

Data Base 2 Course Notes

Elia Ravella

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Contents

I	Transactions	2
1	Definition and Usages of a Transaction	2
2	Cuncurrency Control	3
2.1	Types of Anomalies	3
2.2	Cuncurrency Theory	3
2.2.1	Serial Schedules	4
2.2.2	View Serializability VSR	4
2.2.3	Conflict Serializability CSR	4
2.2.4	Two Phase Locking 2PL	5
2.2.5	Timestamping	6
3	Deadlock Detection - Obermarck Algorithm	7
3.1	Applicability	7
3.2	The Algorithm	7
II	Memory Management	8
4	Memory Layers	8
4.1	Blocks, Tuples	8
5	Access Structures	8
5.1	Data Structures Recap	8
5.1.1	Sequential Structures	9
5.1.2	Hash Based Structures	9
5.1.3	Tree Based Structures	9
6	Query Optimizer	9
6.1	Costs of Various Access Patterns	10
6.1.1	Sequential Scan	10
6.1.2	Lookup	10
6.1.3	Using Statistics	11
6.1.4	Join	11
7	JPA	12
7.1	Objects and Relations	12
7.2	JPA Main Features and Concepts	12
7.2.1	Entity, Managers, Managed Entities	13
7.2.2	Relationships	15
8	Active Databases - Triggers	16

Part I

Transactions

1 Definition and Usages of a Transaction

A transaction is usually defined as an *elementary, atomic unit of work performed by an application*. Transactional systems are those which support transactions. They usually implement a couple of primitives (such as `begin-transaction` and `commit-transaction` or `abort-transaction`) to manage the lifecycle of such an object.

ACID Properties

1. Atomicity (all or nothing)
2. Consistency (only valid states are reachable)
3. Isolation (transaction should overlap only if they not interfere. The output of concurrent transactions should be the same as the output of the very same transactions executed serialized)
4. Durability (changes are permanent)

are the four main properties the are necessary to avoid inconsistent data in a database. Transactional systems tries to embody all these properties in their transaction management mechanism.

DBMS Modules A Data Base Management System is composed of several interconnected modules:

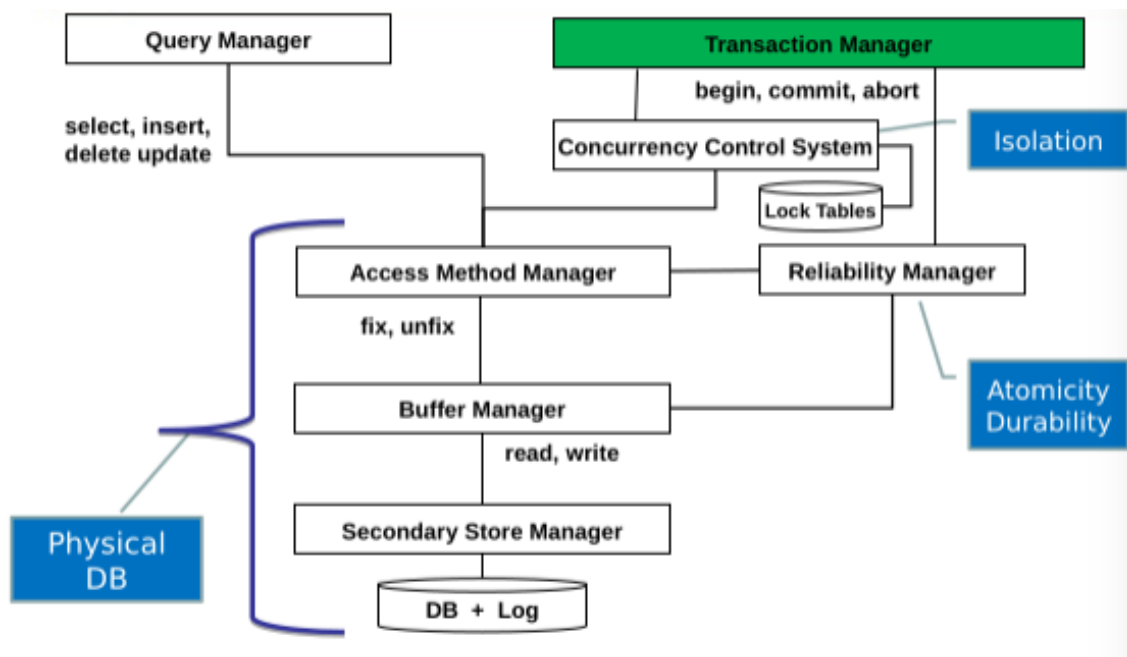


Figure 1: DBMS modules

How do those modules interact in order to deliver transactions successfully and which ACID properties they relate to?

- Atomicity and Durability are handled by the *reliability manager*
- Isolation is guaranteed by the *cuncurrency control system*
- Consistency is managed by the *integrity control system* at query execution time. Obviously in this process the DDL query compiler plays an important role

2 Cuncurrency Control

DBMSs need cuncurrency. It's pretty straightforward how having multiple transactions executing at the same time increases significantly the throughput of the system. On the other hand, managing cuncurrent systems is a mess. If we add 99% of cases a DBMS is also distributed across machines, the mess becomes huge.

2.1 Types of Anomalies

- **Lost Update:** two transaction write a value in a cell that depends from the previous value (think about an update) but they *both read the same initial value* and so one of the two updates is lost. From the system point of view, the pathologic serie is

`r1(x) r2(x) w1(x) w2(x)`

the first write on `x` is "lost"¹.

- **Dirty Read:** when a transaction rollbacks, it's invalidating all the data it has written or updated. All the reads performed on the values changed² by the aborted transaction are *dirty* in the sense that they'e read an invalid value, so they must be reissued.

`r1(x) w1(x) r2(x) abort1`

- **Nonrepeatable Read:** pretty straightforward, this anomaly pops out when two consecutive reads of a value (from a transaction) are allowed to interleave with a write on *that value*. The two read operation will read two different values from the same cell.

`r1(x) r2(x) w2(x) r1(x)`

- **Phantom Update:** Same concept of the Nonrepeatable Read, but this time only a portion of the data read in the first place is modified. The sequence of operation is the same, still.
- **Phantom Insert:** identical as the Nonrepeatable Read, but this time is new data.

The last three anomalies are conceptually the same thing: someone modifies data thet I've previously read before I've finished using it. This is also similar to the concept or the Lost Update anomaly. In fact, these are all cause by the classic cuncurrent model that's implemented on machines, with the assumptions that writes and reads are so basic operations that they can be easily assumed atomic on their own.

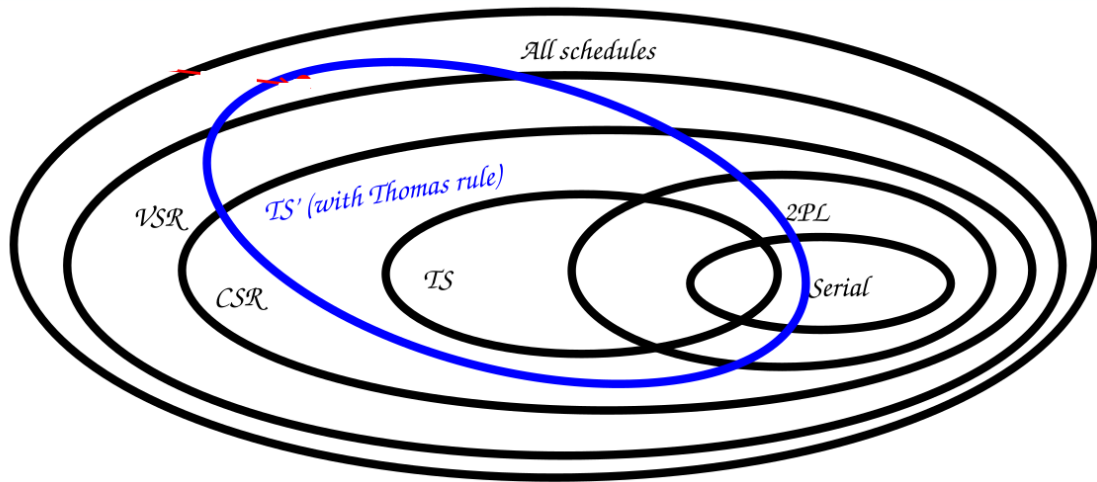
2.2 Cuncurrency Theory

We define as *Schedule* a serie of basic operations³ performed by *one or more processes* on *one or more values*. We can define several sets of schedules, discriminating on which properties each one respects. The overall set is

¹we're assuming that the two updates are dependent.

²written

³reads, writes, aborts, commits



So, there are a lot of sets. Let's analyze all of them

2.2.1 Serial Schedules

Serial schedules are the simplest of all. There's no interleaving of execution threads, all processes start and complete all their operation at once. These kind of schedules do not cause anomalies, but they also *are not concurrent*, so they're useless.

Serializability We love serial schedules from a theoretical point of view. So, a good measure to say if a random schedule is "good" or "bad" for our system can be if we can somehow reduce it to a serial equivalent. To do so, we introduce the concept of Serializability, which is a property that holds in concurrent schedules that can be reduced to a serial one without affecting the initial and final states of the DB. We will define less strict serializability equivalence relations for less strict concurrent models.

2.2.2 View Serializability VSR

This set of schedules comprises all the schedules that are view equivalent to a serial one. Two schedules are view equivalent if

1. They have the same operations
2. The read operations depends from the same write operations
3. All variables in the schedules have the same last write operation (same final write property). This can also be rephrased as "the final value of each object is written by the same transaction as if the transactions were executed serially in some⁴ order"

How to check if a schedule is VSR? Simply construct the serial equivalent of the schedule (if S is composed of transactions T_1, T_2, T_3 then the possible serials are $\{T_1, T_2, T_3\}$ $\{T_1, T_3, T_2\}$ $\{T_2, T_3, T_1\}$ $\{T_2, T_1, T_3\}$ $\{T_3, T_2, T_1\}$ $\{T_3, T_1, T_2\}$) and verify if there's one among those that produces the same state change.

2.2.3 Conflict Serializability CSR

Determining if a random query is view serializable is a NP complete problem. We can redefine the equivalence relation we want in order to obtain a reduced set of possible valid schedules that are easier to verify.

A schedule is Conflict Serializable if it's conflict equivalent to a serial schedule.

Two schedules are conflict equivalent if

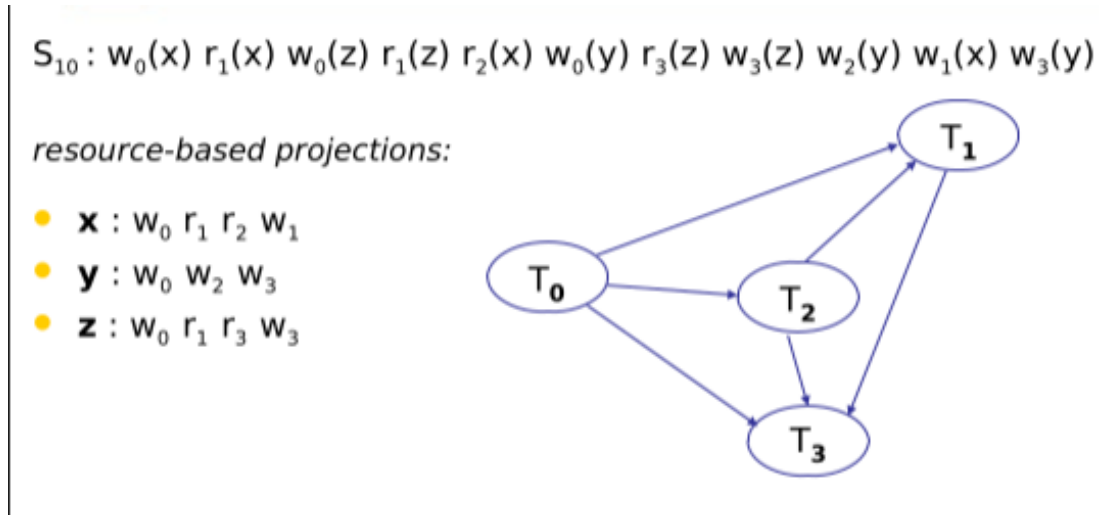
1. They have the same operations

⁴that can also differ from the order they've been issued.

2. All conflicts⁵ occur in the same order

CSR is a proper subset of VSR. This means that all CSR schedule are also view-serializable, but the viceversa does not hold.

To verify the conflict serializability of a schedule we can resort to a *conflict graph*: it's a graph that has as nodes the transactions, and the edges connect transactions which depends one on another (from the variables point of view)



The conflict graph is acyclic if and only if the schedule is conflict serializable. If so, it also shows the serial conflict equivalent schedule.

2.2.4 Two Phase Locking 2PL

It's not realistic to have all the basic read and write operation before the actual execution of the transaction. CSR and VSR checking methods operates on the *history of operations* but we cannot know them a priori (in most cases). We need methods that can react at the arrival of new operations on line. One of this methods exploits a well known technique of concurrency control, locking the resources under request (by a thread / transaction) until the desired operation is finished. The locking mechanism in DBMS uses two locks:

1. Read lock: a transaction requests a read lock when it needs only to read a value from a resource
2. Write lock: pretty auto esplicative

Different lock status are possible, even more locks can be introduced⁶.

2PL is a technique that *imposes a constraint on the usage of locks* in order to provide view (and conflict) serializability. It basically states that **a transaction cannot require any more lock after releasing one**. The set of schedules that respect this policy is called 2PL and is a *proper subset of CSR*.

2PL (being a proper subset of CSR, also) protects against

- Lost update anomalies
- Phantom update anomalies
- Phantom insert anomalies

Strict 2PL A proper subset of 2PL that also protects from dirty reads is the "strict" version of 2PL, that restricts the shrinking phase (the moment when a transaction releases the locks) to a single moment. This way, there is no chance that a modified variable by a transaction could be read before the commit or abort of the transaction itself.

⁵conflicts are defined as "couples of operations on the same resource that contain at least a write operation"

⁶Like update lock, that is a read lock "with the intention of writing".

Other Locking Mechanism A classic problem of lock based resource control is the presence of deadlocks. To avoid so, we can introduce the so called "Update lock" to the pool of possible locking options: the new status-request table will look like

Type of request	Resource status		
	Shared	Update	Exclusive
Shared	Granted	Granted	Denied
Update	Granted	Denied	Denied
Exclusive	Denied	Denied	Denied

Another possible locking mechanism (that addresses the problem of waiting time when waiting for a resource to be unlocked) is **Hierarchical Locking**, that classify the resources in a hierarchical way (so in a DBMS these could be table over tuple over cell, for example). To cope with this additional complexity in the resource model we must specialize the existing locks in order to take into account the hierarchy of resources. We must introduce *intention* locks, that are a modification of a lock primitive that communicates the need of locking a *subelement* in a given mode. The status request table becomes:

Type of request	Resource status				
	ISL	IXL	SL	SIXL	XL
ISL	G	G	G	G	D
IXL	G	G	D	D	D
SL	G	D	G	D	D
SIXL	G	D	D	D	D
XL	D	D	D	D	D

In order to obtain a lock on a certain resource, a transaction must hold an equally or more restrictive lock on its parent. In the case of read lock, the hierarchy is $(ISL, IXL \geq SL, ISL)$ and in the case of write locks it's $(SIXL, IXL \geq IXL, SIXL, XL)$

2.2.5 Timestamping

Timestamping is an alternative approach to locking, usually called optimistic cuncurrency control⁷. The idea behind is really simple and straightforward: we can timestamp (numerate) events in a way to impose a total ordering of them in the system, and then checked at commit time if there are not "time travelling" things.

The implementation makes use of two counters, one for the Last Read Time and one for the Last Write Time. The policy is:

- if we want to read a value that has been written *in the future wrt to us* that read is rejected, otherwise is accepted and the Read Time counter will be updated.
- If we want to write a value that *has been already read or written in the future* the transaction is killed, otherway is accepted and the Write Time counter is updated.

Thomas Rule A modification of the write policy (widely used) that can reduce the kill rate of optimistic cuncurrency control is this:

*on a write operation, we reject it **only if** the value has been read in the future. Otherwise, we just **skip** the write operation, since it is obsolete.*

Multiversioning A simple extension of timestamping uses "verisons" of data to avoid killing transactions; the idea is pretty simple, every write does not delete the previous value, but instead it creates a new copy (with its own Write Time counter). Read operations access only the right value for their needs. *This means that there are no invalid reads.* The only operation that is rejected a priori is a write operation on an object that has already been read in the future.

TimeStamping Multiversioning is a set of schedules that is NOT a proper subset of *any* set we've seen so far. It also comprises non-VSR schedules (much like TimeStamping with Thomas Rule).

⁷while locking is usually referred to as "pessimistic" cuncurrency control

3 Deadlock Detection - Obermarck Algorithm

The Obermarck's algorithm is a distributed algorithm that has the aim of creating a "distributed dependency graph" in order to inform all nodes of possible deadlock situations.

3.1 Applicability

The O's algorithm can be applied only if a certain set of conditions is met:

1. A somewhat centralized control unit exists (that assigns unique labels, at least)
2. Transactions can be decomposed in sub-transactions running on other nodes
3. Super-transactions depend on sub-transactions \Rightarrow "fathers must wait for sons" to proceed and terminate
4. Two kind of wait-for relationships must exists:
 - Father - Son (or brothers) wait for each other *inside a single node* because of locks etc
 - Transactions wait for each other *across nodes* through external network calls

3.2 The Algorithm

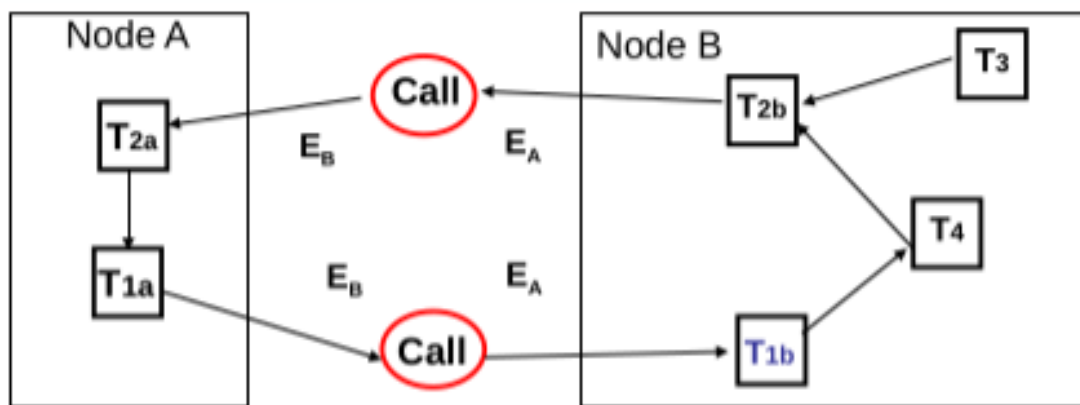
As already said, the idea is that each node exchanges enough information with his neighbours in order to project onto them a sorta global view of all the dependencies.

In practice, Node A sends info to Node B only if these conditions met:

- A has a transaction that *is waited* from a transaction on another node
- A has a transaction that *is waiting* for a transaction on B
- The two linked transactions on A that are also linked to B have "descending indexes" (2 waits from 1)

When a node receives the information about the other transactions from a neighbour, it updates its internal view of the dependency graph and if a cycle is detected, kills a transaction in the cycle.

Example This situation:



Node A sees a descending transaction cycle forming: $E_b \rightarrow T_{2a} \rightarrow T_{1a} \rightarrow E_b = B \rightarrow 2 \rightarrow 1 \rightarrow B$. So it sends it's internal information to B. The same does not happen on B, because the only cycle detected is "in order" ($A \rightarrow 1 \rightarrow 2 \rightarrow A$)

After sending the information to B, A waits. B updates his own dependency graph with all the transactions sent by A, and detects the $1 \rightarrow 2 \rightarrow 1$ circular dependency (that also has T₄ in B,

but it's not relevant in cycle detection; still, it's one of the potential candidate to be killed in order to break the cycle).

Part II

Memory Management

4 Memory Layers

DBMSs run on computers and they have to deal with large datasets. This said, memory management is crucial to deal with DBMS performances. We'll use a reduced model for memory, that only comprises of

- Main memory, structured in *pages*
- Secondary mass storage memory, divided in *blocks*

An I/O operation is defined as "every operation that transfers data from main memory to secondary memory *or viceversa*".

4.1 Blocks, Tuples

How do we map the custom memory structure of a DB schema onto memory? We fix the memory basic block in length (this unit is usually referred as "block" and its read time is usually the basic unit to calculate overall speed of a query) while tuples⁸ are length-varying depending on how data is designed at software level.

This difference in length calls for the definition of a very simple index, the **Block Factor** $B = \lfloor \frac{\text{block size}}{\text{tuple size}} \rfloor$ that simply is the *number of tuples within a block*.

5 Access Structures

Both in main and in secondary memory we have additional data structures (classic data structures as linked lists, trees, hash tables) that are used to speed up the accessing of memory.

Each Table is stored in **exactly one primary data structure** and can have **more than one secondary access structures** but also none.

- The primary structure holds *all the data*. Independently from how the data is accessed, there's no pointer to follow if the access is done through a primary structure.
- The *secondary* structures are used to *index primary structures* and so they contain mostly pointers to tuples. This means that an extra access to the memory is due, if we're accessing data through a secondary memory structure⁹. Secondary data access structures are used to speed up the access / search of data in tables, and are *crucial* when it comes to JOIN operations, for example.

We consider three main categories of data structures: sequential, hashed and tree based structures¹⁰.

5.1 Data Structures Recap

For CS students, data structures should be what we eat at breakfast. I'd like to make a quick recap *considering the fact that we're in the DB realm*, so I'll fly past some aspects while putting the focus on others not in a traditionally way of seeing data structures.¹¹

⁸the actual DB records.

⁹Data can also be stored in secondary structures. We don't care so much though.

¹⁰corresponding to their relative implementations single linked lists, hash maps and black red trees.

¹¹I've killed english in these two lines.

5.1.1 Sequential Structures

We further divide sequential structures in three subcategories:

- Entry sequenced: these are structures organized only by the "arrival order" of tuples
- Arrays: index based access structures¹²
- Sequentially ordered: tuples are ordered (and can be accessed) through a *key*

Let's analyze these structures with regards to the actual operations that must be performed on them

	Entry sequenced	Array	Sequentially ordered
INSERT	Efficient	Efficient	Not Efficient
UPDATE	Efficient	Efficient	Not Efficient for large data sizes
Tuple size	Variable	Fixed	Variable

5.1.2 Hash Based Structures

These structures are *associative structures* that provide access through data by linking it to other piece of data, generally a *well known key* field.

Hash based structures usually make use of "buckets" as a unit of storage (usually the size of a memory block). The buckets of a structure like that are the actual units that are accessed "at once" when an associative access is performed.

Hashing and Stuff Hash based structures make use of an hash function to identify the bucket (so the tuples) they're accessing at a time. Hash functions are prone to *collisions*, so how are that handled? Two main techniques can be deployed:

1. Open addressing: the tuple is inserted (or searched) in *another bucket* (linearly, recursively with the hash function...). This is NOT the method used in DBMSs.
2. Separate chaining: a linked list is of bucket is created, with the first bucket as head. If a tuple has the same hash index as the head BUT it's not in the first bucket, then the list is traversed.

5.1.3 Tree Based Structures

The original implementation of SQL indexes made use of trees. In DBMSs, they're pretty powerful tools because they offer a search possibility, through the key of the tree. The most used version of trees (among Binary, Black Red ecc) in databases is the B+ tree: it's a n-ary tree implementation with *many* children per node and *linked leaves*, so to render the all leaves traversal operation linear in the number of leaves.

The distinction Primary/Secondary structure still holds: if a tree is used as a *primary* data structure the leaves will contain directly the data; otherwise, only the *pointer to the data*. Remember: each leaf has a fixed length and can contain several blocks¹³.

6 Query Optimizer

When a query is issued, a module in the DBMS called *query optimizer* takes care of enhancing the query performance, through relational algebraic transformations and reductions. DB1 course concepts as the relational algebra equivalence are largely used here: the less we access the memory¹⁴ the faster the query will complete.

¹²from an exercise point of view, arrays are not that much used

¹³exercise tips: the "pointers" in a leaf node are just the keys of the tables we're indexing on. So, if the query is selecting that very value, we *don't* have the additional tuple access.

¹⁴that's the only performance enhancement we're looking for

Selectivity Selectivity is a *property of filter predicates*, and indicates the probability that any row will satisfy the predicate specified. This value plays a huge role in query optimization: if we have to apply cascading filters, we may want to apply before the more selective, in order to deal with less tuples afterwards. We always assume, if not specified, that the selectivity of a predicate follows an homogeneous distribution. This means that $S = \frac{1}{N}$ where S is selectivity and N the number of **values**¹⁵ available in that field.

6.1 Costs of Various Access Patterns

6.1.1 Sequential Scan

Well, sequential scan of all blocks. Scan cost = number of blocks (it can eventually be stopped if, for example, we're accessing an ordered field and we have a "less/more than" as a clause).

6.1.2 Lookup

For lookup is intended a query on a field with an equality or interval predicate associated. We'll analyze the cost of lookups on all structures:

Primary Hash If the attribute we're constraining in order to perform our selection¹⁶ is the index of the hash table, we can exploit the associative nature of the structure to speed up the search. We have to distinguish if the lookup is done with an equality predicate or an interval predicate

- Equality lookup: data can be directly accessed from the given constraint, so the overall cost is just 1 + (average overflow chain length). Usually¹⁷ there are no overflow chain, so the cost is just unitary.
- Interval lookup: usually not supported, but the single values in the interval can be accessed one by one: this makes the overall cost of a interval scan on a primary hash equal to the length of the interval itself (with the opportune rescaling to the dimension of the blocks).

Secondary Hash Equality: to the normal 1 + overflow chain length we have to add another access to memory (the access to the actual block in the primary storage).

For intervals, same as Primary plus the access to the memory. Again, interval lookups on associative structures are not the best in term of performance.

Primary B+ Tree All the tree structure must be "traversed downward". This results in an access per level (we are not counting the leaves, for now) Reached the leaves, some scenarios are possible: REMEMBER we're assuming that the tree is built on the field we're using to restrict the search

- Equality lookup: if the *single leaf* is needed (equality constraint on unique field) then the access to the leaf is of weight one.
- Equality lookup BUT multiple selected tuples: then the accesses to the leaf block¹⁸ weights as all the read blocks.
- Interval lookup: exactly as the Equality with multiple tuples to return. The first leaf of the interval is reached and then all the blocks are read sequentially. Total cost: $\{tree\ height\} + \#\{blocks\ accessed\}$.

¹⁵WATCH OUT, it's the number of **values** not the number of **tuples**. This means that a binary field will **always** have a selectivity of 0.5, if assumed homogeneously distributed.

¹⁶the one in the WHERE clause, to be clear

¹⁷in exercises

¹⁸block are grouped in leaves, that are then mapped to the memory

Secondary B+ Tree The cost is similar to the primary B+ tree. The leaves this time are to be considered *independently from the tree itself*, because they just hold the pointer to the memory block. If we're accessing 13 blocks of memory and a leaf only groups 6 of them at a time, in a 4 levels tree, the total cost of the query is

$$\#\{\text{intermediate nodes}\} + \#\{\text{total leaves}\} + \#\{\text{total blocks}\} = 3 + \lceil \frac{13}{6} \rceil + 13 \quad (1)$$

for a total of 20 memory access.

The same reasoning is applicable to the interval queries.

6.1.3 Using Statistics

Calculating the memory accesses needed to complete a query an important role that's not influenced *by the query itself* but rather *by the data* is the actual number of tuples selected by the interrogation. We've indirectly already used them in the last example for the secondary B+ tree: the fact that we know how many blocks are actually to be retrieved (13) must come from an analysis on the contents of the table.

Selectivity is one of the key indexes when optimizing queries: if an interrogation has multiple predicates, the query optimizer will apply first the *most selective ones* in order to reduce the total number of tuples to be taken.

6.1.4 Join

Joins are the most frequent and expensive operations in a database. We can exploit additional data structure to speed them up.

We'll see three main strategies: nested loops, merge scan and hashed join.

Nested Loops Classic bruteforce approach to joining. For each tuple of the first table, all the tuples of the second table must be scanned in order to find the ones to join. In the end, the cost is just the product of the number of blocks of each table. **WATCH OUT:** if it's possible that one of the two tables can be *entirely cached*, then that accesses can be done only one time: the cost reduces to the **sum** of the blocks of the two tables.

A possible improvement of this simple technique is to combine it with lookups: if we have an additional predicate to the query, we can perform a lookup using the predicate to filter out some of the tuples, and then just join the tuples that satisfy the predicate. I'll add an example from the slide down here, keep in mind most of the data is left out. It's the reasoning the important part.

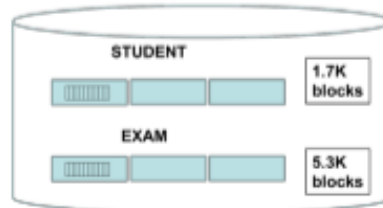
STUDENT (Id, Name, LastName, Email, **City**, Birthdate, Sex)
 EXAM (SID, CourseID, Date, **Grade**)

SELECT STUDENT.*
FROM STUDENT JOIN EXAM ON Id=SID
WHERE City='Milan' AND Grade='30'

val(City)=150
 val(Grade)=17

Option 1: Scan Student, evaluate City='Milan' + Lookup on Exam

- Read all the STUDENT blocks: **1.7K**
- Filter students from Milan → 150K tuples/150 different cities = **1k**
- For 1k students compute the join with the tuples of all the 5.3K blocks in EXAM



Total cost:
 1,7K + 1K*5.3K
 = ~ **7M** I/O accesses

Merge Scan This technique can be exploited if the two merging table are ordered on the index we're using to merge. It's intuitive how the sequential nature of the tuples can be used to *progressively merge* all data. In this case, the cost of the join is linear in the number of blocks to be retrieved¹⁹. If the merge scan approach gives huge performance improvements, one can also sort the tables and then merge-scan join them. This results in an additional cost (the reorder cost) usually in $N \times \log(N)$ where N number of blocks to be accessed.

Hashed Join If two tables are hashed according to the same key of the join operation, then it can be speedupped to cost as a merge scan (so linear in the number of blocks to be read).

7 JPA

7.1 Objects and Relations

Mapping a relational model typical of a DBMS to an object one, typical of object oriented programming, is no easy task. Moreover, it's *inevitable*: 99% of web apps rely on a SQL-like database, and they need to use it in a OOP environment. What is the great and very original idea of JPA? Add another software layer that takes care of the ORM²⁰ and simplifies the passage from one paradigm to another, allowing app programmers to interact with the DB without writing actual DB code.

7.2 JPA Main Features and Concepts

How do people use JPA²¹ and how is it integrated in a plain Java application? JPA adds these concepts to a standard Java application:

- Entity: a *class* that represents a table in the DB. You will get confused: with "entity", we'll refer both to the class and the instance
- Persistence Unit: the *set of classes* (generally a module) that are mapped to the same DB SCHEMA

¹⁹Remember: a B+ is a sequential structure.

²⁰Object Relational Mapping

²¹they don't, by the way

- Persistence Context: all the *actual instanciated objects* of a persistence unit (analogue of the DB INSTANCE)
- Managed Entity: an object of an Entity class, in the persistence context. It will reflect all changes from and to the DB
- Entity Manager: the interface to the the Persistence Context
- Client: a component that can interact with the Persistence Context

all these "new concepts" are added to the good ol Java language through annotations: if a Java class imports `javax.persistence`²² it has access to a set of annotations that in the end extends the language to accomodate the mapping to a relational model. The main concepts to keep in mind, among the one presented, are the Entity, the Entity Manager and the concept of Managed Entity.

7.2.1 Entity, Managers, Managed Entities

When using the "entity" term, we must watch out if it means the entity *class* or the instance in the application. We'll refer to the class as "Entity Class" and to the instance as "Entity" from now on.

An entity is an object²³ that gets associated with a *tuple* in the database. If the entity is also *managed* (being the entity manager aware of his presence and linked it to the corresponding tuple) then every changes made to the tuple are reflected to the entity and viceversa. Remember: an entity **CAN BE UNMANAGED**. An entity class has features from both the object and the relational worlds: it must have a field marked as primary key, it must have fields that are annotated to match the table columns, but it can have methods (as the usual getters, setters, serializers, hashers, iterators-returning ecc ecc). For the correct and complete syntax, use the manual. I'm reporting here just the snippets of²⁴ Java code that is useful to understand how entities are defined and behaves according to relationships.

```
@Entity
@Table(name = "EMP")
public class Employee implements Serializable{

    // the primary key is field-accessed, the identifier is on the field itself
    @Id
    @GeneratedValue(strategy = GenerationType.IDENTITY) // this specifies how to handle new items
    private long EmployeeId

    private int Salary;

    // this field is property-accessed instead, the annotation is on the getter
    @Column(name = "SALARY", nullable = false)
    public int getSalary(){
        return this.Salary
    }

    // many employees in the same department: owning side of a 1:N relationship
    @ManyToOne
    @JoinColumn(name = "DEPARTMENT")
    private Department department;

    // one employee, some parking lots
    // careful here: the mappedBy field tells which FIELD maps these entity, it has nothing
    // to do with the DB
    // the mappedBy should also be present in OneToOne relationships
```

²²maybe you'll have to replace Java with Jakarta, I'll let you discover that yourself

²³classic Java object

²⁴commented

```

@OneToMany(mappedBy = "employee")
private Collection<ParkingLot> parkingLots;

// many employees, many projects
// M:N relationships want a bridge table (as in ER schemas, they are the composition of two
//      1:N relationships)
// this also means that the owner of the relationship is decided by the schema designer
@ManyToMany
@JoinTable(
    name = "EMP_PROJ",
    joinColumns = @JoinColumn(name = "EMP_ID"),
    inverseJoinColumns = @JoinColumn(name = "PROJ_ID")
)
private Collection<Project> projects;
}

```

The Entity Manager The Entity Manager is the interface to ALL the managed objects, and reunites all the functionality to interact with the database. Its interface is super simple:

<code>public void persist(Object entity)</code>	makes an entity managed. Remember: <i>managed</i> \neq <i>written in the DB</i> .
<code>public <T> find(Class <T> entityClass, Object primaryKey)</code>	returns an object from the Persistence Context, searching his primary key. The object will be managed from now on. This triggers also the fetching of the data specified in the fetch policy of the entity class.
<code>public void remove(Object entity)</code>	removes an entity from the persistence context \Rightarrow from the DB. The plain object in the Java application is still present (pay attention to this)
<code>public void refresh(Object entity)</code>	re-sync the entity with the DB tuple
<code>public void flush()</code>	writes all changes to DB

This interface is easy to use and minimal enough to manage the complex mechanism that is a persistence context. Remember: the **new** keyword do not add the entity to the persistence context.

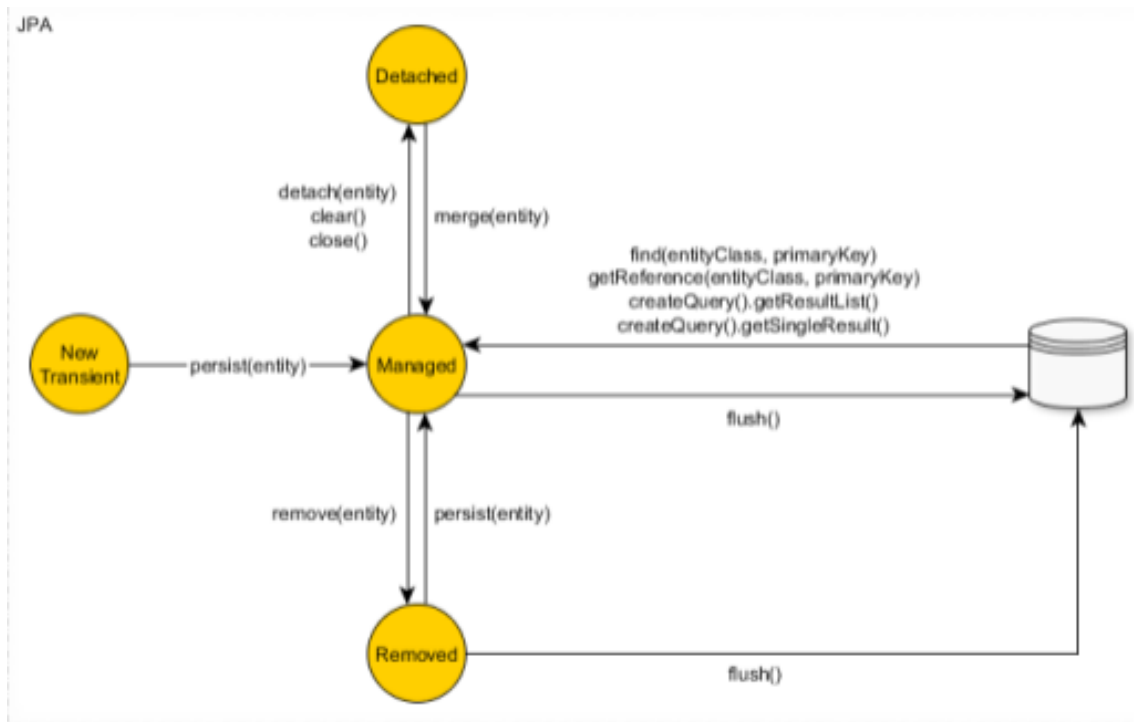


Figure 2: State transitions and associated calls in the JPA framework.

Fraternali spends 53 slides jammed of font 10 JPA code to explain how the various type of entity managers behaves wrt the business Java objects (the EJB, Enterprise Java Beans) and wrt the persistence context and the transactions; it's all summed up here:

EM	Type of	Container Managed Transaction Scoped	Container Managed Extended	Application Managed
Type of client		@stateless EJB	@stateful EJB	Any
Who creates the transaction?		Container	Container	Client via EntityManager or UserTransaction
Who starts/ends the transaction		Container starts a transaction when no transaction is active and EJB method is called Transaction ends when method terminates	Container starts a transaction when no transaction is active and EJB method is called Transaction ends when method terminates	Client explicitly demarcates with begin/commit
Is transaction propagated?		YES Container propagates the transaction across EM references on the same persistence unit and on the call stack	YES Container propagates the transaction across EM references on the same persistence unit and on the call stack	NO, but a method can join the active transaction

Again, refer to the manual to find all the right documentation for this stuff.

7.2.2 Relationships

Mapping relationships between an object oriented worlds and an entity relationship one is kinds strange. Controintuitively, and OO world has *richer* relations between things wrt ER.

An OO relationship is characterized by

- A **direction**: each entity class, in order to access the *other side of the relationship*, may have one or more fields that references the entities on the other side. Practically, the "Author" entity class could reference a "Book" entity class: this is implemented by adding a collection²⁵ of "Books" directly inside the "Author" entity class.
- A **role**: at every end of the relation, entities are playing a role. A role represents if you're the source or the target of some data.
- A **cardinality**: similar to the cardinality in ER, a relation can be single/multiple on each side. In JPA, a many-to-many relationship model is also available: this will be mapped with an intermediate table on the DB.
- A **ownership**: one of the two entity classes at the end of a relationship is said to be the *owner* of it. The owner is the entity class that *in the db* has the foreign key.

8 Active Databases - Triggers

Triggers are event based stored procedures in a DBMS. They are SQL statements (usually enriched with some more programming-language-feeling functionalities) associated to an update/insert/delete event on a specific table (or column of a table) and can be used to validate inputs, check and impose invariants, generate and update metadata and so on. They're very powerful.

Triggers follows the ECA paradigm for their definition:

- **E**vent fires
- **C**ondition is met
- then **A**ction is executed

watch out for the composite (event + condition) combined.

A full trigger expression is composed of:

```
create trigger <TriggerName>
{before | after}
{insert | delete | update [of <Coloumn>]} on <Table>
[referencing
  {
    [old table as <OldTAlias>]
    [new table as <NewTAlias>]
    [old row as <OldRAlias>]
    [new row as <NewRAlias>]
  }
]
[for each {row | statement}]
[when <Condition>]
<Sql Procedural Statement>
```

where angular brackets indicate literals, square brackets optionality²⁶ and curly brackets and vertical bars interchangeability.

Before and After The "before" and "after" modifier determine when the trigger would be executed with regard to the update being observed.

Before an event means *before the update takes place*, that also mean *before a status change in the schema*. Quick reminder: before triggers cannot alter the table. They can only adjust the modified rows.

Also, they cannot exactly modify entirely the updating rows: a before trigger can operate in the space of the "new and old" values and tables (see "New and Old" down here).

After triggers are executed after the state of the table has changed.

²⁵usually an implementation of List

²⁶all the aliases are optional, alongside the specific column for the update table

Row and Statement Pretty straightforward distinction: a statement *could* affect multiple rows. **for each row** triggers are fired at each update on single rows, while **for each statement** triggers are fired only when the event is caught, even if it affects multiple rows.

Old and New When dealing with updates, old values and new values must be taken into consideration. SQL generates two variables (old / new for rows, old table / new table for statements) to store the "before and after" values. These variables and accessible values can be very useful when dealing with **when** statements.

Order of Triggering If several triggers are associated to the same event, they're executed following this order:

1. BEFORE STATEMENT triggers
2. BEFORE ROW triggers
3. state of the table is changed
4. AFTER ROW triggers
5. AFTER STATEMENT triggers