

# Distributed Data Processing

- Introduction
- Distributed DBMS Architecture
- Distributed DB Design
- Semantic Data Control
- Query Processing
- Transaction Management

# Global Query Optimization

### Input: Algebra query on fragments

- Find the best (not necessarily optimal) global schedule, that is to find the best ordering of operations in the fragment query, including communication operations which minimize a cost function
  - Minimize a cost function
    - Available statistics on fragments
  - Distributed join processing
    - Decide on the use of semijoins
    - Join methods

**Output:** optimized (best) algebraic query with communication operations included on fragments



# Local Query Optimization

### Input: Best global execution schedule

- Select the **best access path** by all the site
- Use the **centralized** optimization techniques

# 7. Optimization of Distributed Queries

- Query Optimization
- Centralized Query Optimization
- Join Ordering in Fragment Queries
- Distributed Query Optimization Algorithms
- Local Optimization
- Conclusion

# Objective of optimizer

- Finding an "optimal" ordering of operations for a given query
  - Selecting the optimal execution strategy for a query is NP-hard in the number of relations
    - The actual objective is to find a strategy close to optimal
  - Input to optimizer
    - query on fragments
    - fragment statistics, formulas for estimating the cardinalities of results of relational operations
  - Focus mostly on the ordering of join operation
    - It is a well-understood problem, and queries involving joins, selections, and projection are usually considered to be the most frequent type
    - It is easier to generalize the basic algorithm for other binary operations, such as unions.

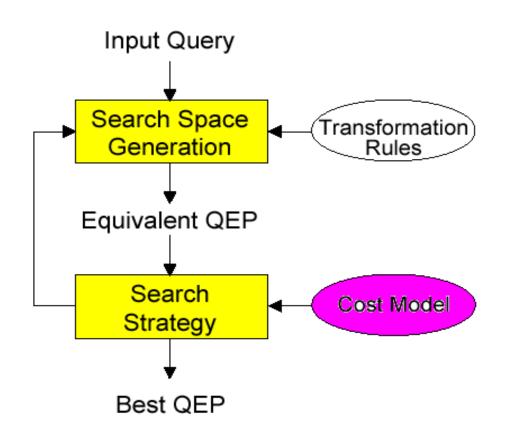


# 7.1 Query Optimization

# **Query Optimization**

- Solution space
  - The set of equivalent algebra expressions (operator trees).
- Cost function (in terms of time)
  - Total time (Response time): I/O cost + CPU cost + communication cost
  - These might have different weights in different distributed environments (LAN vs WAN).
- Search algorithm
  - How do we move inside the solution space?
  - Exhaustive search, heuristic algorithms

### **Query Optimization Process**



# 7.1.1 Search Space

- Search space characterized by alternative execution plans
  - could be expressed as operator trees
- Focus on join trees
  - Operator tree whose operators are join or Cartesian product
  - Because permutations of the join order have the most important effect on performance of queries

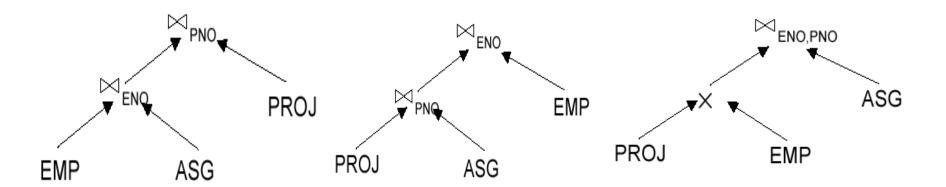


SELECT ENAME, RESP

FROM EMP, ASG, PROJ

WHERE EMP.ENO=ASG.ENO

AND ASG.PNO=PROJ.PNO



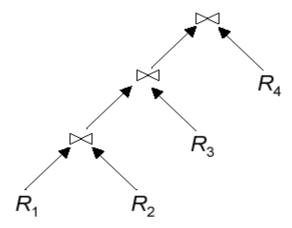


### Restrict the size of the search space

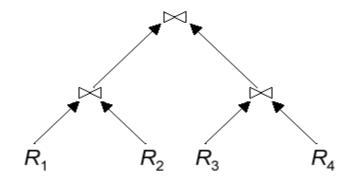
- Restrict by means of heuristics
  - Perform unary operations before binary operations
  - Avoid Cartesian product
- Restrict the shape of the join tree
  - Consider only linear trees, ignore bushy trees

## Linear tree vs. Bushy tree

#### Linear Join Tree



#### **Bushy Join Tree**



## 7.1.2 Search Strategy

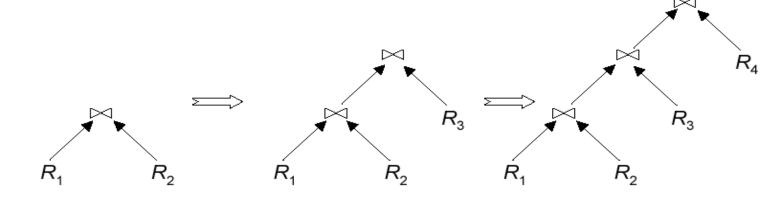
How to "move" in the search space.

- Deterministic
  - Start from base relations and build plans by adding one relation at each step until complete plans are obtained
  - Dynamic programming: breadth-first
  - Greedy: depth-first
- Randomized
  - Search for optimalities around a particular starting point
  - Trade optimization time for execution time
  - Better when > 5-6 relations

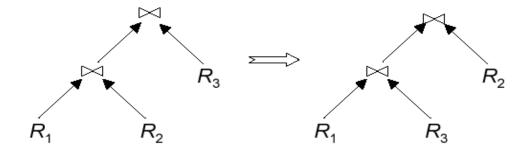
# 1

### Deterministic vs. Randomized

#### Deterministic



Randomized





### 7.1.3 Distributed Cost Model

- Cost function
  - To predict the cost of operators
- Database statistics
  - About the base relations and formulas to evaluate the sizes of intermediate results

### **Cost Function**

- Total Time (or Total Cost)
  - Do as little of each cost component as possible
    - Reduce each cost (in terms of time) component individually
  - Optimizes the utilization of the resources
    - => Increase system throughput
- Response Time
  - Do as many things as possible in parallel
    - May increase total time because of increased total activity
  - Optimizes the degree of parallel execution
    - => Improve user's response time

### Total cost

### Summation of all cost factors

Total cost = CPU cost + I/O cost + communication cost

CPU cost =

unit instruction cost \*number of instructions

I/O cost =

unit disk I/O cost \* number of disk I/Os

communication cost =

message initiation + transmission

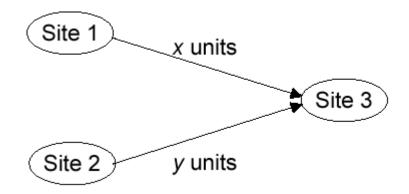
### Response time

# Elapsed time between the initiation and the completion of a query

Response time = CPU time + I/O time + communication time

- CPU time =
  - unit instruction time \* number of *sequential* instructions
- I/O time = unit I/O time \* number of sequential I/Os
- communication time =
  - unit msg initiation time \* number of *sequential* msg + unit transmission time \* number of *sequential* bytes

# Example



#### Assume that only the communication cost is considered

**Total time** = 2 \* message initialization time + unit transmission time \* (x+y)

**Response time** = max {time to send x from 1 to 3, time to send y from 2 to 3}

time to send x from 1 to 3 = message initialization time + unit transmission time \* x

time to send y from 2 to 3 = message initialization time + unit transmission time \* y

# Database statistics

Primary cost factor:

#### size of intermediate relations

- The estimation is based on statistical information about the base relations and formulas to predict the cardinalities of the results of the relational operations
- Make them precise more costly to maintain

### Database statistics – Base relation

- For each relation R[A1, A2, ..., An] fragmented as R1, ..., Rr
  - length of each attribute: length(Ai)
  - the number of distinct values for each attribute in each fragment:  $card(\prod_{Ai} Rj)$
  - maximum and minimum values in the domain of each attribute:
     min(Ai), max(Ai)
  - the cardinalities of each domain: card(dom[Ai])
  - the cardinalities of each fragment: card(Rj )
- Selectivity factor of each operation for relations
  - the proportion of tuples of an operand relation that participate in the result of that operation
  - For joins

$$card(R \infty S)$$
  
 $SF_{\infty}(R,S) = \cdots$   
 $card(R) * card(S)$ 



### Assumptions

- The distribution of attribute values in a relation is supposed to be uniform
- All attributes are independent

### Cardinalities of intermediate results

#### Selection

$$size(R) = card(R) * length(R)$$

$$card(\sigma_F(R)) = SF_{\sigma}(F) * card(R)$$
where
$$SF_{\sigma}(A = value) = \frac{1}{card(\prod_A(R))}$$

$$SF_{\sigma}(A > value) = \frac{max(A) - value}{max(A) - min(A)}$$

$$SF_{\sigma}(A < value) = \frac{value - max(A)}{max(A) - min(A)}$$

$$SF_{\sigma}(p(A_i) \land p(A_j)) = SF_{\sigma}(p(A_i)) * SF_{\sigma}(p(A_j))$$

$$SF_{\sigma}(p(A_i) \lor p(A_j)) = SF_{\sigma}(p(A_i)) + SF_{\sigma}(p(A_j)) - (SF_{\sigma}(p(A_i)) * SF_{\sigma}(p(A_j)))$$

$$SF_{\sigma}(A \in value) = SF_{\sigma}(A = value) * card(\{values\})$$

### Cardinalities of intermediate results

#### Projection

 $card(\Pi_A(R)) = card(R)$  If one of the projected attributes is a key of R

#### Cartesian Product

$$card(R \times S) = card(R) * card(S)$$

#### Union

upper bound:  $card(R \cup S) = card(R) + card(S)$ 

lower bound:  $card(R \cup S) = max\{card(R), card(S)\}$ 

#### Set Difference

upper bound: card(R-S) = card(R)

lower bound: 0

### Cardinalities of intermediate results

#### Join

Special case: A is a key of R and B is a foreign key of S;

$$card(R\bowtie_{A=B} S) = card(S)$$

■ More general:

$$card(R\bowtie S) = SF_{\bowtie}* card(R) * card(S)$$

#### Semijoin

$$card(R \bowtie_{A} S) = SF_{\bowtie}(S.A) * card(R)$$

where

$$SF_{\bowtie}(R\bowtie_{A} S) = SF_{\bowtie}(S.A) = \frac{card(\prod_{A}(S))}{card(dom[A])}$$



# 7.2 Centralized Query Optimization

# Centralized Query Optimization

- Centralized query optimization is a simpler problem
- Distributed query optimization techniques are often extensions of the techniques for centralized system
- A distributed query is translated into local queries, each of which is processed in a centralized way

# Two popular relational database

- INGRES
  - dynamic
- System R
  - static
  - exhaustive search

# 7.2.1 INGRES Algorithm

Dynamic query optimization

Combine the two phase of decomposition and optimization

**Step1:** Decompose each multi-variable query into a sequence of mono-variable queries with a common variable

**Step2:** Process each by a one variable query processor

- Choose an initial execution plan (heuristics)
- Order the rest by considering intermediate relation sizes
- No statistical information is maintained

### Decomposition

Replace an *n* variable query *q* by a series of queries

$$q1 \rightarrow q2 \rightarrow ... \rightarrow qn$$

where  $q_i$  uses the result of  $q_{i-1}$ .

- Detachment
  - Query q decomposed into  $q' \rightarrow q''$  where q' and q'' have a common variable which is the result of q'
- Tuple substitution
  - Replace the value of each tuple with actual values and simplify the query

$$q(V1, V2, ... Vn) \rightarrow (q'(t1, V2, V3, ..., Vn), t1 \in R)$$



### Detachment

```
q: SELECT R2.A2, R3.A3, ..., Rn.An FROM R1, R2, ..., Rn WHERE P1(R1.A1')
AND P2(R1.A1, R2.A2, ..., Rn.An)
```

q': SELECT R1.A1 INTO R1'

FROM R1

**WHERE** P1(R1.A1')

q": SELECT R2.A2, R3.A3, ..., Rn.An

FROM R1', R2, ..., Rn

WHERE P2(R1'.A1, R2.A2, ..., Rn.An)

### Example of detachment

Names of employees working on CAD/CAM project

 $q_1$ : **SELECT** EMP.ENAME

FROM EMP, ASG, PROJ

WHERE EMP.ENO=ASG.ENO

**AND** ASG.PNO=PROJ.PNO

**AND** PROJ. PNAME="CAD/CAM"

 $q_{11}$ : **SELECT** PROJ.PNO **INTO** JVAR

FROM PROJ

WHERE PROJ. PNAME="CAD/CAM"

q': SELECT EMP.ENAME

FROM EMP, ASG, JVAR

WHERE EMP.ENO=ASG.ENO

**AND** ASG.PNO=JVAR.PNO

### Example of detachment

q': SELECT EMP.ENAME

FROM EMP, ASG, JVAR

WHERE EMP.ENO=ASG.ENO

**AND** ASG.PNO=JVAR.PNO

 $q_{12}$ : **SELECT** ASG.ENO **INTO** GVAR

FROM ASG, JVAR

WHERE ASG.PNO=JVAR.PNO

 $q_{13}$ : SELECT EMP.ENAME

FROM EMP, GVAR

WHERE EMP.ENO=GVAR.ENO

### Example of tuple substitution

 $q_{11}$  is a mono-variable query  $q_{12}$  and  $q_{13}$  is subject to tuple substitution Assume GVAR has two tuples only: <E1> and <E2> Then  $q_{13}$  becomes

 $q_{131}$ : SELECT EMP.ENAME

FROM EMP

WHERE EMP. ENO="E1"

 $q_{132}$ : **SELECT** EMP.ENAME

FROM EMP

WHERE EMP. ENO= "E2"

# **INGRES Algorithm**

- Applying the selections and projections as soon as possible by detachment
- Results of the monorelation queries are **stored** in data structures that are capable of optimizing the later queries (such as joins)
- The irreducible queries that remain after detachment must be processed by tuple substitution
- For the irreducible query, the smallest relation whose cardinality is known from the result of the preceding query is chosen for substitution
- Monorelation queries generated by the reduction algorithm are processed by the OVQP that chooses the best existing access path to the relation, according to the query qualification

# 7.2.2 System R Algorithm

Static query optimization based on the exhaustive search of the solution space

- INPUT: relational algebra tree resulting from the query decomposition
- OUTPUT: an execution plan that implements the "optimal" relational algebra tree

Statistical information is maintained

## System R Algorithm

- To limit the overhead of optimization, the number of alternative trees is reduced using dynamic programming
  - The set of alternative strategies is constructed dynamically so that, when two joins are equivalent by commutatively, only the cheapest one is kept
  - The strategies that include Cartesian products are eliminated whenever possible

# System R Algorithm

- Step1: the best access path to each individual relation based on a select predicate is predicted
- Step2: the best join ordering is estimated for each relation R
  - Determine the possible ordering of joins
  - Determine the cost of each ordering
  - Choose the join ordering with minimal cost

#### **Alternatives for Join**

#### Nested loops

```
for each tuple of external relation (cardinality n1)
for each tuple of internal relation (cardinality n2)
join two tuples if the join predicate is true
end
end
```

- Complexity: n1\* n2
- Merge join

```
sort relations
merge relations
```

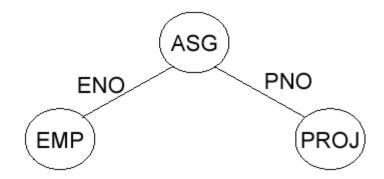
Complexity: n1 + n2 if relations are previously sorted and equijoin



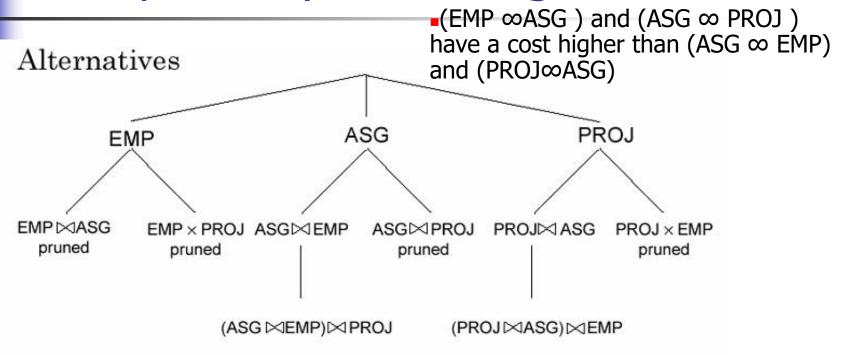
Query: Names of employees working on the CAD/CAM project

#### Assume

- EMP has an index on ENO,
- ASG has an index on PNO,
- PROJ has an index on PNO and an index on PNAME
- (EMP ∞ASG) and (ASG ∞ PROJ) have a cost higher than (ASG ∞ EMP) and (PROJ∞ASG)



- Choose the best access paths to each relation
  - EMP: sequential scan (no selection on EMP based on ENO)
  - ASG: sequential scan (no selection on ASG based on PNO)
  - PROJ: index on PNAME (there is a selection on PROJ based on PNAME)
- Determine the best join ordering
  - EMP ∞ASG ∞PROJ
  - ASG ∞ PROJ ∞ EMP
  - PROJ∞ASG ∞ FMP
  - ASG ∞ EMP ∞ PROJ
  - EMP x PROJ ∞ ASG
  - PROJ x EMP ∞ ASG
  - Select the best ordering based on the join costs evaluated according to the two methods



Best total join order is one of ((ASG⋈EMP)⋈PROJ) ((PROJ⋈ASG)⋈EMP)

- ((PROJ ∞ ASG) ∞ EMP) has a useful index on the select attribute and direct access to the join attributes of ASG and EMP
- Therefore, chose it with the following access methods:
  - select PROJ using index on PNAME
  - then join with ASG using index on PNO
  - then join with EMP using index on ENO

### System R Algorithm

- To select the best single-relation access method to each relation in the query
- To examine all possible permutations of join orders and select the best access strategy for the query
  - First, the join of each relation with every other relation is considered
  - Then, joins of three relations are optimized
  - The continues until joins of n relations are optimized



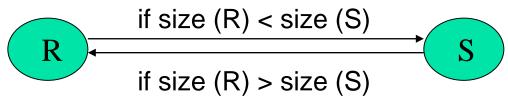
# 7.3 Join Ordering in Fragment Queries

### Ordering Joins in Fragment Queries

- Ordering joins
  - Distributed INGRES
  - System R\*
- Semijoin ordering
  - SDD-1

### 7.3.1 Join Ordering

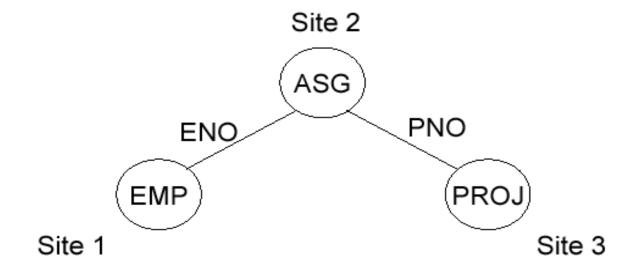
- Consider two relations only
  - Send the smaller relation to the site of the large one

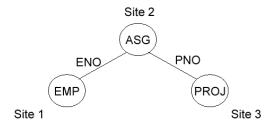


- Multiple relations more difficult because too many alternatives.
  - Compute the cost of all alternatives and select the best one.
    - Join may reduce or increase the size of the intermediate results
    - Necessary to compute the size of intermediate relations which is difficult
  - Use heuristics

### **Example of Join Ordering**

 $\begin{array}{c} Consider \\ PROJ \bowtie_{PNO} ASG \bowtie_{ENO} EMP \end{array}$ 





#### Example of Join Ordering

#### Execution alternatives:

1. EMP  $\rightarrow$  Site 2

Site 2 computes EMP'=EMP™ASG

 $EMP' \rightarrow Site 3$ 

Site 3 computes EMP™ PROJ

2. ASG  $\rightarrow$  Site 1

Site 1 computes EMP'=EMP™ASG

EMP'  $\rightarrow$  Site 3

Site 3 computes EMP'™ PROJ

 $3. ASG \rightarrow Site 3$ 

 $ASG' \rightarrow Site 1$ 

Site 1 computes ASG'™ EMP

4. PROJ  $\rightarrow$  Site 2

Site 3 computes ASG'=ASG ⋈ PROJ Site 2 computes PROJ'=PROJ ⋈ ASG

 $PROJ' \rightarrow Site 1$ 

Site 1 computes PROJ'™ EMP

5. EMP  $\rightarrow$  Site 2

 $PROJ \rightarrow Site 2$ 

Site 2 computes EMP $\bowtie$  PROJ  $\bowtie$  ASG

### 7.3.2 Semijoin Based Algorithms

- Consider the join of two relations:
  - R[A] (located at site 1)
  - S[A](located at site 2)
- Alternatives:
  - 1. Do the join  $R \sim_A S$
  - 2. Perform one of the semijoin equivalents

$$R \sim_{A} S \Leftrightarrow (R \sim_{A} S) \sim_{A} S$$
  
 $\Leftrightarrow R \sim_{A} (S \sim_{A} R)$   
 $\Leftrightarrow (R \sim_{A} S) \sim_{A} (S \sim_{A} R)$ 

#### Join vs. Semijoin Algorithms

- Perform the join
  - $\implies$  send R to Site 2
  - $\blacksquare$  Site 2 computes  $R\bowtie_A S$
- Consider semijoin  $(R \bowtie_A S) \bowtie_A S$ 
  - $\Rightarrow S' \leftarrow \prod_A(S)$
  - $\longrightarrow$  Site 1
  - $\blacksquare$  Site 1 computes  $R' = R \bowtie_A S'$
  - $R' \rightarrow \text{Site } 2$
  - $\Longrightarrow$  Site 2 computes  $R' \bowtie_A S$

Semijoin is better if

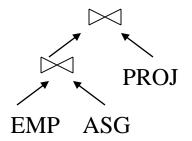
$$size(\Pi_A(S)) + size(R \bowtie_A S)) < size(R)$$

#### Join vs. Semijoin

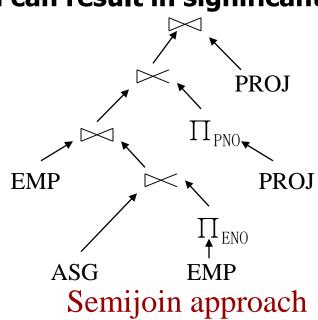
- Semijoin may increase the local processing time
- However, if
  - The join attribute length is smaller than the length of an entire tuple
  - The semijoin has good selectivity

Then the semijoin approach can result in significant savings in

communication time



Join approach





# 7.4 Distributed Query Optimization Algorithms

# Three Basic Distributed Query Optimization algorithms

- Distributed INGRES
  - Ordering joins
- <u>R\*</u>
  - Ordering joins
- <u>SDD-1</u>
  - Semijoin ordering

# Distributed Query Optimization Algorithms

Algorithms	Opt. Timing	Objective Function	Opt. Factors	Network Topology	Semijoin	Stats	Fragments
Dist. INGRES	Dynamic	Resp. time or Total time	Msg. Size, Proc. Cost	General or Broadcast	No	1	Horizontal
R*	Static	Total time	No. Msg., Msg. Size, IO, CPU	General or Local	No	1, 2	No
SDD-1	Static	Total time	Msg. Size	General	Yes	1,3,4, 5	No

<sup>1:</sup> relation cardinality; 2: number of unique values per attribute; 3: join selectivity factor; 4: size of projection on each join attribute; 5: attribute size and tuple size

## 7.4.1 Distributed INGRES Algorithm

- Same as the centralized version except
  - Movement of relations (and fragments) need to be considered
  - Optimization with respect to communication cost or response time possible



#### Distributed INGRES Algorithm

- Dynamic
- Considers only joins
- The algorithm also takes advantage of fragmentation, but only horizontal fragmentation is handled for simplicity
- Both general and broadcast networks are considered

#### Distributed INGRES Algorithm

```
Input: MRQ: multirelation query
Output: result of the last multirelation query
Begin
    Run all detachable one-relation queries in MRQ;
    Replace MRQ by a list of n irreducible queries MRQ'_list;
    Repeat
       choose next irreducible query MRQ' from MRQ'_list involving the
       smallest fragments;
       Determine fragments to transfer and processing site for MRQ';
       Move the selected fragments to the selected sites;
       Run MRO';
    Until MRQ' list empty
End
```

# Optimization in Distributed INGRES Algorithm

- choose next irreducible query
  - Having no predecessor and involving the smaller fragments
- Determine fragments to transfer and processing site
  - based on communication cost
  - Assuming that one relation Rp is the remaining fragmented, and K sites participate in processing, then the fragments of Ri have to be moved as follows:
    - For a processing site j, Rij is moved to k-1 other sites
    - For a non-processing site j, Rij is moved to k other sites and any fragments of Rpj are moved to any one processing site
  - broadcast
    - CTk(#bytes) = CT1(#bytes)
  - point—to—point
    - CTk(#bytes) = k\*CT1(#bytes)

# -

#### Example of Distributed INGRES Algorithm

PROJ ∞ ASG

	Site 1	Site 2	Site 3	Site 4
PROJ	1000	1000	1000	1000
ASG			2000	

- Point-to-point
  - Send each PROJi to site 3 (cost:3000)
- Broadcast
  - Send ASG (in a single transfer) to 1,2,4 (cost:2000)

### 7.4.2 R\* Algorithm

- Exhaustive search
- Compilation
- Considers only joins
- Cost function includes local processing as well as transmission
- Published papers provide solutions to handling horizontal and vertical fragmentations but the implemented prototype does not

### R\* Algorithm

```
Input: QT: Query tree
Output: minimum cost strategy
Begin
   For each relation R_i \in QT do
       Get the best_AP; from all access path to R;
    For each order do
       Get the best join order from all the orders
   For each site k storing a relation involved in QT do
       LS_k \leftarrow local strategy
       Send(LS<sub>k</sub>, site k)
end
```

#### R\* Algorithm

- Select the join ordering
- Join algorithm
  - Nested loop
  - Merge join
- Access path for each fragment
  - Index
  - Sequential scan
- Select the sites of join results
- Method of transferring data between sites

# Performing joins – transfer method

#### Ship whole

- larger data transfer
- smaller number of messages
- better if relations are small

#### Fetch as needed

- number of messages = O(cardinality of external relation)
- data transfer per message is minimal
- better if relations are large and the selectivity is good

- Move entire external relation to the site of the internal relation
  - Retrieve external relation R tuples
  - Send them to the internal relation S site
  - Join them as they arrive

```
Total Cost = LT(retrieve card(R) tuples from R)
```

- + CT(size(R))
- + LT(retrieve card( $S \propto_A R$ ) tuples from S)

- Ship the entire internal relation to the site of the external relation
  - Cannot join as they arrive; they need to be stored

```
Total Cost = LT(retrieve card(S) tuples from S)
+ CT(size(S))
+LT(store card(S) tuple in T)
+LT(retrieve card(R) tuples from R)
+ LT(retrieve card(S\propto_AR) tuples from T)
```

- Fetch internal tuples as needed
  - Retrieve qualified tuples at external relation site
  - Send request containing join column values of external relation to internal relation site
  - Retrieve matching internal tuples at internal relation site
  - Send the matching internal tuples to external relation site
  - Join as they arrive

```
Total Cost = LT(retrieve card(R) tuples from R)
+ CT(length(A)*card(R))
+ LT(retrieve card(S \propto_A R) tuples from T)
+ CT(card(S \propto_A R) *length(S))
```

Move both relations to another site

```
Total Cost = LT(retrieve card(S) tuples from S)
+ CT(size(S))
+LT(store card(S) tuple in T)
+LT(retrieve card(R) tuples from R)
+ CT(size(R))
+ LT(retrieve card(S\propto_AR) tuples from T)
```

# R\* Algorithm

- Predict the total time of each strategy
- Selects the cheapest.

# 4

#### Example of R\* Algorithm

PROJ  $\infty_{PNO}$  ASG

(Assume: two site; ASG: Index on PNO)

- 1. Ship whole PROJ to site of ASG
- 2. Ship whole ASG to site of PROJ
- 3. Fetch ASG tuples as needed for each tuple of PROJ
- 4. Move ASG and PROJ to a third site

## 7.4.3 SDD-1 Algorithm

- Based on the Hill Climbing Algorithm
  - Static
  - Semijoins
  - No replication
  - No fragmentation
  - Cost of transferring the result to the user site from the final result site is not considered

## Idea – Greedy Algorithm

Refinements of an initial feasible solution are recursively computed until no more cost improvement can be made

## Hill Climbing Algorithm

The first distributed query processing algorithm

#### Assume join is between three relations.

Step 1: Do initial local processing

Step 2: Select initial feasible solution (ESO)

- Determine the candidate result sites sites where a relation referenced in the query exist
- Compute the cost of transferring all the other referenced relations to each candidate site
- ESO = candidate site with minimum cost

# Hill Climbing Algorithm

Step 3: Determine candidate splits of ESO into {ES1, ES2 }

- ES1 consists of sending one of the relations to the other relation's site
- ES2 consists of sending the join of the relations to the final result site

Step 4: Replace ESO with the split schedule which gives cost(ES1) + cost(local join) + cost(ES2) < cost(ES0)

Step 5: Recursively apply steps 3-4 on ES1 and ES2 until no such plans can be found

Step 6: Check for redundant transmissions in the final plan and eliminate them.

What are the salaries of engineers who work on the CAD/CAM project?

$$\Pi_{SAL}(PAY \bowtie_{TITLE}(EMP \bowtie_{ENO}(ASG \bowtie_{PNO}(\sigma_{PNAME="CAD/CAM"}(PROJ)))))$$

<u>Relation</u>	$\underline{\text{Size}}$	Site
$\mathbf{EMP}$	8	1
PAY	4	2
PROJ	1	3
ASG	10	4

#### Assume:

- Size of relations is defined as their cardinality
- Minimize total cost
- ➡ Transmission cost between two sites is 1
- Ignore local processing cost

```
Based on join selectivities,

size(EMP \infty PAY) = size(EMP),

size(PROJ \infty ASG) = 2*size(PROJ),

size(ASG \infty EMP) = size(ASG)
```



#### Step 1:

Selection on PROJ; result has cardinality 1

<u>Relation</u>	$\underline{\text{Size}}$	Site
EMP	8	1
PAY	4	2
PROJ	1	3
ASG	10	4

Step 2: Initial feasible s
----------------------------

Alternative 1: Resulting site is Site 1

Total cost = 
$$cost(PAY \rightarrow Site 1) + cost(ASG \rightarrow Site 1) + cost(PROJ \rightarrow Site 1)$$
  
=  $4 + 10 + 1 = 15$ 

Relation

**EMP** 

PAY

**PROJ** 

ASG

Size

8

4

10

Site

1

4

Alternative 2: Resulting site is Site 2

Total cost = 
$$8 + 10 + 1 = 19$$

Alternative 3: Resulting site is Site 3

Total cost = 
$$8 + 4 + 10 = 22$$

Alternative 4: Resulting site is Site 4

Total cost = 
$$8 + 4 + 1 = 13$$

Therefore  $ES_0 = \{ \text{EMP} \rightarrow \text{Site 4}; \text{PAY} \rightarrow \text{Site 4}; \text{PROJ} \rightarrow \text{Site 4} \}$ 

Relation	Size	Site
EMP	8	1
PAY	4	2
PROJ	1	3
ASG	10	4

Step 3: Determine candidate splits

Alternative 1:  $\{ES_1, ES_2, ES_3\}$  where

 $ES_1$ : EMP  $\rightarrow$  Site 2

 $ES_2$ : (EMP $\bowtie$  PAY)  $\rightarrow$  Site 4

 $ES_3$ : PROJ  $\rightarrow$  Site 4

Alternative 2:  $\{ES_1, ES_2, ES_3\}$  where

 $ES_1: PAY \rightarrow Site 1$ 

 $ES_2$ : (PAY  $\bowtie$  EMP)  $\rightarrow$  Site 4

 $ES_3$ : PROJ  $\rightarrow$  Site 4

#### Step 4: Determine costs of each split alternative

$$cost(\text{Alternative 1}) = cost(\text{EMP} \rightarrow \text{Site 2}) + cost((\text{EMP} \bowtie \text{PAY}) \rightarrow \text{Site 4}) + \\ cost(\text{PROJ} \rightarrow \text{Site 4})$$

$$= 8 + 8 + 1 = 17$$

$$cost(\text{Alternative 2}) = cost(\text{PAY} \rightarrow \text{Site 1}) + cost((\text{PAY} \bowtie \text{EMP}) \rightarrow \text{Site 4}) + \\ cost(\text{PROJ} \rightarrow \text{Site 4})$$

- 4+8+1-13	Iteration	DIZC	Ditt
D ' ' DO NOW CDI IM	EMP	8	1
Decision : DO NOT SPLIT	PAY	4	2
Character F.C. in the office of?	PROJ	1	3
Step 5: $ES_0$ is the "best".	ASG	10	4

 $4 \pm 9 \pm 1 - 19$ 

Step 6: No redundant transmissions.

Based on join selectivities,  $size(EMP \infty PAY) = size(EMP)$ ,  $size(PROJ \infty ASG) = 2*size(PROJ)$ ,  $size(ASG \infty EMP) = size(ASG)$ 

Siza

Relation

Sita

### Hill Climbing Algorithm

#### Problems:

- Greedy algorithm → determines an initial feasible solution and iteratively tries to improve it
- ② If there are local minimas, it may not find global minima
- If the optimal schedule has a high initial cost, it won't find it since it won't choose it as the initial feasible solution
  Relation

#### Example: A better schedule is

$$PROJ \rightarrow Site 4$$

$$ASG' = (PROJ \bowtie ASG) \rightarrow Site 1$$

$$(ASG' \bowtie EMP) \rightarrow Site 2$$

$$Total cost = 1 + 2 + 2 = 5$$

Based on join selectivities, size(EMP∞PAY) = size(EMP), size(PROJ∞ASG) = 2\*size(PROJ), size(ASG∞EMP) = size(ASG)

EMP PAY

PROJ ASG Size

10

Site

# SDD-1

- Extensive use of semijoins
- Objective function is expressed in terms of total communication time
- The algorithm uses statistics on the database, called database profiles
- Like its predecessor hill-climbing algorithm, the SDD-1 algorithm selects locally optimal strategies

#### Initialization

- Step 1: In the execution strategy (call it ES), include all the local processing
- Step 2: Reflect the effects of local processing on the database profile
- Step 3: Construct a set of beneficial semijoin operations (BS) as follows:

$$BS = \emptyset$$

For each semijoin  $SJ_i$ 

$$BS \leftarrow BS \cup SJ_i \text{ if } cost(SJ_i) < benefit(SJ_i)$$

#### Cost vs. Benefit of Semi-Join

Cost(
$$R \sim_A S$$
) =  
 $T_{MSG} + T_{TR} * size(\Pi_A(S))$ 

Benefit(
$$R \sim_A S$$
) = (1 -  $SF_{\infty}(S.A)$ ) \* size(R) \*  $T_{TR}$ 

#### Iterative Process

- Step 4: Remove the most beneficial  $SJ_i$  from BS and append it to ES
- Step 5: Modify the database profile accordingly
- Step 6: Modify BS appropriately
  - compute new benefit/cost values
  - check if any new semijoin need to be included in BS
- Step 7: If  $BS \neq \emptyset$ , go back to Step 4.

#### Assembly Site Selection

Step 8: Find the site where the largest amount of data resides and select it as the assembly site

#### Postprocessing

Step 9: For each  $R_i$  at the assembly site, find the semijoins of the type

$$R_i \bowtie R_j$$

where the total cost of ES without this semijoin is smaller than the cost with it and remove the semijoin from ES.

Note: There might be indirect benefits.

Step 10: Permute the order of semijoins if doing so would improve the total cost of ES.

# 4

# Example of SDD-1 (SQL)

Select R3. C
From R1, R2, R3
Where R1.A=R2.A
and R2.B=R3.B



relatio n	card	tuple size	relatio n size
R1	30	50	1500
R2	100	30	3000
R3	50	40	2000

	Semi- join	benefit	cost
SJ1	R2∝R1	(1-0.3)*3000=2100	36
SJ2	R2∝R3	(1-0.4)*3000=1800	80
SJ3	R1∝R2	(1-0.8)*1500=300	320
SJ4	R3∝R2	(1-1)*2000=0	400

Attribute	SFsj	Size (Πattribute)
R1.A	0.3	36
R2.A	0.8	320
R2.B	1.0	400
R3.B	0.4	80

# Example of SDD-1 (DB Profile)

relation	card	tuple size	relation size
R1	30	50	1500
R2	100	30	3000
R3	50	40	2000

relation	card	tuple size	relation size
R1	30	50	1500
R2	30	30	900
R3	50	40	2000

Attribute	SFsj	Size
		(Пattribute)
R1.A	0.3	36
R2.A	0.8	320
R2.B	1.0	400
R3.B	0.4	80

Attribute	SFsj	Size
		(Пattribute)
R1.A	0.3	36
R2.A	0.24	96
R2.B	0.3	120
R3.B	0.4	80

- Initialization
- Selection of beneficial semijoins
- Assembly site selection
- Postoptimization
- Like its predecessor hill-climbing algorithm, the SDD-1 algorithm selects locally optimal strategies
  - It ignores the higher-cost semijoins which would result in increasing the benefits and decreasing the costs of other semijoins. Thus this algorithm may not be able to select the global minimum cost solution



# 7.5 Local Optimization



#### Input: Best global execution schedule

- Select the best access path by all the site
- Use the centralized optimization techniques



# 7.6 Conclusion

### Conclusion

- Objective function
  - Communication cost
  - Local processing cost
- Input
  - Database statistics
  - Formulas
- Join vs. Semijoin

### Conclusion

- Larger set of queries
  - optimization only on select-project-join queries
  - also need to handle complex queries (e.g., unions, disjunctions, aggregations and sorting)
- Optimization cost vs execution cost tradeoff
  - heuristics to cut down on alternatives
  - controllable search strategies
- Optimization/reoptimization interval
  - extent of changes in database profile before reoptimization is necessary

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