cont.)

- **Query Optimization:** Amongst all equivalent evaluation plans choose the one with lowest cost.
  - Cost is estimated using statistical information from the database catalog
    - e.g.. number of tuples in each relation, size of tuples, etc.
- In this chapter we study
  - How to measure query costs
  - Algorithms for evaluating relational algebra operations
  - How to combine algorithms for individual operations in order to evaluate a complete expression
- In Chapter 16
  - We study how to optimize queries, that is, how to find an evaluation plan with lowest estimated cost



# Measures of Query Cost

- Many factors contribute to time cost
  - *disk access, CPU, and network communication*
- Cost can be measured based on
  - **response time**, i.e. total elapsed time for answering query, or
  - total **resource consumption**
- We use total resource consumption as cost metric
  - Response time harder to estimate, and minimizing resource consumption is a good idea in a shared database
- We ignore CPU costs for simplicity
  - Real systems do take CPU cost into account
  - Network costs must be considered for parallel systems
- We describe how estimate the cost of each operation
  - We do not include cost to writing output to disk



# Measures of Query Cost

- Disk cost can be estimated as:
  - Number of seeks \* average-seek-cost
  - Number of blocks read \* average-block-read-cost
  - Number of blocks written \* average-block-write-cost
- For simplicity we just use the **number of block transfers from disk and the number of seeks** as the cost measures
  - $t_T$  – time to transfer one block
    - Assuming for simplicity that write cost is same as read cost
  - $t_S$  – time for one seek
  - Cost for  $b$  block transfers plus  $S$  seeks
$$b * t_T + S * t_S$$
- $t_S$  and  $t_T$  depend on where data is stored; with 4 KB blocks:
  - High end magnetic disk:  $t_S = 4$  msec and  $t_T = 0.1$  msec
  - SSD:  $t_S = 20\text{-}90$  microsec and  $t_T = 2\text{-}10$  microsec for 4KB



# Measures of Query Cost (Cont.)

- Required data may be buffer resident already, avoiding disk I/O
  - But hard to take into account for cost estimation
- Several algorithms can reduce disk IO by using extra buffer space
  - Amount of real memory available to buffer depends on other concurrent queries and OS processes, known only during execution
- Worst case estimates assume that no data is initially in buffer and only the minimum amount of memory needed for the operation is available
  - But more optimistic estimates are used in practice



# Selection Operation

- **File scan**
- Algorithm **A1 (linear search)**. Scan each file block and test all records to see whether they satisfy the selection condition.
  - Cost estimate =  $b_r$  block transfers + 1 seek
    - $b_r$  denotes number of blocks containing records from relation  $r$
  - If selection is on a key attribute, can stop on finding record
    - cost =  $(b_r/2)$  block transfers + 1 seek
  - Linear search can be applied regardless of
    - selection condition or
    - ordering of records in the file, or
    - availability of indices
- Note: binary search generally does not make sense since data is not stored consecutively
  - except when there is an index available,
  - and binary search requires more seeks than index search



# Selections Using Indices

- **Index scan** – search algorithms that use an index
  - selection condition must be on search-key of index.
- **A2 (clustering index, equality on key)**. Retrieve a single record that satisfies the corresponding equality condition
  - $Cost = (h_i + 1) * (t_T + t_S)$
- **A3 (clustering index, equality on nonkey)** Retrieve multiple records.
  - Records will be on consecutive blocks
    - Let b = number of blocks containing matching records
  - $Cost = h_i * (t_T + t_S) + t_S + t_T * b$



# Selections Using Indices

- A4 (secondary index, equality on key/non-key).
  - Retrieve a single record if the search-key is a candidate key
    - $Cost = (h_i + 1) * (t_T + t_S)$
  - Retrieve multiple records if search-key is not a candidate key
    - each of  $n$  matching records may be on a different block
    - $Cost = (h_i + n) * (t_T + t_S)$ 
      - Can be very expensive!



# Selections Involving Comparisons

- Can implement selections of the form  $\sigma_{A \leq v}(r)$  or  $\sigma_{A \geq v}(r)$  by using
  - a linear file scan,
  - or by using indices in the following ways:
- **A5 (clustering index, comparison).** (Relation is sorted on A)
  - For  $\sigma_{A \geq v}(r)$  use index to find first tuple  $\geq v$  and scan relation sequentially from there
  - For  $\sigma_{A \leq v}(r)$  just scan relation sequentially till first tuple  $> v$ ; do not use index
- **A6 (clustering index, comparison).**
  - For  $\sigma_{A \geq v}(r)$  use index to find first index entry  $\geq v$  and scan index sequentially from there, to find pointers to records.
  - For  $\sigma_{A \leq v}(r)$  just scan leaf pages of index finding pointers to records, till first entry  $> v$
  - In either case, retrieve records that are pointed to
  - requires an I/O per record; Linear file scan may be cheaper!



# Implementation of Complex Selections

- **Conjunction:**  $\sigma_{\theta_1 \wedge \theta_2 \wedge \dots \wedge \theta_n}(r)$
- **A7 (conjunctive selection using one index).**
  - Select a combination of  $\theta_i$  and algorithms A1 through A7 that results in the least cost for  $\sigma_{\theta_i}(r)$ .
  - Test other conditions on tuple after fetching it into memory buffer.
- **A8 (conjunctive selection using composite index).**
  - Use appropriate composite (multiple-key) index if available.
- **A9 (conjunctive selection by intersection of identifiers).**
  - Requires indices with record pointers.
  - Use corresponding index for each condition, and take intersection of all the obtained sets of record pointers.
  - Then fetch records from file
  - If some conditions do not have appropriate indices, apply test in memory.



# Algorithms for Complex Selections

- **Disjunction:**  $\sigma_{\theta_1 \vee \theta_2 \vee \dots \vee \theta_n}(r)$ .
- **A10 (disjunctive selection by union of identifiers).**
  - Applicable if *all* conditions have available indices.
    - Otherwise use linear scan.
  - Use corresponding index for each condition, and take union of all the obtained sets of record pointers.
  - Then fetch records from file
- **Negation:**  $\sigma_{\neg\theta}(r)$ 
  - Use linear scan on file
  - If very few records satisfy  $\neg\theta$ , and an index is applicable to  $\theta$ 
    - Find satisfying records using index and fetch from file



# Sorting

- We may build an index on the relation, and then use the index to read the relation in sorted order. May lead to one disk block access for each tuple.
- For relations that fit in memory, techniques like quicksort can be used.
  - For relations that don't fit in memory, **external sort-merge** is a good choice.



# Join Operation

- Several different algorithms to implement joins
  - Nested-loop join
  - Block nested-loop join
  - Indexed nested-loop join
  - Merge-join
  - Hash-join
- Choice based on cost estimate
- Examples use the following information
  - Number of records of *student*: 5,000    *takes*: 10,000
  - Number of blocks of *student*: 100    *takes*: 400



# Nested-Loop Join

- To compute the theta join  $r \bowtie_{\theta} s$   
**for each tuple  $t_r$  in  $r$  do begin**  
    **for each tuple  $t_s$  in  $s$  do begin**  
        test pair  $(t_r, t_s)$  to see if they satisfy the join condition  $\theta$   
        if they do, add  $t_r \cdot t_s$  to the result.  
    **end**  
**end**
- $r$  is called the **outer relation** and  $s$  the **inner relation** of the join.
- Requires no indices and can be used with any kind of join condition.
- Expensive since it examines every pair of tuples in the two relations.



# Nested-Loop Join (Cont.)

- In the worst case, if there is enough memory only to hold one block of each relation, the estimated cost is
$$n_r * b_s + b_r \text{ block transfers, plus } n_r + b_r \text{ seeks}$$
- If the smaller relation fits entirely in memory, use that as the inner relation.
  - Reduces cost to  $b_r + b_s$  block transfers and 2 seeks
- Assuming worst case memory availability cost estimate is
  - with *student* as outer relation:
    - $5000 * 400 + 100 = 2,000,100$  block transfers,
    - $5000 + 100 = 5100$  seeks
  - with *takes* as the outer relation
    - $10000 * 100 + 400 = 1,000,400$  block transfers and 10,400 seeks
- If smaller relation (*student*) fits entirely in memory, the cost estimate will be 500 block transfers.
- Block nested-loops algorithm (next slide) is preferable.



# Indexed Nested-Loop Join

- Index lookups can replace file scans if
  - join is an equi-join or natural join and
  - an index is available on the inner relation's join attribute
    - Can construct an index just to compute a join.
- For each tuple  $t_r$  in the outer relation  $r$ , use the index to look up tuples in  $s$  that satisfy the join condition with tuple  $t_r$ .
- Worst case: buffer has space for only one page of  $r$ , and, for each tuple in  $r$ , we perform an index lookup on  $s$ .
- Cost of the join:  $b_r(t_T + t_S) + n_r * c$ 
  - Where  $c$  is the cost of traversing index and fetching all matching  $s$  tuples for one tuple of  $r$
  - $c$  can be estimated as cost of a single selection on  $s$  using the join condition.
- If indices are available on join attributes of both  $r$  and  $s$ , use the relation with fewer tuples as the outer relation.



# Merge-Join

1. Sort both relations on their join attribute (if not already sorted on the join attributes).
2. Merge the sorted relations to join them
  1. Join step is similar to the merge stage of the sort-merge algorithm.
  2. Main difference is handling of duplicate values in join attribute — every pair with same value on join attribute must be matched
  3. Detailed algorithm in book

	$a1$	$a2$
$pr \rightarrow$	a	3
	b	1
	d	8
	d	13
	f	7
	m	5
	q	6

$r$

	$a1$	$a3$
$ps \rightarrow$	a	A
	b	G
	c	L
	d	N
	m	B

$s$

$$\#(A) = n, \#(B) = m$$

$$n * m$$

$$n * \log(n) + m * \log(m) + n + m$$



## Merge-Join (Cont.)

- Can be used only for equi-joins and natural joins
- Each block needs to be read only once (assuming all tuples for any given value of the join attributes fit in memory)
- Thus the cost of merge join is:  
 $b_r + b_s$  block transfers +  $\lceil b_r/b_b \rceil + \lceil b_s/b_b \rceil$  seeks  
+ the cost of sorting if relations are unsorted.
- **hybrid merge-join:** If one relation is sorted, and the other has a secondary B<sup>+</sup>-tree index on the join attribute
  - Merge the sorted relation with the leaf entries of the B<sup>+</sup>-tree .
  - Sort the result on the addresses of the unsorted relation's tuples
  - Scan the unsorted relation in physical address order and merge with previous result, to replace addresses by the actual tuples
    - Sequential scan more efficient than random lookup



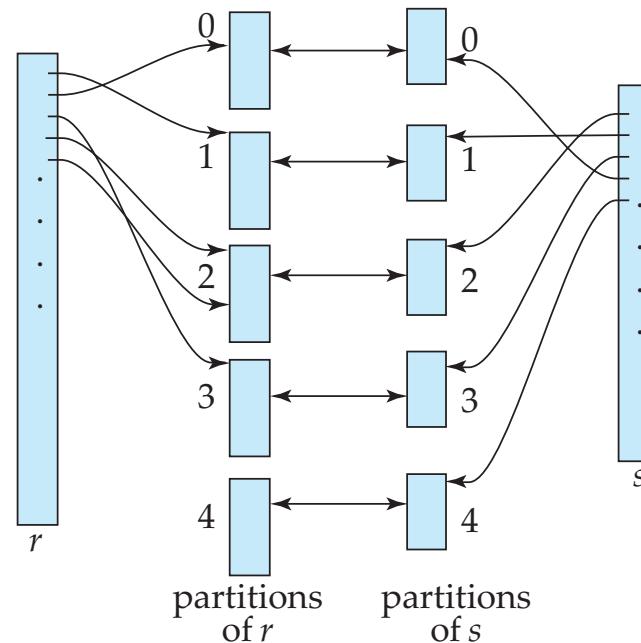
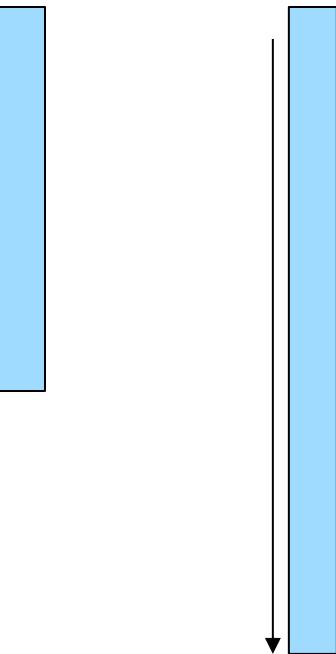
# Hash-Join

- Applicable for equi-joins and natural joins.
- A hash function  $h$  is used to partition tuples of both relations
- $h$  maps  $JoinAttrs$  values to  $\{0, 1, \dots, n\}$ , where  $JoinAttrs$  denotes the common attributes of  $r$  and  $s$  used in the natural join.
  - $r_0, r_1, \dots, r_n$  denote partitions of  $r$  tuples
    - Each tuple  $t_r \in r$  is put in partition  $r_i$  where  $i = h(t_r[JoinAttrs])$ .
  - $r_0, r_1, \dots, r_n$  denotes partitions of  $s$  tuples
    - Each tuple  $t_s \in s$  is put in partition  $s_i$ , where  $i = h(t_s[JoinAttrs])$ .
- Note: In book, Figure 12.10  $r_i$  is denoted as  $H_{ri}$ ,  $s_i$  is denoted as  $H_{si}$  and  $n$  is denoted as  $n_h$ .



## Hash-Join (Cont.)

On  $a.X=b.Z$







# Hash-Join Algorithm

The hash-join of  $r$  and  $s$  is computed as follows.

1. Partition the relation  $s$  using hashing function  $h$ . When partitioning a relation, one block of memory is reserved as the output buffer for each partition.
2. Partition  $r$  similarly.
3. For each  $i$ :
  - (a) Load  $s_i$  into memory and build an in-memory hash index on it using the join attribute. This hash index uses a different hash function than the earlier one  $h$ .
  - (b) Read the tuples in  $r_i$  from the disk one by one. For each tuple  $t_r$ , locate each matching tuple  $t_s$  in  $s_i$  using the in-memory hash index. Output the concatenation of their attributes.

Relation  $s$  is called the **build input** and  $r$  is called the **probe input**.



## Hash-Join algorithm (Cont.)

- The value  $n$  and the hash function  $h$  is chosen such that each  $s_i$  should fit in memory.
  - Typically  $n$  is chosen as  $\lceil b_s/M \rceil * f$  where  $f$  is a “**fudge factor**”, typically around 1.2
  - The probe relation partitions  $s_i$  need not fit in memory
- **Recursive partitioning** required if number of partitions  $n$  is greater than number of pages  $M$  of memory.
  - instead of partitioning  $n$  ways, use  $M - 1$  partitions for  $s$
  - Further partition the  $M - 1$  partitions using a different hash function
  - Use same partitioning method on  $r$
  - Rarely required: e.g., with block size of 4 KB, recursive partitioning not needed for relations of < 1GB with memory size of 2MB, or relations of < 36 GB with memory of 12 MB



# Other Operations

- **Duplicate elimination** can be implemented via hashing or sorting.
  - On sorting duplicates will come adjacent to each other, and all but one set of duplicates can be deleted.
  - *Optimization:* duplicates can be deleted during run generation as well as at intermediate merge steps in external sort-merge.
  - Hashing is similar – duplicates will come into the same bucket.
- **Projection:**
  - perform projection on each tuple
  - followed by duplicate elimination.



# Other Operations : Aggregation

- **Aggregation** can be implemented in a manner similar to duplicate elimination.
  - **Sorting** or **hashing** can be used to bring tuples in the same group together, and then the aggregate functions can be applied on each group.
  - Optimization: **partial aggregation**
    - combine tuples in the same group during run generation and intermediate merges, by computing partial aggregate values
    - For count, min, max, sum: keep aggregate values on tuples found so far in the group.
      - When combining partial aggregate for count, add up the partial aggregates
    - For avg, keep sum and count, and divide sum by count at the end



# Evaluation of Expressions

- So far: we have seen algorithms for individual operations
- Alternatives for evaluating an entire expression tree
  - **Materialization:** generate results of an expression whose inputs are relations or are already computed, **materialize** (store) it on disk. Repeat.
  - **Pipelining:** pass on tuples to parent operations even as an operation is being executed
- We study above alternatives in more detail

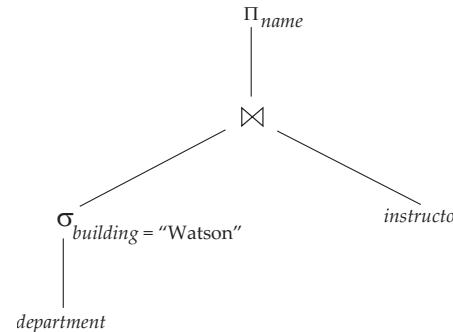


# Materialization

- **Materialized evaluation:** evaluate one operation at a time, starting at the lowest-level. Use intermediate results materialized into temporary relations to evaluate next-level operations.
- E.g., in figure below, compute and store

$$\sigma_{building="Watson"}(department)$$

then compute the store its join with *instructor*, and finally compute the projection on *name*.





# Materialization (Cont.)

- Materialized evaluation is always applicable
- Cost of writing results to disk and reading them back can be quite high
  - Our cost formulas for operations ignore cost of writing results to disk, so
    - Overall cost = Sum of costs of individual operations + cost of writing intermediate results to disk
- **Double buffering:** use two output buffers for each operation, when one is full write it to disk while the other is getting filled
  - Allows overlap of disk writes with computation and reduces execution time



# Pipelining

- **Pipelined evaluation:** evaluate several operations simultaneously, passing the results of one operation on to the next.
- E.g., in previous expression tree, don't store result of
$$\sigma_{building = "Watson"}(department)$$
  - instead, pass tuples directly to the join.. Similarly, don't store result of join, pass tuples directly to projection.
- Much cheaper than materialization: no need to store a temporary relation to disk.
- Pipelining may not always be possible – e.g., sort, hash-join.
- For pipelining to be effective, use evaluation algorithms that generate output tuples even as tuples are received for inputs to the operation.
- Pipelines can be executed in two ways: **demand driven** and **producer driven**



# Pipelining (Cont.)

- In **demand driven** or **lazy** evaluation
  - system repeatedly requests next tuple from top level operation
  - Each operation requests next tuple from children operations as required, in order to output its next tuple
  - In between calls, operation has to maintain “**state**” so it knows what to return next
- In **producer-driven** or **eager** pipelining
  - Operators produce tuples eagerly and pass them up to their parents
    - Buffer maintained between operators, child puts tuples in buffer, parent removes tuples from buffer
    - if buffer is full, child waits till there is space in the buffer, and then generates more tuples
  - System schedules operations that have space in output buffer and can process more input tuples
- Alternative name: **pull** and **push** models of pipelining



# Pipelining (Cont.)

- Implementation of demand-driven pipelining
  - Each operation is implemented as an **iterator** implementing the following operations
    - **open()**
      - E.g., file scan: initialize file scan
        - state: pointer to beginning of file
      - E.g., merge join: sort relations;
        - state: pointers to beginning of sorted relations
    - **next()**
      - E.g., for file scan: Output next tuple, and advance and store file pointer
      - E.g., for merge join: continue with merge from earlier state till next output tuple is found. Save pointers as iterator state.
    - **close()**

# *Query Optimization*



# Chapter 16: Query Optimization

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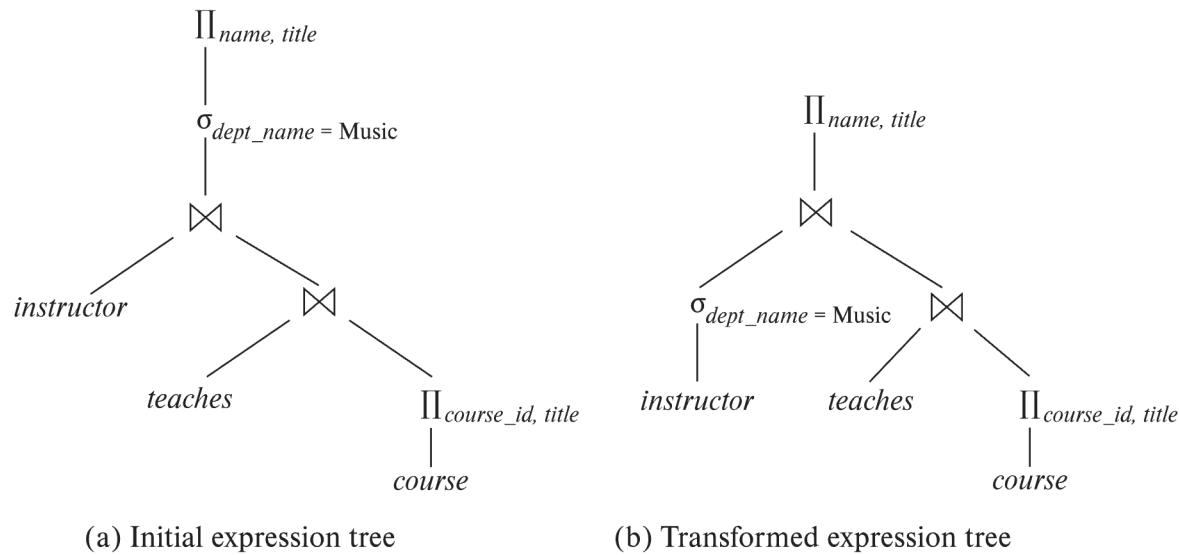
# Outline

- Introduction
- Transformation of Relational Expressions
- Catalog Information for Cost Estimation
- Statistical Information for Cost Estimation
- Cost-based optimization
- Dynamic Programming for Choosing Evaluation Plans
- Materialized views



# Introduction

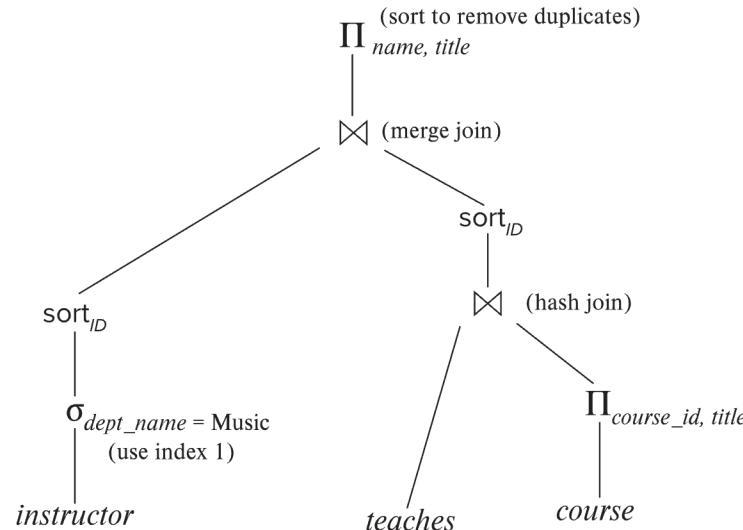
- Alternative ways of evaluating a given query
  - Equivalent expressions
  - Different algorithms for each operation





# Introduction (Cont.)

- An **evaluation plan** defines exactly what algorithm is used for each operation, and how the execution of the operations is coordinated.



- Find out how to view query execution plans on your favorite database



# Introduction (Cont.)

- Cost difference between evaluation plans for a query can be enormous
  - E.g., seconds vs. days in some cases
- Steps in **cost-based query optimization**
  1. Generate logically equivalent expressions using **equivalence rules**
  2. Annotate resultant expressions to get alternative query plans
  3. Choose the cheapest plan based on **estimated cost**
- Estimation of plan cost based on:
  - Statistical information about relations. Examples:
    - number of tuples, number of distinct values for an attribute
  - Statistics estimation for intermediate results
    - to compute cost of complex expressions
  - Cost formulae for algorithms, computed using statistics



# Viewing Query Evaluation Plans

- Most database support **explain <query>**
  - Displays plan chosen by query optimizer, along with cost estimates
  - Some syntax variations between databases
    - Oracle: **explain plan for <query>** followed by **select \* from table (dbms\_xplan.display)**
    - SQL Server: **set showplan\_text on**
- Some databases (e.g. PostgreSQL) support **explain analyse <query>**
  - Shows actual runtime statistics found by running the query, in addition to showing the plan
- Some databases (e.g. PostgreSQL) show cost as **f./l**
  - $f$  is the cost of delivering first tuple and  $l$  is cost of delivering all results



# Generating Equivalent Expressions

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# Transformation of Relational Expressions

- Two relational algebra expressions are said to be **equivalent** if the two expressions generate the same set of tuples on every *legal* database instance
  - Note: order of tuples is irrelevant
  - we don't care if they generate different results on databases that violate integrity constraints
- In SQL, inputs and outputs are multisets of tuples
  - Two expressions in the multiset version of the relational algebra are said to be equivalent if the two expressions generate the same multiset of tuples on every legal database instance.
- An **equivalence rule** says that expressions of two forms are equivalent
  - Can replace expression of first form by second, or vice versa



# Equivalence Rules

1. Conjunctive selection operations can be deconstructed into a sequence of individual selections.

$$\sigma_{\theta_1 \wedge \theta_2}(E) \equiv \sigma_{\theta_1}(\sigma_{\theta_2}(E))$$

2. Selection operations are commutative.

$$\sigma_{\theta_1}(\sigma_{\theta_2}(E)) \equiv \sigma_{\theta_2}(\sigma_{\theta_1}(E))$$

3. Only the last in a sequence of projection operations is needed, the others can be omitted.

$$\Pi_{L_1}(\Pi_{L_2}(\dots(\Pi_{L_n}(E))\dots)) \equiv \Pi_{L_1}(E)$$

where  $L_1 \subseteq L_2 \dots \subseteq L_n$

4. Selections can be combined with Cartesian products and theta joins.

a.  $\sigma_{\theta}(E_1 \times E_2) \equiv E_1 \bowtie_{\theta} E_2$

b.  $\sigma_{\theta_1}(E_1 \bowtie_{\theta_2} E_2) \equiv E_1 \bowtie_{\theta_1 \wedge \theta_2} E_2$



## Equivalence Rules (Cont.)

5. Theta-join operations (and natural joins) are commutative.

$$E_1 \bowtie E_2 \equiv E_2 \bowtie E_1$$

6. (a) Natural join operations are associative:

$$(E_1 \bowtie E_2) \bowtie E_3 \equiv E_1 \bowtie (E_2 \bowtie E_3)$$

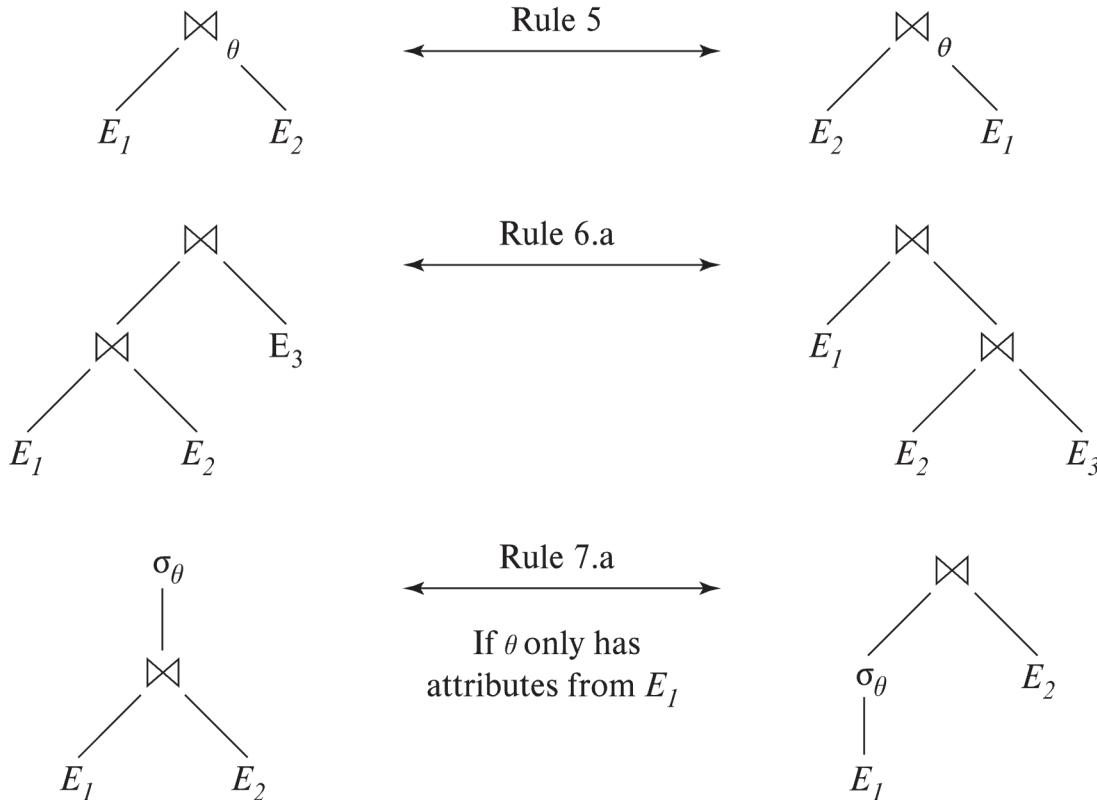
- (b) Theta joins are associative in the following manner:

$$(E_1 \bowtie_{\theta_1} E_2) \bowtie_{\theta_2 \wedge \theta_3} E_3 \equiv E_1 \bowtie_{\theta_1 \wedge \theta_3} (E_2 \bowtie_{\theta_2} E_3)$$

where  $\theta_2$  involves attributes from only  $E_2$  and  $E_3$ .



## Pictorial Depiction of Equivalence Rules





## Equivalence Rules (Cont.)

7. The selection operation distributes over the theta join operation under the following two conditions:
  - (a) When all the attributes in  $\theta_0$  involve only the attributes of one of the expressions ( $E_1$ ) being joined.

$$\sigma_{\theta_0}(E_1 \bowtie_{\theta} E_2) \equiv (\sigma_{\theta_0}(E_1)) \bowtie_{\theta} E_2$$

- (b) When  $\theta_1$  involves only the attributes of  $E_1$  and  $\theta_2$  involves only the attributes of  $E_2$ .

$$\sigma_{\theta_1 \wedge \theta_2}(E_1 \bowtie_{\theta} E_2) \equiv (\sigma_{\theta_1}(E_1)) \bowtie_{\theta} (\sigma_{\theta_2}(E_2))$$



## Equivalence Rules (Cont.)

8. The projection operation distributes over the theta join operation as follows:

(a) if  $\theta$  involves only attributes from  $L_1 \cup L_2$ :

$$\Pi_{L_1 \cup L_2}(E_1 \bowtie_{\theta} E_2) \equiv \Pi_{L_1}(E_1) \bowtie_{\theta} \Pi_{L_2}(E_2)$$

(b) In general, consider a join  $E_1 \bowtie_{\theta} E_2$ .

- Let  $L_1$  and  $L_2$  be sets of attributes from  $E_1$  and  $E_2$ , respectively.
- Let  $L_3$  be attributes of  $E_1$  that are involved in join condition  $\theta$ , but are not in  $L_1 \cup L_2$ , and
- let  $L_4$  be attributes of  $E_2$  that are involved in join condition  $\theta$ , but are not in  $L_1 \cup L_2$ .

$$\Pi_{L_1 \cup L_2}(E_1 \bowtie_{\theta} E_2) \equiv \Pi_{L_1 \cup L_2}(\Pi_{L_1 \cup L_3}(E_1) \bowtie_{\theta} \Pi_{L_2 \cup L_4}(E_2))$$

Similar equivalences hold for outerjoin operations:  $\bowtie$ ,  $\bowtie_l$ , and  $\bowtie_r$



## Equivalence Rules (Cont.)

9. The set operations union and intersection are commutative

$$E_1 \cup E_2 \equiv E_2 \cup E_1$$

$$E_1 \cap E_2 \equiv E_2 \cap E_1$$

(set difference is not commutative).

10. Set union and intersection are associative.

$$(E_1 \cup E_2) \cup E_3 \equiv E_1 \cup (E_2 \cup E_3)$$

$$(E_1 \cap E_2) \cap E_3 \equiv E_1 \cap (E_2 \cap E_3)$$

11. The selection operation distributes over  $\cup$ ,  $\cap$  and  $-$ .

a.  $\sigma_{\theta}(E_1 \cup E_2) \equiv \sigma_{\theta}(E_1) \cup \sigma_{\theta}(E_2)$

b.  $\sigma_{\theta}(E_1 \cap E_2) \equiv \sigma_{\theta}(E_1) \cap \sigma_{\theta}(E_2)$

c.  $\sigma_{\theta}(E_1 - E_2) \equiv \sigma_{\theta}(E_1) - \sigma_{\theta}(E_2)$

d.  $\sigma_{\theta}(E_1 \cap E_2) \equiv \sigma_{\theta}(E_1) \cap E_2$

e.  $\sigma_{\theta}(E_1 - E_2) \equiv \sigma_{\theta}(E_1) - E_2$

preceding equivalence does not hold for  $\cup$

12. The projection operation distributes over union

$$\Pi_L(E_1 \cup E_2) \equiv (\Pi_L(E_1)) \cup (\Pi_L(E_2))$$



## Equivalence Rules (Cont.)

13. Selection distributes over aggregation as below

$$\sigma_{\theta}(G\gamma_A(E)) \equiv G\gamma_A(\sigma_{\theta}(E))$$

provided  $\theta$  only involves attributes in  $G$

14. a. Full outerjoin is commutative:

$$E_1 \bowtie E_2 \equiv E_2 \bowtie E_1$$

- b. Left and right outerjoin are not commutative, but:

$$E_1 \bowtie L E_2 \equiv E_2 \bowtie R E_1$$

15. Selection distributes over left and right outerjoins as below, provided  $\theta_1$  only involves attributes of  $E_1$

a.  $\sigma_{\theta_1}(E_1 \bowtie_0 E_2) \equiv (\sigma_{\theta_1}(E_1)) \bowtie_0 E_2$

b.  $\sigma_{\theta_1}(E_1 \bowtie_0 E_2) \equiv E_2 \bowtie_0 (\sigma_{\theta_1}(E_1))$

16. Outerjoins can be replaced by inner joins under some conditions

a.  $\sigma_{\theta_1}(E_1 \bowtie_0 E_2) \equiv \sigma_{\theta_1}(E_1 \bowtie_0 E_2)$

b.  $\sigma_{\theta_1}(E_1 \bowtie_0 E_2) \equiv \sigma_{\theta_1}(E_1 \bowtie_0 E_2)$

provided  $\theta_1$  is null rejecting on  $E_2$



## Equivalence Rules (Cont.)

Note that several equivalences that hold for joins do not hold for outerjoins

- $\sigma_{\text{year}=2017}(\text{instructor} \bowtie \text{teaches}) \not\equiv \sigma_{\text{year}=2017}(\text{instructor} \bowtie \text{teaches})$
- Outerjoins are not associative  
 $(r \bowtie s) \bowtie t \not\equiv r \bowtie (s \bowtie t)$ 
  - e.g. with  $r(A,B) = \{(1,1)\}$ ,  $s(B,C) = \{(1,1)\}$ ,  $t(A,C) = \{\}$

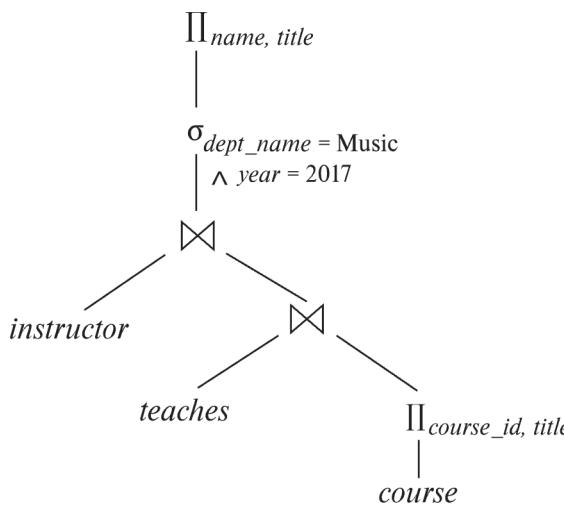


# Transformation Example: Pushing Selections

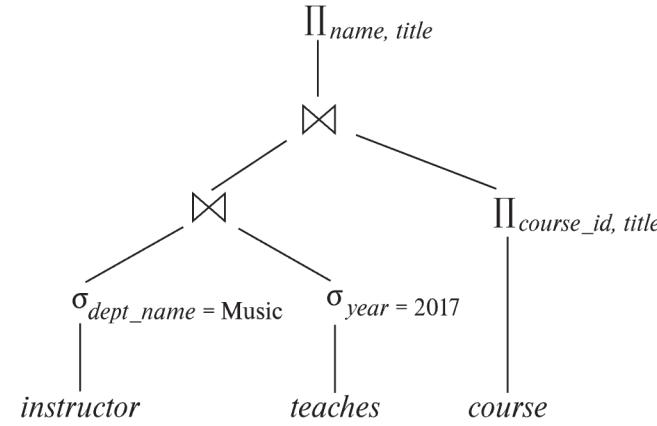
- Query: Find the names of all instructors in the Music department, along with the titles of the courses that they teach
  - $\Pi_{name, title}(\sigma_{dept\_name = \text{Music}}(instructor \bowtie (teaches \bowtie \Pi_{course\_id, title}(course))))$
- Transformation using rule 7a.
  - $\Pi_{name, title}((\sigma_{dept\_name = \text{Music}}(instructor)) \bowtie (teaches \bowtie \Pi_{course\_id, title}(course)))$
- Performing the selection as early as possible reduces the size of the relation to be joined.



# Multiple Transformations (Cont.)



(a) Initial expression tree



(b) Tree after multiple transformations



# Join Ordering Example

- For all relations  $r_1, r_2$ , and  $r_3$ ,  
$$(r_1 \bowtie r_2) \bowtie r_3 = r_1 \bowtie (r_2 \bowtie r_3)$$
(Join Associativity)  $\bowtie$
- If  $r_2 \bowtie r_3$  is quite large and  $r_1 \bowtie r_2$  is small, we choose

$$(r_1 \bowtie r_2) \bowtie r_3$$

so that we compute and store a smaller temporary relation.



## Join Ordering Example (Cont.)

- Consider the expression

$$\begin{aligned}\Pi_{name, title}(\sigma_{dept\_name= \text{'Music'}}(instructor) \bowtie teaches \\ \bowtie \Pi_{course\_id, title}(course)))\end{aligned}$$

- Could compute  $teaches \bowtie \Pi_{course\_id, title}(course)$  first, and join result with  
 $\sigma_{dept\_name= \text{'Music'}}(instructor)$   
but the result of the first join is likely to be a large relation.
- Only a small fraction of the university's instructors are likely to be from the Music department
  - it is better to compute

$$\sigma_{dept\_name= \text{'Music'}}(instructor) \bowtie teaches$$

first.



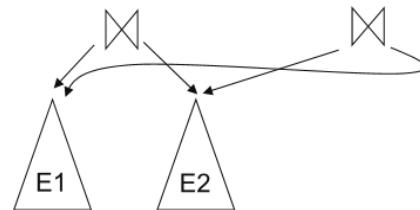
# Enumeration of Equivalent Expressions

- Query optimizers use equivalence rules to **systematically** generate expressions equivalent to the given expression
- Can generate all equivalent expressions as follows:
  - Repeat
    - apply all applicable equivalence rules on every subexpression of every equivalent expression found so far
    - add newly generated expressions to the set of equivalent expressions
- Until no new equivalent expressions are generated above
- The above approach is very expensive in space and time
  - Two approaches
    - Optimized plan generation based on transformation rules
    - Special case approach for queries with only selections, projections and joins



# Implementing Transformation Based Optimization

- Space requirements reduced by sharing common sub-expressions:
  - when E1 is generated from E2 by an equivalence rule, usually only the top level of the two are different, subtrees below are the same and can be shared using pointers
    - E.g., when applying join commutativity



- Same sub-expression may get generated multiple times
  - Detect duplicate sub-expressions and share one copy
- Time requirements are reduced by not generating all expressions
  - Dynamic programming
    - We will study only the special case of dynamic programming for join order optimization



# Cost Estimation

- Cost of each operator computer as described in Chapter 15
  - Need statistics of input relations
    - E.g., number of tuples, sizes of tuples
- Inputs can be results of sub-expressions
  - Need to estimate statistics of expression results
  - To do so, we require additional statistics
    - E.g., number of distinct values for an attribute
- More on cost estimation later



# Choice of Evaluation Plans

- Must consider the interaction of evaluation techniques when choosing evaluation plans
  - choosing the cheapest algorithm for each operation independently may not yield best overall algorithm. E.g.
    - merge-join may be costlier than hash-join, but may provide a sorted output which reduces the cost for an outer level aggregation.
    - nested-loop join may provide opportunity for pipelining
- Practical query optimizers incorporate elements of the following two broad approaches:
  1. Search all the plans and choose the best plan in a cost-based fashion.
  2. Uses heuristics to choose a plan.



# Cost-Based Optimization

- Consider finding the best join-order for  $r_1 \bowtie r_2 \bowtie \dots \bowtie r_n$ .
- There are  $(2(n - 1))!/(n - 1)!$  different join orders for above expression. With  $n = 7$ , the number is 665280, with  $n = 10$ , the number is greater than 176 billion!
- No need to generate all the join orders. Using dynamic programming, the least-cost join order for any subset of  $\{r_1, r_2, \dots, r_n\}$  is computed only once and stored for future use.



# Statistics for Cost Estimation

Database System Concepts, 7<sup>th</sup> Ed.

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See [www.db-book.com](http://www.db-book.com) for conditions on re-use



# Statistical Information for Cost Estimation

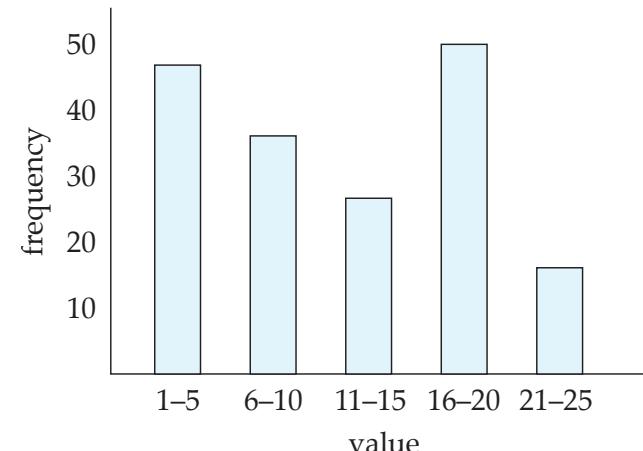
- $n_r$ : number of tuples in a relation  $r$ .
- $b_r$ : number of blocks containing tuples of  $r$ .
- $l_r$ : size of a tuple of  $r$ .
- $f_r$ : blocking factor of  $r$  — i.e., the number of tuples of  $r$  that fit into one block.
- $V(A, r)$ : number of distinct values that appear in  $r$  for attribute  $A$ ; same as the size of  $\Pi_A(r)$ .
- If tuples of  $r$  are stored together physically in a file, then:

$$b_r = \left\lceil \frac{n_r}{f_r} \right\rceil$$



# Histograms

- Histogram on attribute *age* of relation *person*



- **Equi-width** histograms
- **Equi-depth** histograms break up range such that each range has (approximately) the same number of tuples
  - E.g. (4, 8, 14, 19)
- Many databases also store  $n$  **most-frequent values** and their counts
  - Histogram is built on remaining values only



# Histograms (cont.)

- Histograms and other statistics usually computed based on a **random sample**
- Statistics may be out of date
  - Some database require a **analyze** command to be executed to update statistics
  - Others automatically recompute statistics
    - e.g., when number of tuples in a relation changes by some percentage



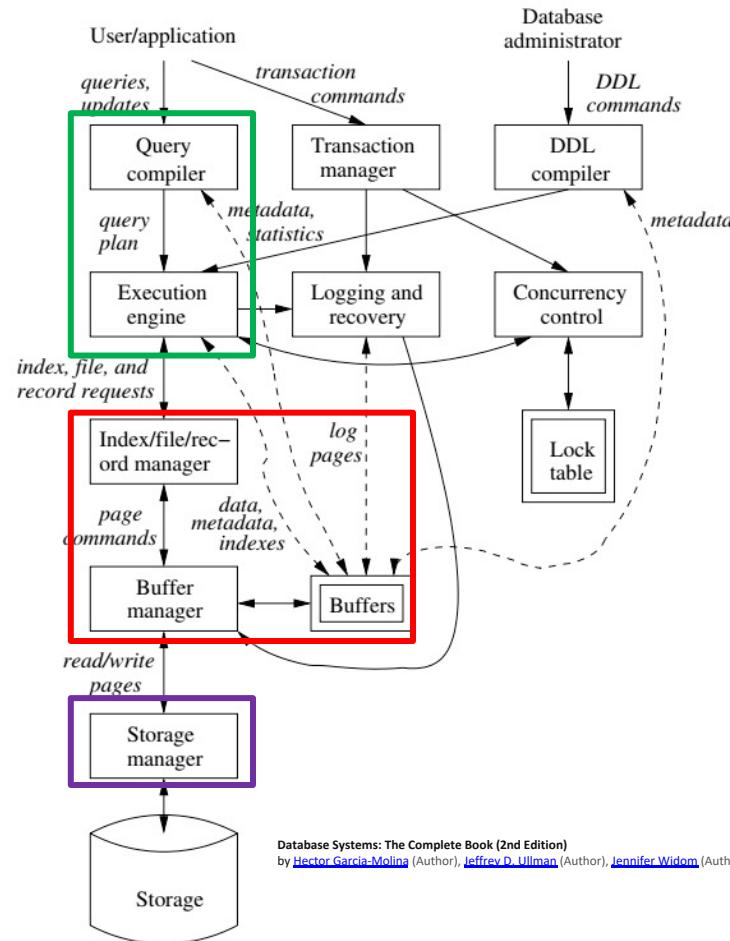
# Selection Size Estimation

- $\sigma_{A=v}(r)$ 
  - $n_r / V(A,r)$  : number of records that will satisfy the selection
  - Equality condition on a key attribute: *size estimate* = 1
- $\sigma_{A \leq v}(r)$  (case of  $\sigma_{A \geq v}(r)$  is symmetric)
  - Let  $c$  denote the estimated number of tuples satisfying the condition.
  - If  $\min(A,r)$  and  $\max(A,r)$  are available in catalog
    - $c = 0$  if  $v < \min(A,r)$
  - $c = n_r \cdot \frac{v - \min(A,r)}{\max(A,r) - \min(A,r)}$
  - If histograms available, can refine above estimate
  - In absence of statistical information  $c$  is assumed to be  $n_r/2$ .

# Data Management

## Today

- ~~Load/save things quickly.~~
- Find things quickly.
- Query/Change Data



# *NoSQL (2)*

# *Reminder*

# Simplistic Classification

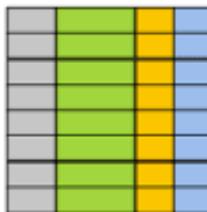
(<https://medium.com/swlh/4-types-of-nosql-databases-d88ad21f7d3b>)

Relational is the foundational model.

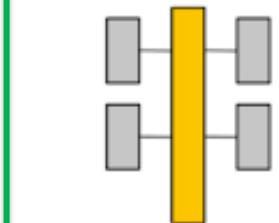
Begin Today

SQL Database

Relational



Analytical (OLAP)

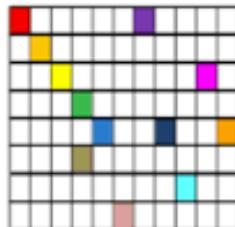


We will see OLAP in a future lecture.

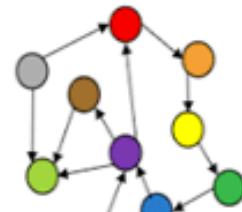
Continue Today

NoSQL Database

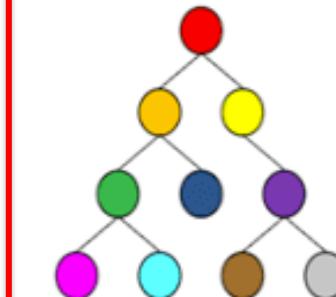
Column-Family



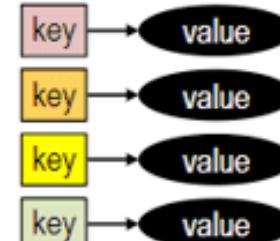
Graph



Document



Key-Value

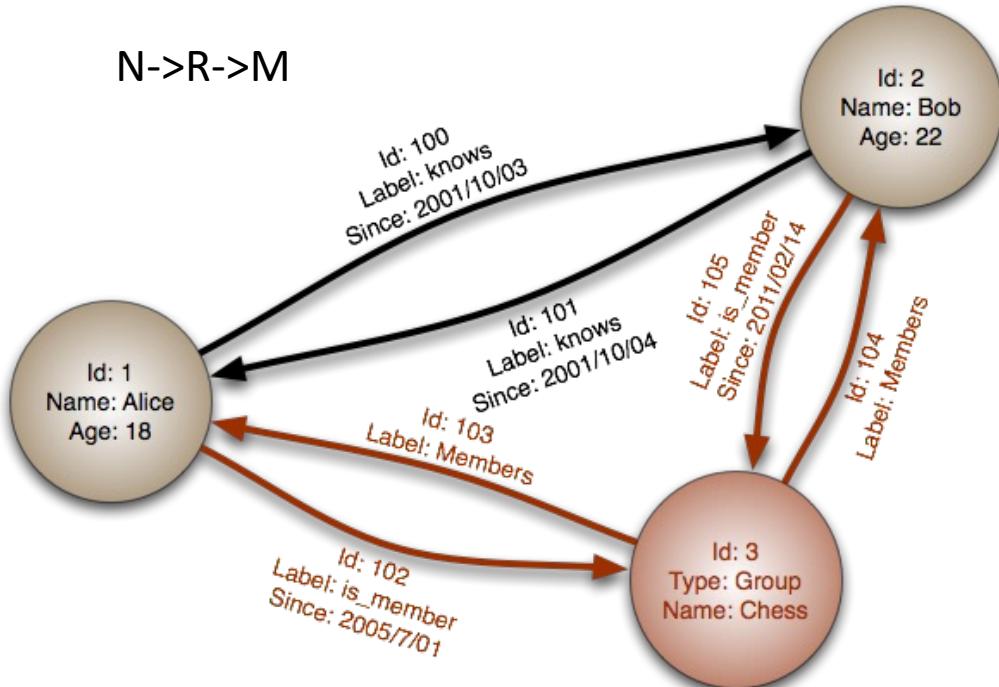


# *Graph Databases*

## *Neo4j*

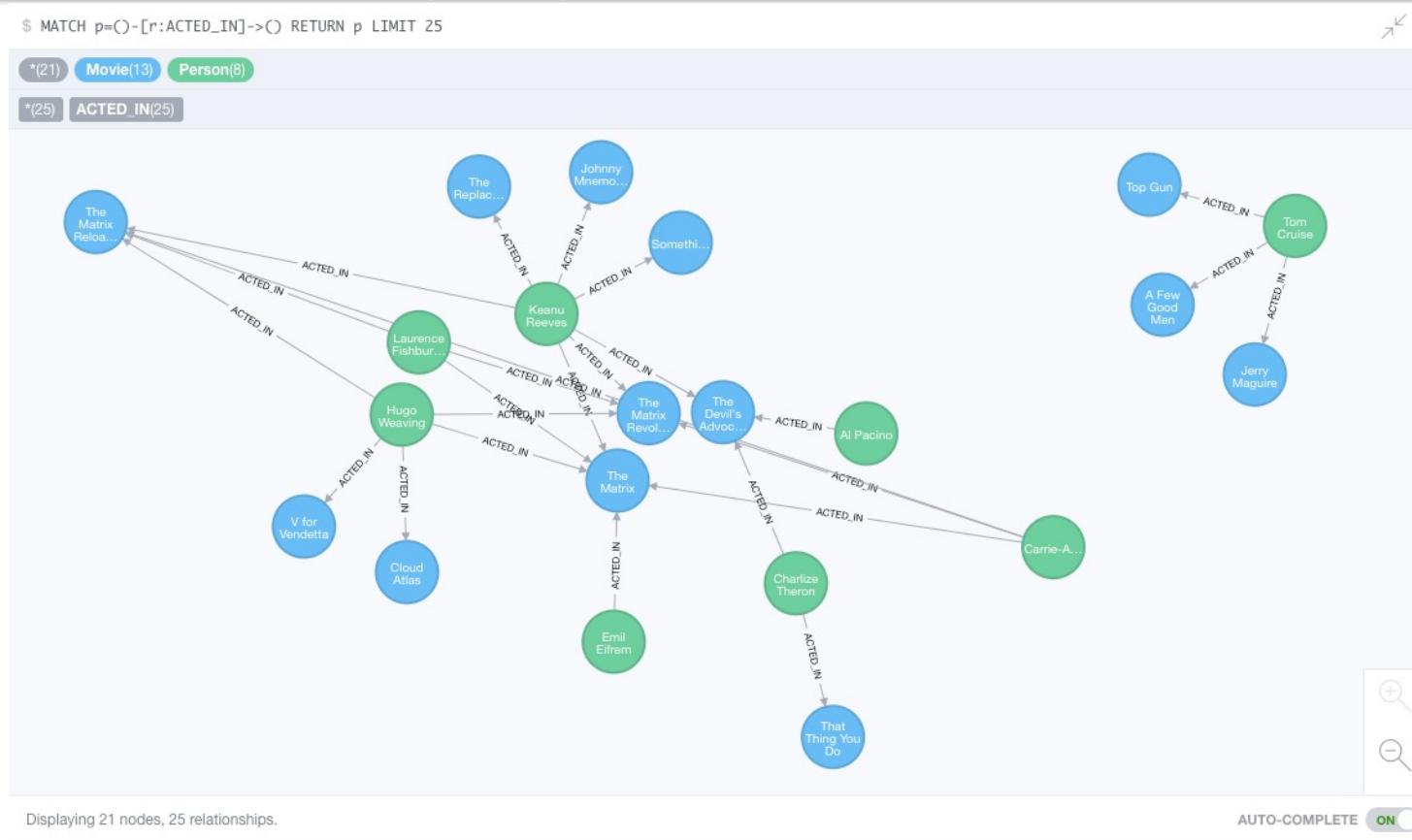
# Graph Database

- Exactly what it sounds like
- Two core types
  - Node
  - Edge (link)
- Nodes and Edges have
  - Label(s) = “Kind”
  - Properties (free form)
- Query is of the form
  - $p1(n)-p2(e)-p3(m)$
  - n, m are nodes; e is an edge
  - $p1, p2, p3$  are predicates on labels



# Neo4J Graph Query

```
$ MATCH p=(:Movie)-[r:ACTED_IN]->(:Person) RETURN p LIMIT 25
```



# Why Graph Databases?

- Schema Less and Efficient storage of Semi Structured Information
- No O/R mismatch – very natural to map a graph to an Object Oriented language like Ruby.
- Express Queries as Traversals. Fast deep traversal instead of slow SQL queries that span many table joins.
- Very natural to express graph related problem with traversals (recommendation engine, find shortest path etc..)
- Seamless integration with various existing programming languages.
- ACID Transaction with rollbacks support.
- Whiteboard friendly – you use the language of node, properties and relationship to describe your domain (instead of e.g. UML) and there is no need to have a complicated O/R mapping tool to implement it in your database. You can say that Neo4j is “Whiteboard friendly” !(<http://video.neo4j.org/JHU6F/live-graph-session-how-allison-knows-james/>)



# Social Network “path exists” Performance

- Experiment:
  - ~1k persons
  - Average 50 friends per person
  - `pathExists(a, b)` limited to depth 4

	# persons	query time
Relational database	1000	2000ms
Neo4j	1000	2ms
Neo4j	1000000	2ms

Graph databases are

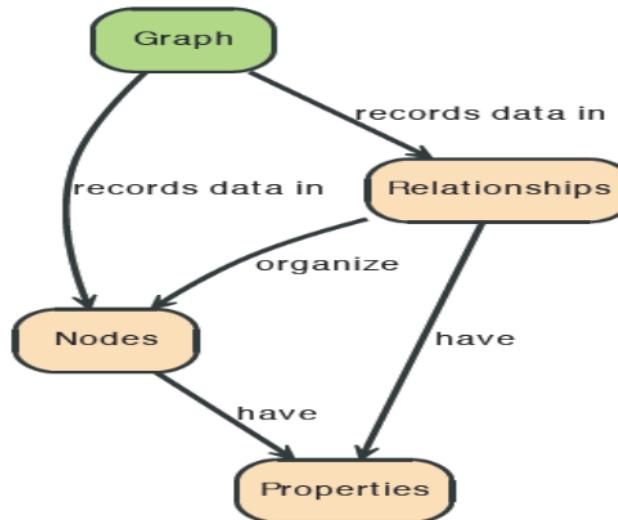
- Extremely fast for some queries and data models.
- Implement a language that vastly simplifies writing queries.

# What are graphs good for?

- Recommendations
- Business intelligence
- Social computing
- Geospatial
- Systems management
- Web of things
- Genealogy
- Time series data
- Product catalogue
- Web analytics
- Scientific computing (especially bioinformatics)
- Indexing your *slow* RDBMS
- And much more!

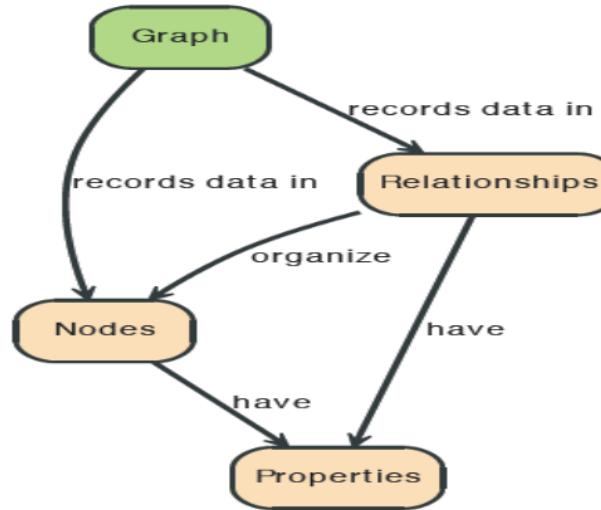
# Graphs

- “A Graph —records data in → Nodes —which have → Properties”



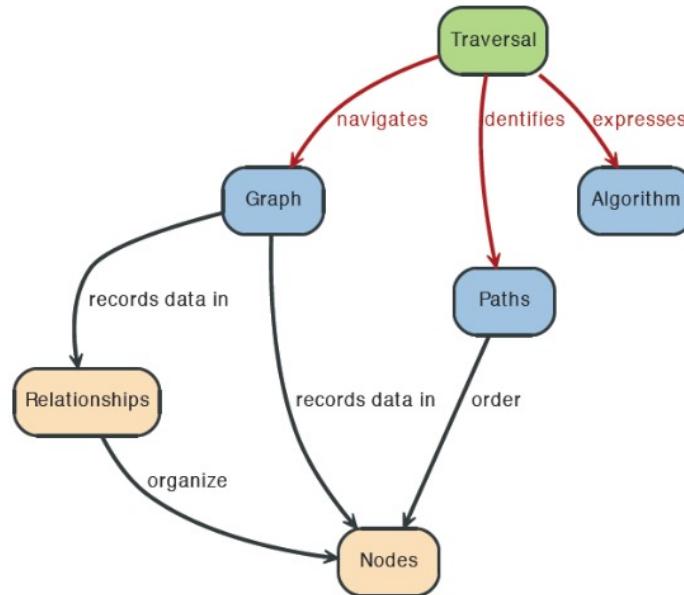
# Graphs

- “Nodes —are organized by→ Relationships — which also have→ Properties”



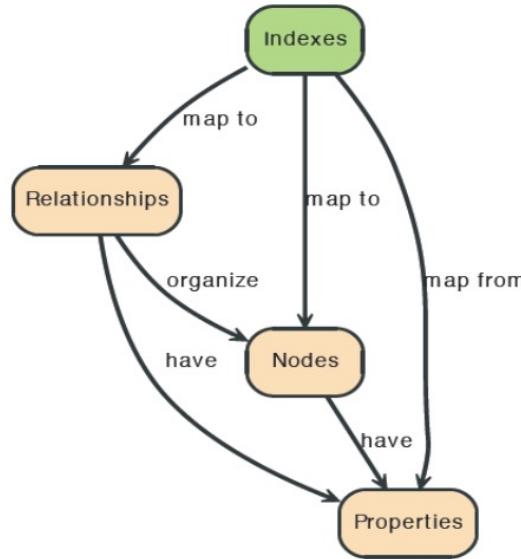
# Query a graph with Traversal

- “A Traversal —navigates→ a Graph; it — identifies→ Paths —which order→ Nodes”

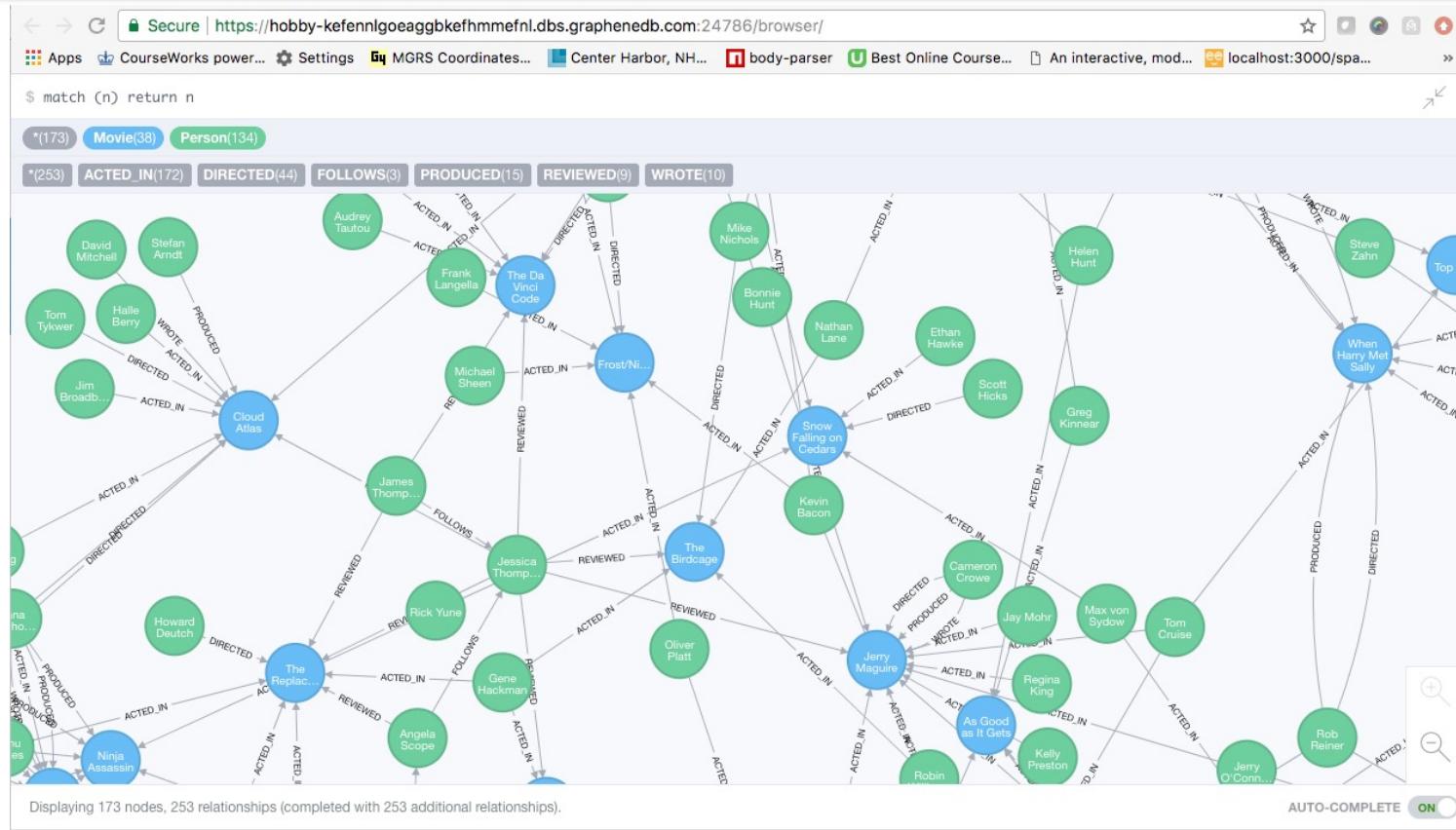


# Indexes

- “An Index —maps from → Properties —to either → Nodes or Relationships”



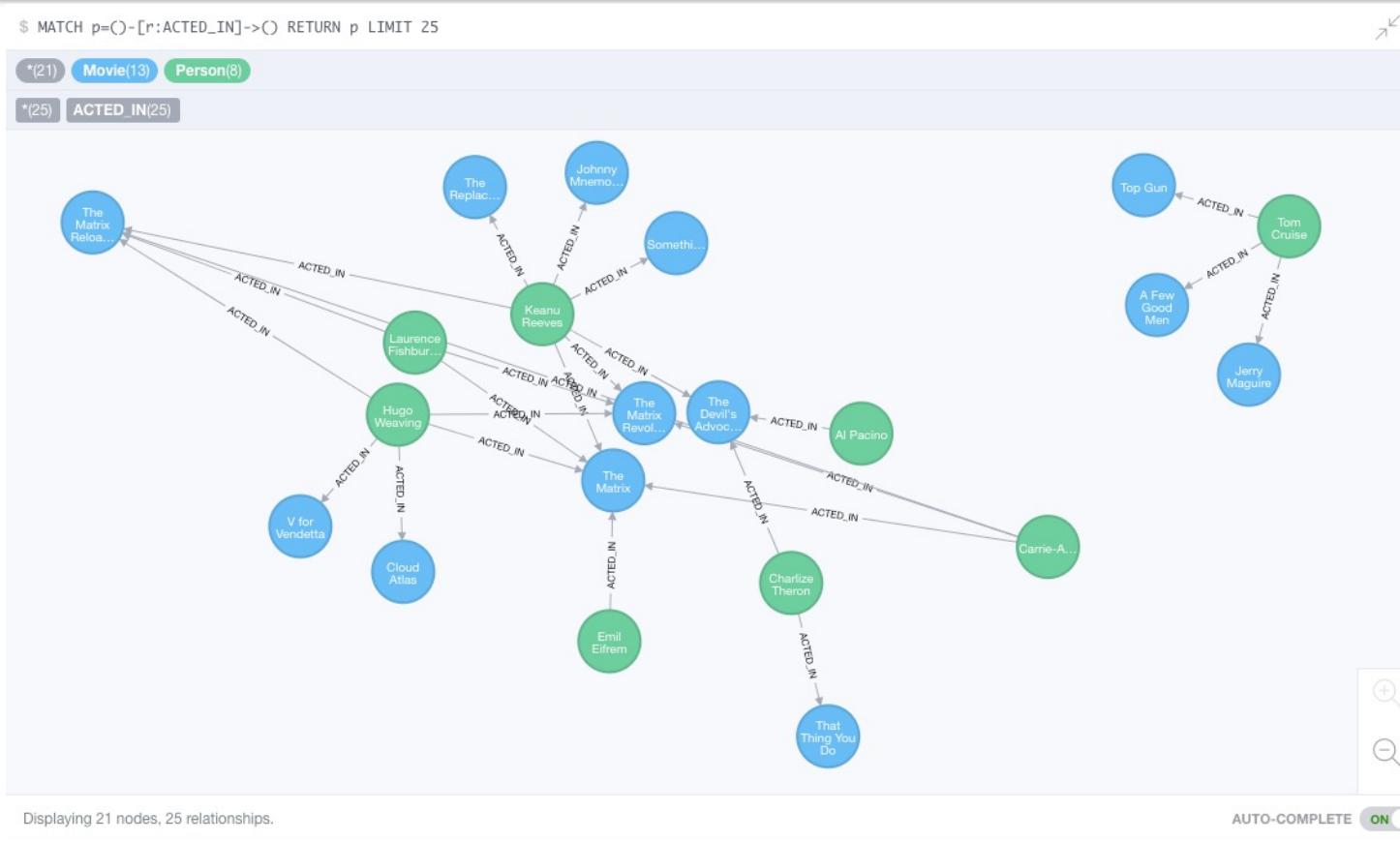
# A Graph Database (Sample)



# Neo4J Graph Query

Who acted in which movies?

```
$ MATCH p=(n)-[r:ACTED_IN]->(m) RETURN p LIMIT 25
```



# Big Deal. That is just a JOIN.

- Yup. But that is simple.
- Try writing the queries below in SQL.

## The Movie Graph Recommend

Let's recommend new co-actors for Tom Hanks. A basic recommendation approach is to find connections past an immediate neighborhood which are themselves well connected.

For Tom Hanks, that means:

1. Find actors that Tom Hanks hasn't yet worked with, but his co-actors have.
2. Find someone who can introduce Tom to his potential co-actor.

Extend Tom Hanks co-actors, to find co-co-actors who haven't work with Tom Hanks...

```
MATCH (tom:Person {name:"Tom Hanks"})-[:ACTED_IN]-(m)<-[ACTED_IN]-(coActors),  
      (coActors)-[:ACTED_IN]-(m2)<-[ACTED_IN]-(cocoActors)  
WHERE NOT (tom)-[:ACTED_IN]-(m2)  
RETURN cocoActors.name AS Recommended, count(*) AS Strength ORDER BY Strength DESC
```

Find someone to introduce Tom Hanks to Tom Cruise

```
MATCH (tom:Person {name:"Tom Hanks"})-[:ACTED_IN]-(m)<-[ACTED_IN]-(coActors),  
      (coActors)-[:ACTED_IN]-(m2)<-[ACTED_IN]-(cruise:Person {name:"Tom Cruise"})  
RETURN tom, m, coActors, m2, cruise
```

# Recommend

```
1 MATCH (tom:Person {name: "Tom Hanks"})-[:ACTED_IN]->(m)<-[:ACTED_IN]-(coActors),  
2     (coActors)-[:ACTED_IN]->(m2)<-[:ACTED_IN]-(cocoActors)  
3 WHERE NOT (tom)-[:ACTED_IN]->(m2)  
4 RETURN cocoActors.name AS Recommended, count(*) AS Strength ORDER BY Strength DESC
```



```
$ MATCH (tom:Person {name: "Tom Hanks"})-[:ACTED_IN]->(m)<-[:ACTED_IN]-(coActors), (coActors)-[:ACTED_IN]->(m2)<-[:ACTED_IN]-(cocoActors) ...
```



	Recommended	Strength
Rows	Tom Cruise	5
A	Zach Grenier	5
Text	Helen Hunt	4
</>	Cuba Gooding Jr.	4
Code	Keanu Reeves	4
	Tom Skerritt	3
	Carrie-Anne Moss	3
	Val Kilmer	3
	Bruno Kirby	3
	Philip Seymour Hoffman	3
	Billy Crystal	3
	Carrie Fisher	3

```

1 MATCH (tom:Person {name:"Tom Hanks"})-[:ACTED_IN]->(m)<-[:ACTED_IN]-(coActors),
2   (coActors)-[:ACTED_IN]->(m2)<-[:ACTED_IN]-(cruise:Person {name:"Tom Cruise"})
3 RETURN tom, m, coActors, m2, cruise

```



\$ MATCH (tom:Person {name:"Tom Hanks"})-[:ACTED\_IN]->(m)<-[:ACTED\_IN]-(coActors), (coActors)-[:ACTED\_IN]->(m2)<-[:ACTED\_IN]-(cruise:Person {name:"Tom Cruise"})



Graph

\*(13) Movie(8) Person(5)

\*(16) ACTED\_IN(16)

Rows

A Text

</> Code



Which actors have  
worked with both  
Tom Hanks and  
Tom Cruise?

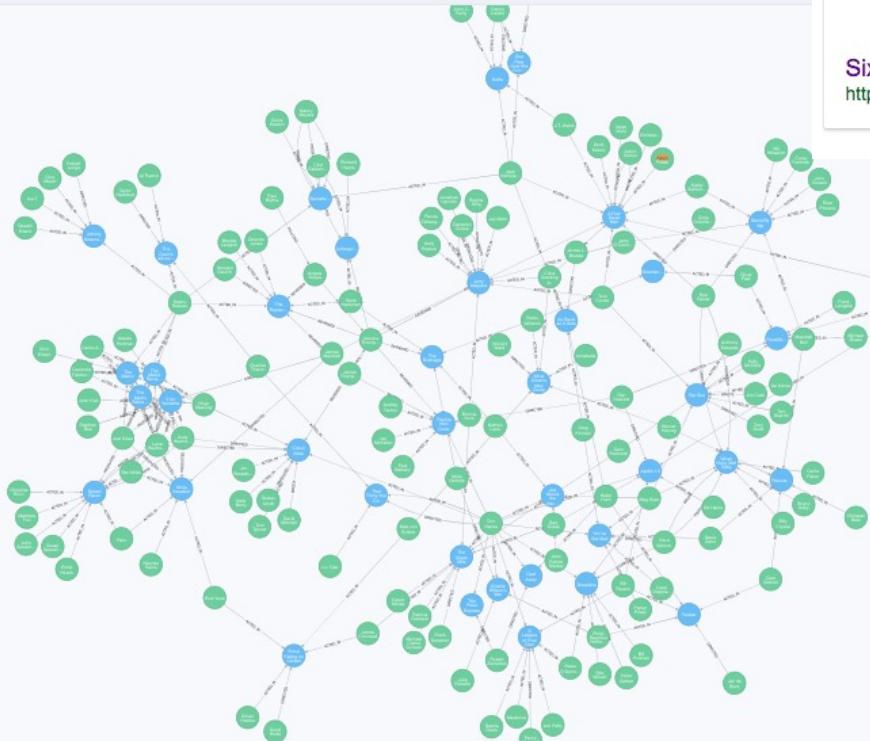
Displaying 13 nodes, 16 relationships (completed with 16 additional relationships).

AUTO-COMPLETE

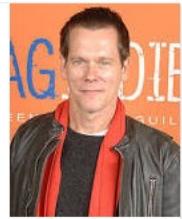
```
$ MATCH (s:Person { name: 'Kevin Bacon' })-[*0..6]-(m) return s,m
```

\*(171) Movie(38) Person(133)

(253) ACTED\_IN(172) DIRECTED(44) FOLLOWS(3) PRODUCED(15) REVIEWED(9) WROTE(10)



**Six Degrees of Kevin Bacon** is a parlour game based on the "six degrees of separation" concept, which posits that any two people on Earth are six or fewer acquaintance links apart. Movie buffs challenge each other to find the shortest path between an arbitrary actor and prolific actor **Kevin Bacon**.



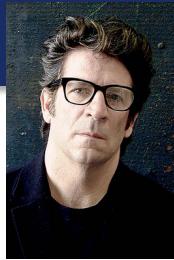
**Six Degrees of Kevin Bacon - Wikipedia**  
[https://en.wikipedia.org/wiki/Six\\_Degrees\\_of\\_Kevin\\_Bacon](https://en.wikipedia.org/wiki/Six_Degrees_of_Kevin_Bacon)

About this result Feedback

## Six Degrees of Kevin Bacon

Game





# How do you get from Kevin Bacon to Robert Longo?

```
$ MATCH (kevin:Person { name: 'Kevin Bacon' }), (robert:Person { name: 'Robert Longo' }), p = shortestPath((kevin)-[*..15]-(robert)) RETURN p
```



Graph  
Rows  
Text  
Code



# *Backup*

# Cloud Concepts – One Perspective

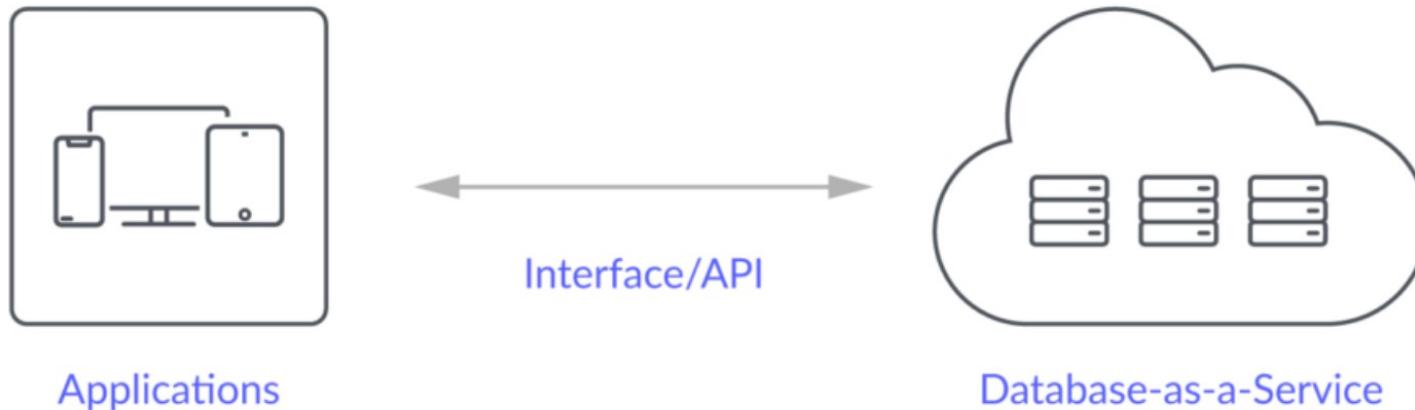
## Categorizing and Comparing the Cloud Landscape

<http://www.theenterprisearchitect.eu/blog/2013/10/12/the-cloud-landscape-described-categorized-and-compared/>

6	SaaS	Applications			<i>End-users</i>
5	App Services	App Services	Communication and Social Services	Data-as-a-Service	<i>Citizen Developers</i>
4	Model-Driven PaaS	Model-Driven aPaaS, bpmPaaS	Model-Driven iPaaS	Data Analytics, baPaaS	<i>Rapid Developers</i>
3	PaaS	aPaaS	iPaaS	dbPaaS	<i>Developers / Coders</i>
2	Foundational PaaS	Application Containers	Routing, Messaging, Orchestration	Object Storage	<i>DevOps</i>
1	Software-Defined Datacenter	Virtual Machines	Software-Defined Networking (SDN), NFV	Software-Defined Storage (SDS), Block Storage	<i>Infrastructure Engineers</i>
0	Hardware	Servers	Switches, Routers	Storage	
		Compute	Communicate	Store	

# Database-as-a-Service

"A cloud database is a database that typically runs on a cloud computing platform and access to the database is provided as-a-service. There are two common deployment models: users can run databases on the cloud independently, using a virtual machine image, or they can purchase access to a database service, maintained by a cloud database provider. Of the databases available on the cloud, some are SQL-based and some use a NoSQL data model. Database services take care of scalability and high availability of the database. Database services make the underlying software-stack transparent to the user." (Wikipedia)





# Cloud Computing Layers

