

DB 2

14 – Query Optimization

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1 : One Query — Millions of Plans

Q: Given a SQL query Q , what is *the optimal* (a reasonable)¹ plan to evaluate it? — **A:** It depends:

- Can we **simplify** (flatten, unnest) Q ?
- How can we **access the tables** referenced in Q ?
- How do **CPU and (sequential, random) I/O cost** compare?
- What is the **selectivity of the predicates** used in Q ?
- Which plan **operator implementations** are applicable?
- Can we **regroup/reorder the joins** in Q ?

¹ Here: focus on reducing the overall query evaluation time. The optimum is, generally, not reached.

Excerpt of the TPC-H Benchmark (at Scale Factor SF)

<u>o_orderkey</u>	o_custkey	o_totalprice	o_clerk	...
o	c			

orders ($\approx SF \times 1.5 \times 10^6$ rows)

<u>l_orderkey</u>	<u>l_linenum</u>	l_partkey	l_quantity	l_extendedprice	...
o					

lineitem ($\approx SF \times 6 \times 10^6$ rows)

<u>c_custkey</u>	c_name	c_acctbal	c_nationkey	...
c			n	

customer ($\approx SF \times 150000$ rows)

<u>n_nationkey</u>	n_name	n_regionkey	...
n		r	

nation (25 rows)

<u>r_regionkey</u>	r_name	...
r		

region (5 rows)



Q_{14} : Three-Way Join Against a TPC-H Instance

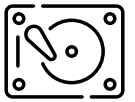
Price and quantity of parts ordered by customer #001:

```

SELECT 1.1_partkey, 1.1_quantity, 1.1_extendedprice
FROM   lineitem AS l JOIN orders AS o      -- } l ⋈ o
      ON (l.1_orderkey = o.o_orderkey)    -- }
      JOIN customer AS c                  -- } ⋈ c
      ON (o.o_custkey = c.c_custkey)      -- }
WHERE  c.c_name = 'Customer#001';

```

- Above SQL syntax suggests the **join order** $(l \bowtie o) \bowtie c$.
- Commutativity and associativity of \bowtie enable the RDBMS to **reorder** the joins—based on *estimated evaluation costs*.
 - ... unless we insist on the syntactic order. 🧐



2 : Pre-Processing: Query Normalization

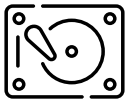
Transform the input SQL query such that it features **SELECT-FROM-WHERE** (SFW) blocks of the following shape:

```

SELECT [ DISTINCT ] e, ..., e
FROM       $\Delta$ , ...,  $\Delta$ 
[ WHERE    p AND ... AND p ]
[ GROUP BY g, ..., g
  [ HAVING  p AND ... AND p ] ]
[ ORDER BY o, ..., o ]
[ OFFSET   n ]
[ LIMIT    m ]
--  $\Delta \equiv$  base table or (query)
-- p  $\equiv$  predicate in DNF
-- { e, p, g, o  $\equiv$ 
--   atomic expression or
--   scalar (subquery)
-- { n, m  $\equiv$  integer literal

```

- Query clauses in [...] may be missing.

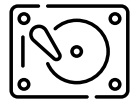


3 | Pre-Processing: Query Unnesting

Nested SQL queries suggest a (naïve, inefficient) nested-loop-style evaluation strategy. Consider:

<pre> SELECT c.c_name FROM customer AS c, ⚠ { (SELECT n.n_nationkey, n.n_name { FROM nation AS n) AS t WHERE c.c_nationkey = t.n_nationkey AND strpos(c.c_address, t.n_name) > 0 </pre>		<pre> SELECT o.o_orderkey FROM orders AS o WHERE o.o_custkey IN ⚠ { (SELECT c.c_custkey { FROM customer AS c WHERE c.c_name = '...') </pre>
---	--	--

- 💡 If possible, **unnest** ⚠ queries and “inline” into parent query \Rightarrow ⚠ can participate in join reordering.



Pre-Processing: Query Unnesting

Perform **query unnesting** on the level of

- the operator-based plan representation of the query,² or
- the internal AST representation of SQL. Re ②:

<pre> SELECT e₁ FROM q₁, ..., q_i WHERE p₁ AND e₂ IN (SELECT e₃ FROM q_{i+1}, ..., q_n WHERE p₃) </pre>	\cong	<pre> SELECT DISTINCT e₁ FROM q₁, ..., q_i, q_{i+1}, ..., q_n WHERE p₁ AND e₂ = e₃ AND p₃ </pre>
--	---------	--

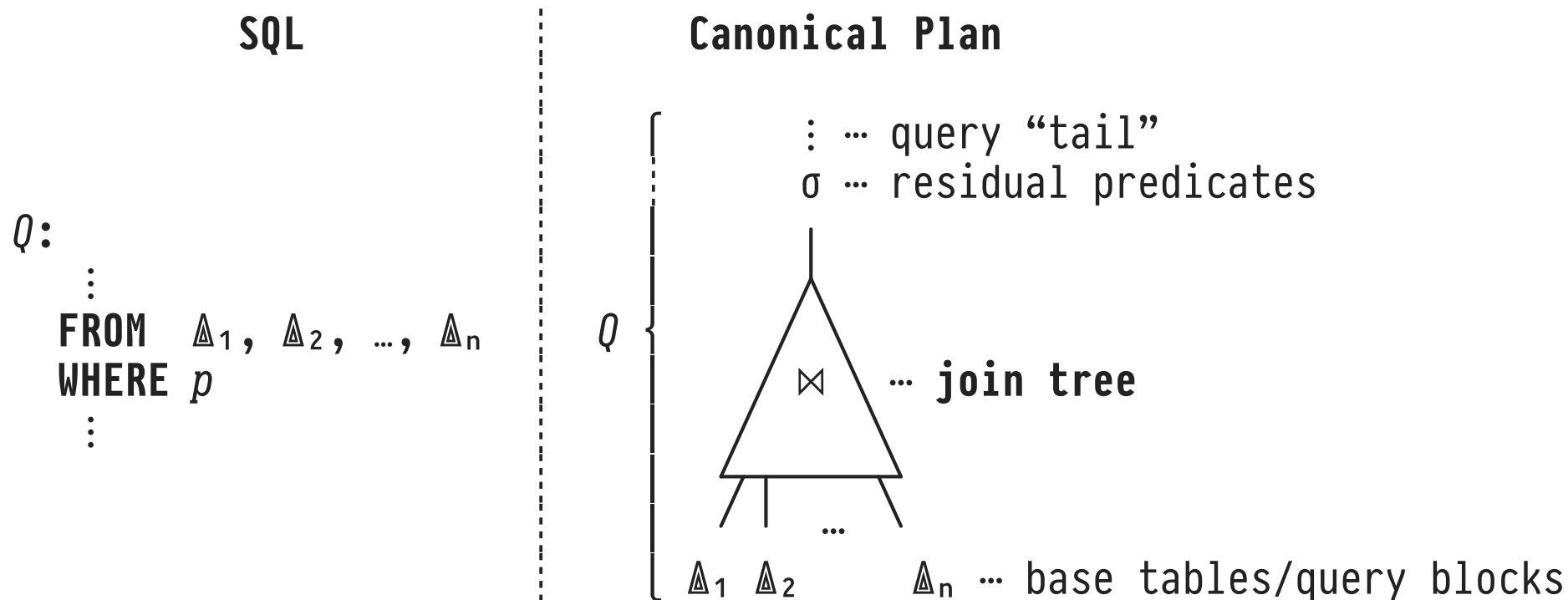
* Precondition: e_1 is key in the left-hand side query

² See *Unnesting Arbitrary Queries*, Thomas Neumann, Alfons Kemper. BTW 2015, Hamburg, Germany.



4 : Join Tree Optimization

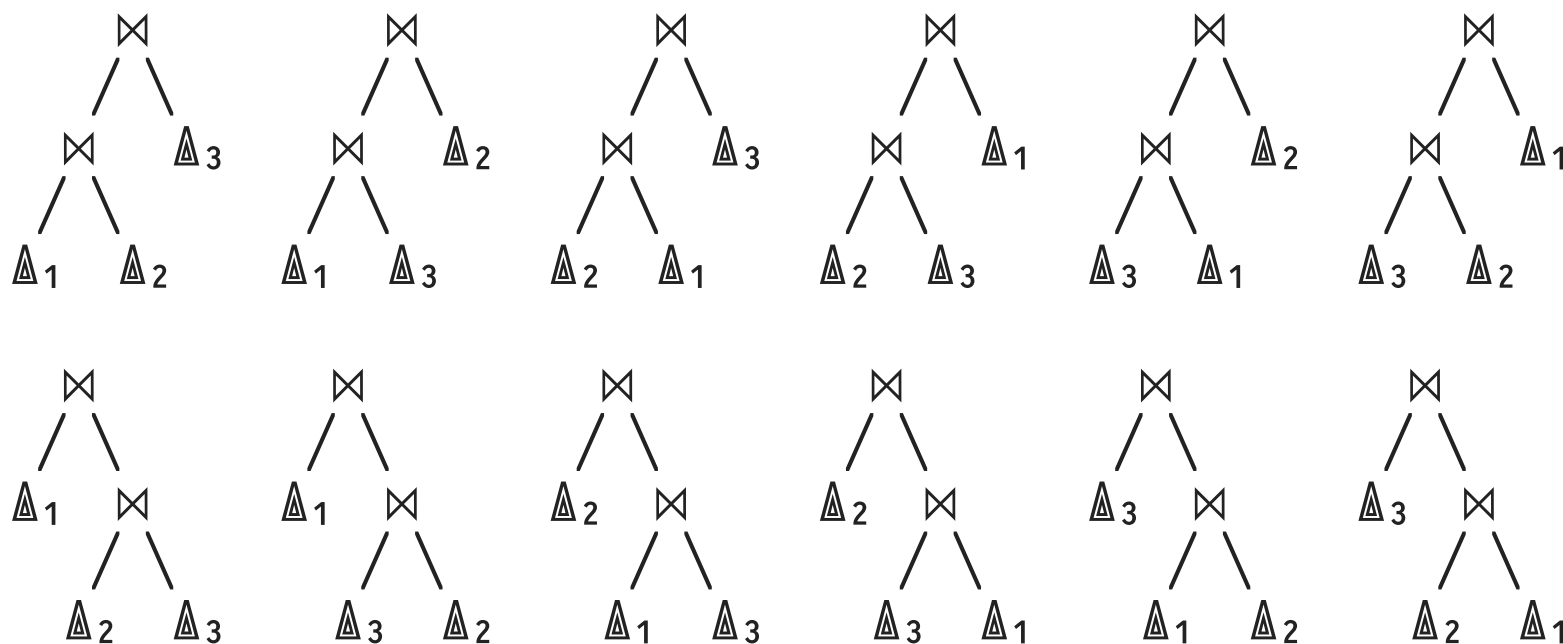
Processing a SQL query Q starts out with its **FROM** and **WHERE** clauses which describe a **join tree** over Q 's inputs:





Join Tree Optimization

Given n join inputs, the number of possible **join tree shapes** is *huge*. Consider $n = 3$:

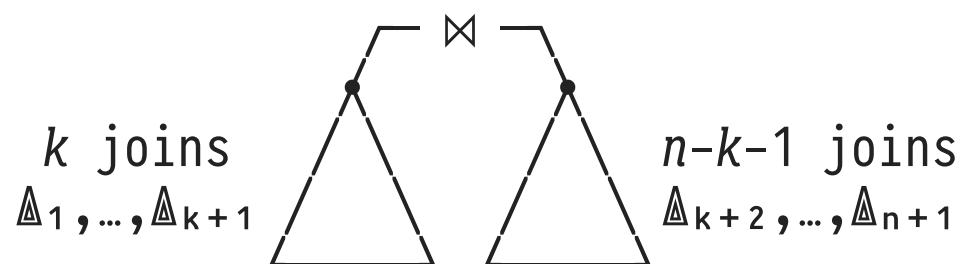


- Shapes based on associativity and commutativity of \bowtie .



How Many Possible Join Trees Are There?

1. A join of $n+1$ inputs Δ requires n binary joins. The root \bowtie combines subtrees of k and $n-k-1$ joins ($0 \leq k \leq n-1$):³



of join tree shapes:

$$C_n = \sum_{k=0}^{n-1} C_k \times C_{n-k-1}$$

2. Orderings of the Δ at the join tree leaf level: $(n+1)!$.
3. Join algorithm choices (a available algorithms): a^n .

³ C_n are the *Catalan numbers*, the number of ordered binary trees with $n+1$ leaves. $C_0 = 1$.



How Many Possible Join Trees Are There?

Number of possible join trees given n binary joins with $a = 3$ implementation choices:

# of $\Delta (n+1)$	C_n	# of join trees
2	1	6
3	2	108
4	5	3240
5	14	136080
6	42	7384320
7	132	484989120
8	429	37829151360
9	1430	3404623622400
10	4862	347271609484800

- A search space of this size is impossible to fully explore for any query optimizer.



Join Plan Generation Through Dynamic Programming

- **Problem:** Find optimal query plan $opt[\{\Delta_1, \dots, \Delta_n\}]$ that joins n inputs $\Delta_1, \dots, \Delta_n$.
 1. **Iteration 1:** For each Δ_j , find and memorize **best 1-input plan** $opt[\{\Delta_j\}]$ that accesses Δ_j only.
 2. **Iteration $k > 1$:** Find and memorize **best k -input plans** that join $k \leq n$ inputs by combining (for $1 \leq i < k$)
 - the best i -input plans and \setminus simple lookups in
 - the best $(k-i)$ -input plans. $\int opt[\cdot]$ memo 👍



Bottom-Up Dynamic Programming ($n = 3$)

- k Possible k -input Access/Join Plans** if Δ_i is complex
- 1** $opt[\{\Delta_1\}] \leftarrow \text{prune}(\{\text{Seq Scan } \Delta_1, \text{Index Scan } \Delta_1, \text{Bitmap Scan } \Delta_1, \overbrace{\Delta_1}^{\text{complex}}\})$
 $opt[\{\Delta_2\}] \leftarrow \text{prune}(\{\text{Seq Scan } \Delta_2, \text{Index Scan } \Delta_2, \text{Bitmap Scan } \Delta_2, \Delta_2\})$
 $opt[\{\Delta_3\}] \leftarrow \text{prune}(\{\text{Seq Scan } \Delta_3, \text{Index Scan } \Delta_3, \text{Bitmap Scan } \Delta_3, \Delta_3\})$
-
- 2** $opt[\{\Delta_1, \Delta_2\}] \leftarrow \text{prune}(opt[\{\Delta_1\}] \otimes opt[\{\Delta_2\}])$
 $opt[\{\Delta_1, \Delta_3\}] \leftarrow \text{prune}(opt[\{\Delta_1\}] \otimes opt[\{\Delta_3\}])$
 $opt[\{\Delta_2, \Delta_3\}] \leftarrow \text{prune}(opt[\{\Delta_2\}] \otimes opt[\{\Delta_3\}])$
-
- 3** $opt[\{\Delta_1, \Delta_2, \Delta_3\}] \leftarrow \text{prune}(opt[\{\Delta_1\}] \otimes opt[\{\Delta_2, \Delta_3\}] \cup$
 $opt[\{\Delta_2\}] \otimes opt[\{\Delta_1, \Delta_3\}] \cup$
 $opt[\{\Delta_3\}] \otimes opt[\{\Delta_1, \Delta_2\}])$

$\text{prune}(P) \equiv \text{best (= minimal cost + interestingly ordered) plans in set } P$

$$l \otimes r \equiv \{l \bowtie^{n1} r, r \bowtie^{n1} l, l \bowtie^{mj} r, r \bowtie^{mj} l, l \bowtie^{hj} r, r \bowtie^{hj} l\}$$



Join Plan Generation (Notes)

- **Access plan choices** (*access*(·)):
 - Consider sequential/index scans if Δ is a base table, otherwise simply consume Δ 's rows.
- **Join plan choices** ($_ \otimes _$):
 - Considers all viable join algorithms (given θ , available indexes, ...) and left/right input orders.
- **Principle of Optimality** (*prune*(·)): A globally optimal plan is built from optimal subplans. Thus:
 - 💡 For each subset of $\{\Delta_1, \dots, \Delta_n\}$, memorize in *opt*[·]
 1. ... its overall best plan and
 2. ... its best plan satisfying each **interesting order**.



(Bushy) Join Plan Generation: Pseudo Code

```

JoinPlan( $\{\Delta_1, \dots, \Delta_n\}$ ):
  foreach  $p \in \{\Delta_1, \dots, \Delta_n\}$  } 1-input plans
  |  $opt[\{p\}] \leftarrow prune(access(p))$ ;

  for  $k$  in  $2, \dots, n$  }  $k$ -input plans
  |   foreach  $S \subseteq \{\Delta_1, \dots, \Delta_n\}$  with  $|S| = k$  } enumerate subsets
  |   |    $opt[S] \leftarrow \phi$ ;
  |   |   foreach  $T \subset S$  with  $T \neq \phi$   $\Join^a \setminus$ 
  |   |   |    $opt[S] \leftarrow opt[S] \cup \{ opt[T] \Join^a opt[S \setminus T] \}$ ;
  |   |   |    $opt[S] \leftarrow prune(opt[S])$ ;

  return  $opt[\{\Delta_1, \dots, \Delta_n\}]$ ;

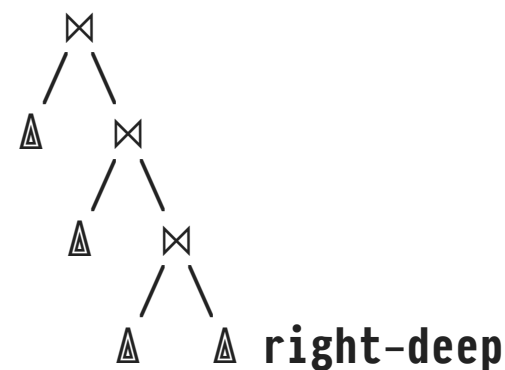
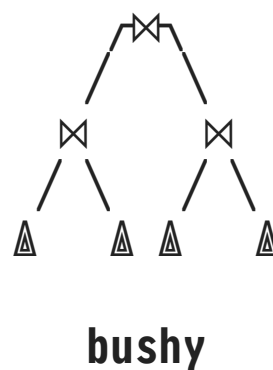
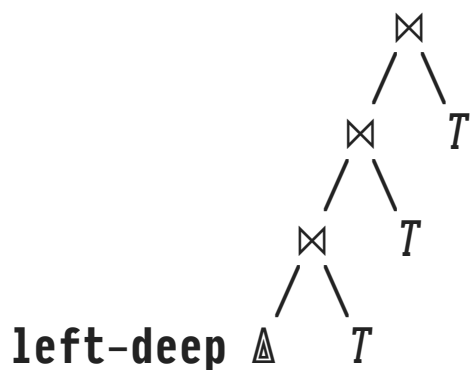
```

- $access(\cdot)$, $prune(\cdot)$ defined as above,
 $\Join^a \setminus$ builds all join algorithm choices ($a \in \{nl, mj, hj\}$).



Reducing the Search Space

- Avoid generating costly **Cartesian products**: don't form joins between inputs w/o join predicate (`_ θ _ = true`).
- Generate **left-deep** join plans only: right join input (NL \bowtie : inner input) is a scan over base table T .
 - Admits use of Index Nested Loop Join.
 - Straightforward Volcano-style execution (reset inner).





5 : Estimating Plan Cost

The query optimizer explores the vast plan search space to find the **optimal** (“best”, “cheapest”) plan.

- Typically, RDBMSs measure **plan cost** in terms of *total execution time* (time until last result row delivered).
- These total plan costs are **estimated** *before* plan execution begins (`EXPLAIN: ... cost=c1..c2← ...`).
- A **cost model**—measured in abstract “space\$”—reflects the true costs (measured in *ms*, CPU time, # I/O ops, ...) of plans p_1 , p_2 :

$$\text{space}\$(p_1) < \text{space}\$(p_2) \Rightarrow \text{true cost}(p_1) < \text{true cost}(p_2)$$



PostgreSQL: Plan Cost

EXPLAIN shows estimated costs (unit: space\$) and cardinalities (# of rows):

QUERY PLAN	
startup cost	total cost
	↓
Hash Join	(cost=299.00..15443.31 rows=505183 width=50)
⋮	↑
	cardinality

- **run cost** $\stackrel{\text{def}}{=} \text{total cost} - \text{startup cost}$ ⁴ (not shown).
- Optimizer decisions are based on estimated **total cost**.

⁴ To implement `set enable_<op> = off`, PostgreSQL sets the operator's **startup cost** to 10^{10} ($\equiv \infty$).



Cost Model Configuration

Model Configuration	Default	Description
<code>seq_page_cost</code>	1.0	I/O cost of one sequential page access
<code>random_page_cost</code>	4.0	I/O cost of one random page access
<code>cpu_tuple_cost</code>	0.01	CPU cost to process a heap file row
<code>cpu_index_tuple_cost</code>	0.005	CPU cost to process an index leaf entry
<code>cpu_operator_cost</code>	0.0025	CPU function/operator evaluation cost
<code>parallel_tuple_cost</code>	0.1	Cost of passing one row worker→leader
<code>parallel_setup_cost</code>	1000.0	Cost of spawning a parallel worker

- Parameters are configurable:
 - Seek cost, thus `random_page_cost` » `seq_page_cost`. But...
 - ... if DB fits in RAM, `random_page_cost` = `seq_page_cost` may be more appropriate.



Cost of Seq Scan 1

Given an occurrence of **Seq Scan** with arguments

- *in*: input table,
- *pred*: (optional) filter predicate on *in*,
- *expr*: **SELECT** clause expression(s),

how does PostgreSQL derive *startup_cost* and *total_cost*?

<i>in</i>	QUERY PLAN	<i>total_cost</i>
Seq Scan on public.indexed i	(cost=0.00..22.75 rows=100 width=4)	
Output: (a + 1)	↑ <i>expr</i>	↑
Filter: (i.a <= 100) ← <i>pred</i>	<i>startup_cost</i>	<i>#rows(out)</i>



Cost of Seq Scan 2

Cost calculation depends on the following parameters, mostly available in PostgreSQL's internal `pg_*` meta data tables:

Parameter	Description	Available as...
<code>#rows(<i>in</i>)</code>	# rows (cardinality) of table <i>in</i>	<code>pg_class.reltuples</code>
<code>#pages(<i>in</i>)</code>	# pages in heap file of <i>in</i>	<code>pg_class.relpages</code>
<code>sel(<i>pred</i>)</code>	selectivity of filter <i>pred</i> ⁵	see below

- Meta data like `#rows(in)`, `#pages(in)` and others are updated whenever the system performs an `ANALYZE` run on table *in*.
- Predicate selectivity `sel(pred)` is estimated based on sampled table data and the syntactic structure of *pred*.

⁵ `sel(pred)` $\in \{0, \dots, 1\}$ with `sel(pred) = 0` \equiv no row satisfies filter *pred*.



Cost of Seq Scan 3

\nwarrow typically = 0 \swarrow
startup_cost $\stackrel{\text{def}}{=}$ $\text{startup_cost}(\text{pred}) + \text{startup_cost}(\text{expr})$

$\text{cpu_run_cost} \stackrel{\text{def}}{=}$ $\underbrace{\#rows(in) \times (\text{cpu_tuple_cost} + \text{run_cost}(\text{pred}))}_{\text{decode heap row}} + \underbrace{\#rows(in) \times \text{sel}(\text{pred}) \times \text{run_cost}(\text{expr})}_{\text{evaluate filter}}$
 $\quad \quad \quad = \#rows(out) \quad \quad \quad \text{evaluate SELECT clause}$

disk_run_cost $\stackrel{\text{def}}{=}$ $\underbrace{\#pages(in) \times \text{seq_page_cost}}_{\text{sequentially read entire input heap file}}$

total_cost $\stackrel{\text{def}}{=}$ $\text{startup_cost} + \underbrace{\text{cpu_run_cost} + \text{disk_run_cost}}_{= \text{run_cost}}$



Cost of Index Scan 1

Modeling the cost for an **Index Scan** has to reflect that *two* data structures (heap file & B+Tree) are involved:

<i>idx</i>	<i>in</i>	QUERY PLAN
↓	↓	
Index Scan using indexed_a on indexed i (cost=0.42..443.12 rows=10885 ...		
Output: (c + '1'::numeric) ← <i>expr</i>		↑
Index Cond: (i.a ≤ 10000) ← <i>pred</i>		#rows(<i>out</i>)

The model separately accounts for

1. the B+Tree descent (startup of the **Index Scan**),
2. the index leaf level scan, and
3. heap file access (clustered *vs.* non-clustered).



Cost of Index Scan

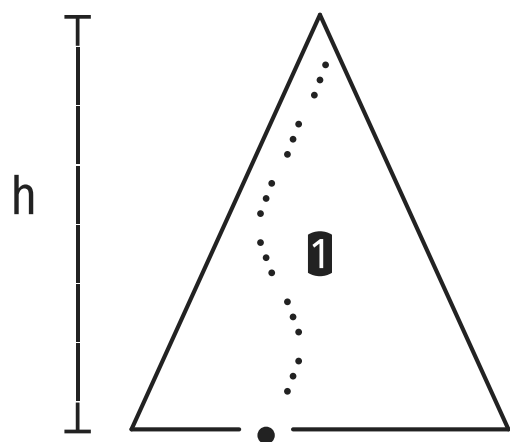
Cost model parameters:

Parameter	Description	Available as...
<code>#rows(<i>in</i>)</code>	# rows (cardinality) of table <i>in</i>	<code>pg_class.reltuples</code>
<code>#pages(<i>in</i>)</code>	# pages in heap file of <i>in</i>	<code>pg_class.relpages</code>
<code>sel(<i>pred</i>)</code>	selectivity of filter <i>pred</i>	see below
<code>h(<i>idx</i>)</code>	height of B+Tree <i>idx</i>	<code>bt_metap(.)</code>
<code>#rows(<i>idx</i>)</code>	# leaf entries in index <i>idx</i>	<code>pg_class.reltuples</code>
<code>#pages(<i>idx</i>)</code>	# pages in leaf level of <i>idx</i>	<code>pg_class.relpages</code>
<code>corr(<i>idx</i>)</code>	\approx clustering factor for index <i>idx</i>	<code>pg_stats.correlation</code>

- `corr(idx)` $\in \{-1.0, \dots, 1.0\}$ characterizes how much the physical orderings of index leaves and heap file deviate.
 - After `CLUSTER in ON idx`, we have `corr(idx) = 1.0`.



Cost of Index Scan **E** (B+Tree Descent)



- B+Tree height $h = \log_{2 \times o}(\#rows(idx))$

⇒ # of key comparisons during B+Tree descent **1**:

$$\underbrace{[\log_2(2 \times o) \times h]}_{\text{binary search in inner B+Tree node with fan-out } F = 2 \times o} = [\log_2(\#rows(idx))]$$

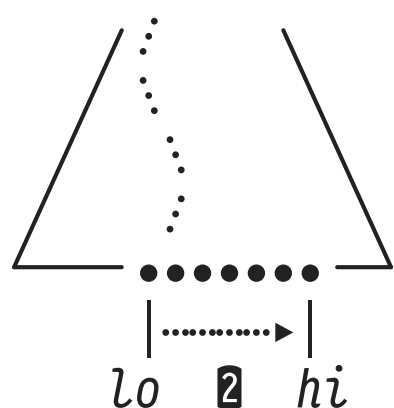
binary search in inner B+Tree
node with fan-out $F = 2 \times o$

$$\begin{aligned} \text{startup_cost} &\stackrel{\text{def}}{=} \text{startup_cost}(pred) + \text{startup_cost}(expr) \\ &+ \underbrace{([\log_2(\#rows(idx))])}_{\text{B+Tree descent}} + \underbrace{(h + 1)}_{\text{index node processing}} \times \underbrace{50}_{\text{index node processing}} \times \text{cpu_operator_cost} \end{aligned}$$



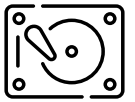
Cost of Index Scan ④ (Leaf Level Scan)

The index leaf level (sequence set) scan ② incurs CPU as well as I/O cost that contribute to the overall **run_cost**:



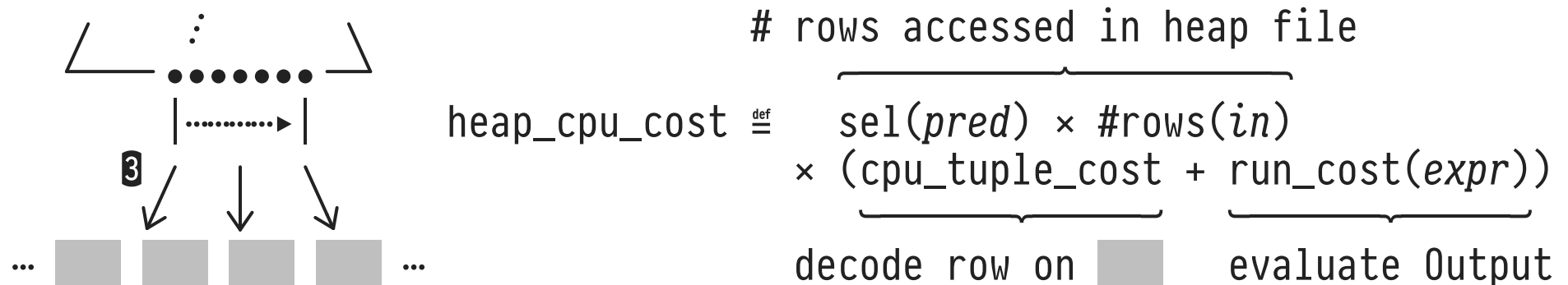
$$\text{index_cpu_cost} \stackrel{\text{def}}{=} \underbrace{\text{sel}(\text{pred}) \times \text{\#rows}(\text{idx})}_{\text{\# rows in scanned range } |\text{.....}\rangle|} \times \underbrace{(\text{cpu_index_tuple_cost})}_{\text{decode index leaf entry}} + \underbrace{\text{run_cost}(\text{pred})}_{\text{evaluate } \leq \text{hi}}$$

$$\text{index_IO_cost} \stackrel{\text{def}}{=} \underbrace{[\text{sel}(\text{pred}) \times \text{\#pages}(\text{idx})]}_{\text{\# of pages } \bullet \text{ in scanned range}} \times \underbrace{\text{random_page_cost}}_{\text{B+Tree leaves not clustered}}$$



Cost of Index Scan 5 (Heap File Access)

Heap file accesses ③ incur additional CPU and I/O costs (no I/O cost if we perform an **Index Only Scan**):

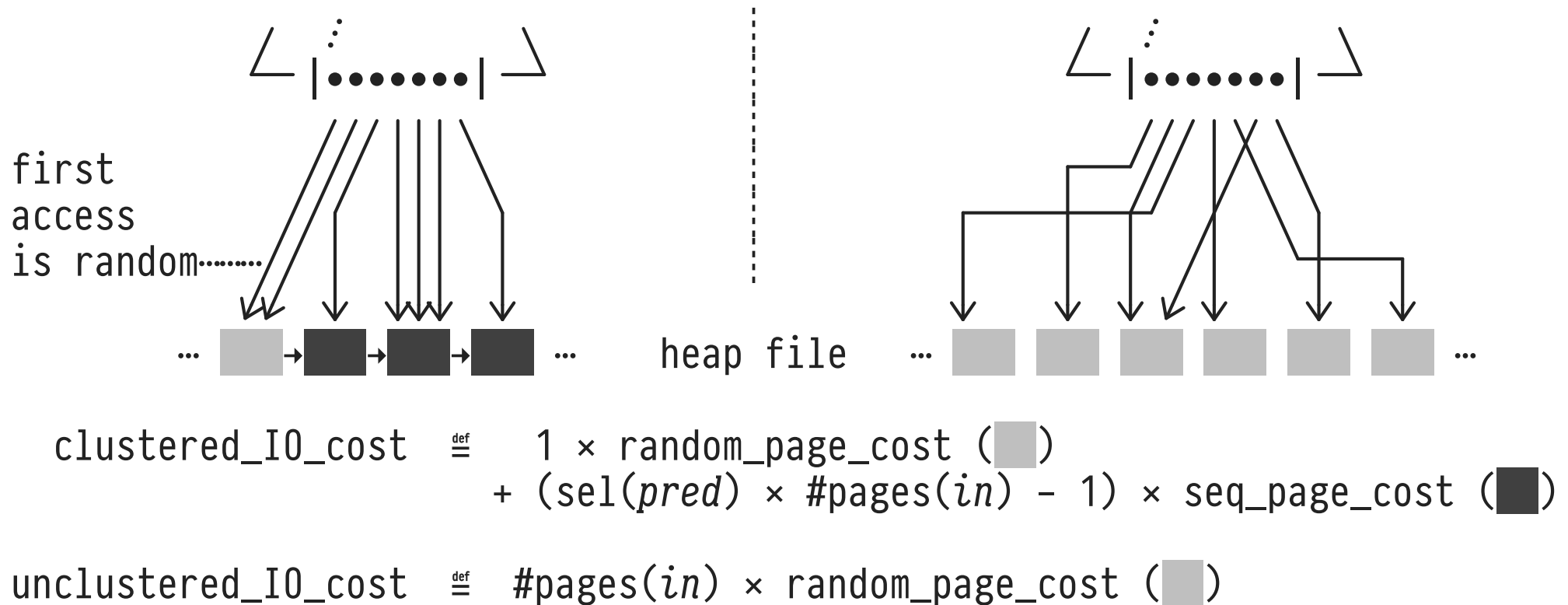


- The more **clustered** the index, the cheaper the heap I/O. Linearly interpolate between the clustered and non-clustered scenarios:

$$\text{heap_IO_cost} \stackrel{\text{def}}{=} \text{unclustered_IO_cost} + \text{corr}(\text{idx})^2 \times (\text{clustered_IO_cost} - \text{unclustered_IO_cost})$$

$\text{corr}(\text{idx}) \approx \text{clustering factor} \in \{0, \dots, 1\}$

Cost of Index Scan ([Non-]Clustered Heap File Access)



$$\text{total_cost} \stackrel{\text{def}}{=} \text{startup_cost} + \text{index_cpu_cost} + \text{index_I0_cost} + \text{heap_cpu_cost} + \text{heap_I0_cost}$$



Index Correlation (Clustering Factor)

Given ordered index idx over column A with values $a_1 \leq a_2 \leq \dots \leq a_n$, where $pos(a_i) \in \{1, \dots, n\}$ gives the position of a_i in the heap file for A .⁶

- **Index Correlation** $corr(idx) \in \{-1, \dots, 1\}$ measures how far $[pos(a_1), \dots, pos(a_n)]$ deviates from $[1, \dots, n]$, i.e., idx 's clustering degree:

$$corr(idx) = \frac{n \times (\sum_{i=1 \dots n} i \times pos(a_i)) - (\sum_{i=1 \dots n} i)^2}{n \times (\sum_{i=1 \dots n} i \times i) - (\sum_{i=1 \dots n} i)^2}$$

⁶ After `CLUSTER <table> USING idx`, we have $pos(a_i) = i$ and thus $corr(idx) = 1$.