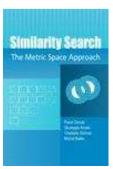
# SIMILARITY SEARCH The Metric Space Approach

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### Table of Content

Part I: Metric searching in a nutshell

- Foundations of metric space searching
- Survey of existing approaches

### Part II: Metric searching in large collections

- Centralized index structures
- Approximate similarity search
- Parallel and distributed indexes

# Features of "good" index structures

### Dynamicity

support insertions and deletions and minimize their costs

#### Disk storage

for dealing with large collections of data

#### CPU & I/O optimization

 support different distance measures with completely different CPU requirements, e.g., L<sub>2</sub> and quadratic-form distance.

### Extensibility

 similarity queries, i.e., range query, k-nearest neighbors query Centralized Index Structures for Large Databases

- 1. M-tree family
- 2. hash-based metric indexing
- 3. performance trials

# M-tree Family

### The M-tree

- Bulk-Loading Algorithm
- Multi-Way Insertion Algorithm
- The Slim Tree
- Slim-Down Algorithm
  - Generalized Slim-Down Algorithm
- Pivoting M-tree
- The M<sup>+</sup>-tree

### The M<sup>2</sup>-tree

### The M-tree

- Inherently dynamic structure
- Disk-oriented (fixed-size nodes)
- Built in a bottom-up fashion
  - Inspired by R-trees and B-trees
- All data in *leaf nodes*
- Internal nodes: pointers to subtrees and additional information
- Similar to GNAT, but objects are stored in leaves.

### M-tree: Internal Node

Internal node consists of an entry for each subtree

- Each entry consists of:
  - Pivot: p
  - Covering radius of the sub-tree: r<sup>c</sup>
  - □ Distance from *p* to *parent* pivot  $p^p$ :  $d(p,p^p)$
  - Pointer to sub-tree: ptr

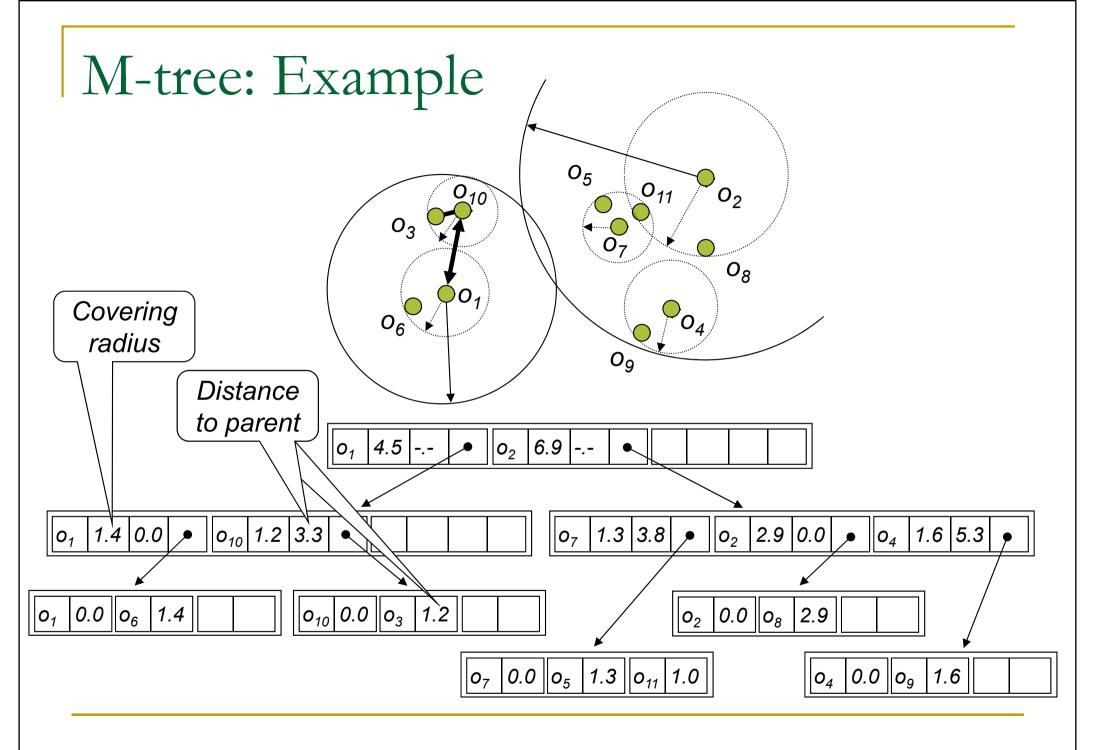
 $|\langle p_1, r_1^c, d(p_1, p^p), ptr_1 \rangle| |\langle p_2, r_2^c, d(p_2, p^p), ptr_2 \rangle| \cdots |\langle p_m, r_m^c, d(p_m, p^p), ptr_m \rangle|$ 

• All objects in subtree *ptr* are within the distance  $r^c$  from *p*.

### M-tree: Leaf Node

- leaf node contains data entries
- each entry consists of pairs:
  - object (its identifier): o
  - distance between o and its parent pivot:  $d(o, o^p)$

$$\langle o_1, d(o_1, o^p) \rangle \langle o_2, d(o_2, o^p) \rangle \cdots \langle o_m, d(o_m, o^p) \rangle$$



### M-tree: Insert

- Insert a new object o<sub>N</sub>:
- recursively descend the tree to locate the most suitable leaf for o<sub>N</sub>
- in each step enter the subtree with pivot p for which:
  - □ no enlargement of radius  $r^c$  needed, i.e.,  $d(o_N, p) \le r^c$ 
    - in case of ties, choose one with the nearest p to  $o_N$
  - □ minimize the enlargement of *r*<sup>c</sup>

### M-tree: Insert (cont.)

when reaching leaf node N then:

• if *N* is not full then store  $o_N$  in *N* 

• else **Split**( $N, o_N$ ).

# M-tree: Split

**Split**(*N*,*o*<sub>*N*</sub>):

- Let S be the set containing all entries of N and  $o_N$
- Select pivots p<sub>1</sub> and p<sub>2</sub> from S
- Partition S to  $S_1$  and  $S_2$  according to  $p_1$  and  $p_2$
- Store S<sub>1</sub> in N and S<sub>2</sub> in a new allocated node N'
- If N is root
  - Allocate a new root and store entries for  $p_1$ ,  $p_2$  there
- else (let N<sup>p</sup> and p<sup>p</sup> be the parent node and parent pivot of N)
  - Replace entry  $p^p$  with  $p_1$
  - If  $N^p$  is full, then **Split**( $N^p$ , $p_2$ )
  - else store  $p_2$  in node  $N^p$

### M-tree: Pivot Selection

Several pivots selection policies

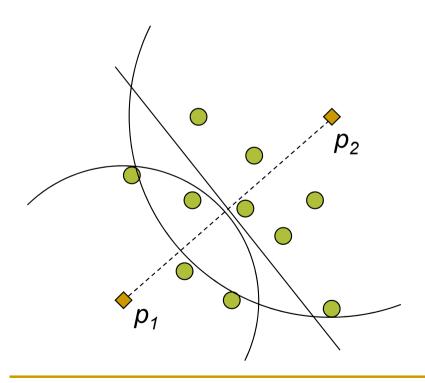
- **RANDOM** select pivots  $p_1$ ,  $p_2$  randomly
- □  $m\_RAD$  select  $p_1$ ,  $p_2$  with minimum  $(r_1^c + r_2^c)$
- □  $mM\_RAD$  select  $p_1$ ,  $p_2$  with minimum  $max(r_1^c, r_2^c)$
- □ *M*\_*LB*\_*DIST* − let  $p_1 = p^p$  and  $p_2 = o_i | \max_i \{ d(o_i, p^p) \}$

Uses the pre-computed distances only

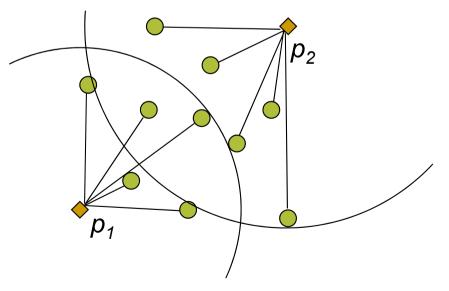
- Two versions (for most of the policies):
  - □ **Confirmed** reuse the original pivot *p<sup>p</sup>* and select only one
  - Unconfirmed select two pivots (notation: RANDOM\_2)
- In the following, the *mM\_RAD\_2* policy is used.

# M-tree: Split Policy

- Partition S to  $S_1$  and  $S_2$  according to  $p_1$  and  $p_2$
- Unbalanced
  - Generalized hyperplane



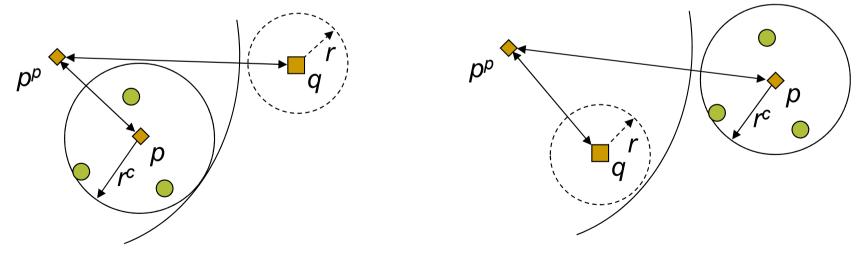
- Balanced
  - Larger covering radii
  - Worse than unbalanced one



# M-tree: Range Search

### Given R(q,r):

- Traverse the tree in a depth-first manner
- In an internal node, for each entry (p,r<sup>c</sup>,d(p,p<sup>p</sup>),ptr)
  - □ Prune the subtree if  $|d(q,p^p) d(p,p^p)| r^c > r$
  - Application of the pivot-pivot constraint

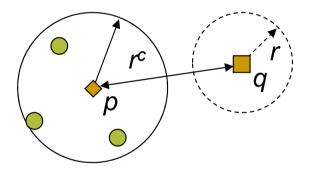


### M-tree: Range Search (cont.)

If not discarded, compute d(q,p) and

• Prune the subtree if  $d(q,p) - r^c > r$ 

Application of the range-pivot constraint



All non-pruned entries are searched recursively.

# M-tree: Range Search in Leaf Nodes

- In a leaf node, for each entry  $\langle o, d(o, o^p) \rangle$ 
  - □ Ignore entry if  $|d(q,o^p) d(o,o^p)| > r$
  - □ else compute d(q,o) and check  $d(q,o) \le r$
  - Application of the object-pivot constraint

### M-tree: k-NN Search

### Given *k*-*NN*(*q*):

- Based on a *priority queue* and the pruning mechanisms applied in the range search.
- Priority queue:
  - Stores pointers to sub-trees where qualifying objects can be found.
  - Considering an entry  $E = \langle p, r^c, d(p, p^p), ptr \rangle$ , the pair  $\langle ptr, d_{min}(E) \rangle$  is stored.
  - $\Box \quad d_{min}(E) = max \{ d(p,q) r^{c}, 0 \}$
- Range pruning: instead of fixed radius r, use the distance to the k-th current nearest neighbor.

# M-tree Family

The M-tree

### Bulk-Loading Algorithm

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### The M<sup>2</sup>-tree

# Bulk-Loading Algorithm

- first extension of M-tree
- improved tree-building (insert) algorithm
- requires the dataset to be given in advance

#### Notation:

- Dataset  $X = \{o_1, \dots, o_n\}$
- □ Number of entries per node: *m*

### Bulk-Loading Algorithm:

- □ First phase: build the M-tree
- Second phase: refinement of unbalanced tree

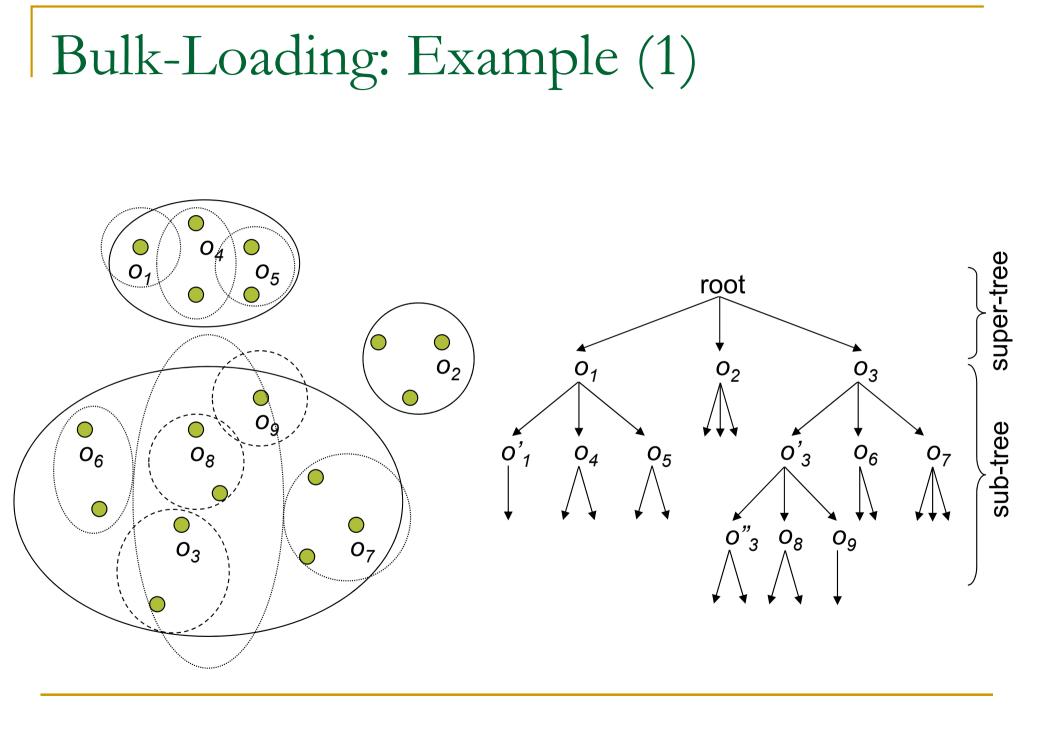
# Bulk-Loading: First Phase

randomly select / pivots P={p<sub>1</sub>,...,p<sub>l</sub>} from X
 Usually *I=m*

- objects from X are assic
- objects from X are assigned to the *nearest pivot* producing *I* subsets *P*<sub>1</sub>,...,*P*<sub>1</sub>
- recursively apply the bulk-loading algorithm to the subsets and obtain / sub-trees T<sub>1</sub>,...,T<sub>1</sub>

leaf nodes with maximally / objects

create the root node and connect all the sub-trees to it.



# Bulk-Loading: Discussion

Problem of choosing pivots  $P=\{p_1,...,p_l\}$ 

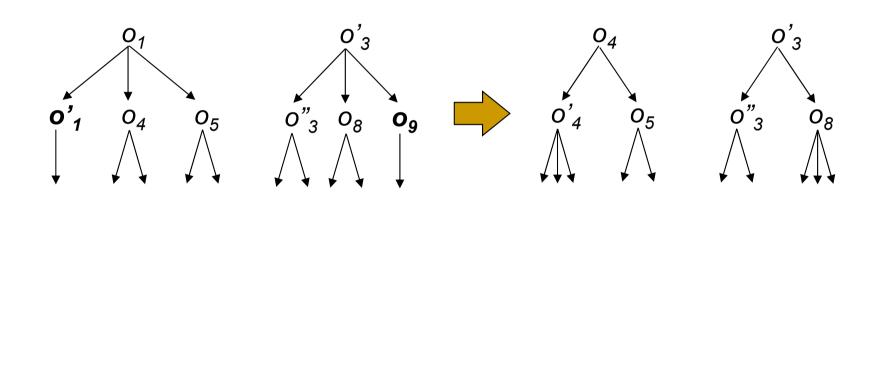
- sparse region → shallow sub-tree
  - far objects assigned to other pivots
- dense region  $\rightarrow$  deep sub-tree
- observe this phenomenon in the example

# Bulk-Loading: Second Phase

- refinement of the unbalanced M-tree
- apply the following two techniques to adjust the set of pivots P={p<sub>1</sub>,...,p<sub>l</sub>}
  - under-filled nodes reassign to other pivots and corresponding pivots deleted from P
  - deeper subtrees *split* into shallower ones and *add* the obtained pivots to *P*

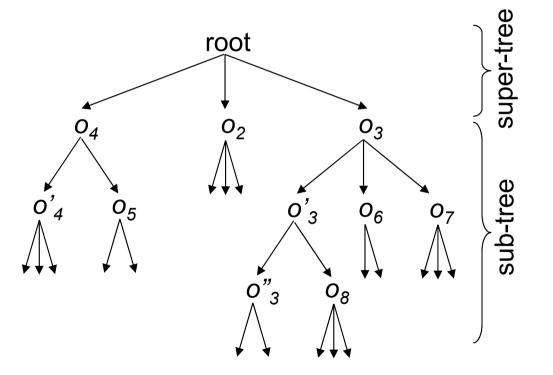
# Bulk-Loading: Example (2)

• Under-filled nodes in the example:  $o'_1, o_9$ 



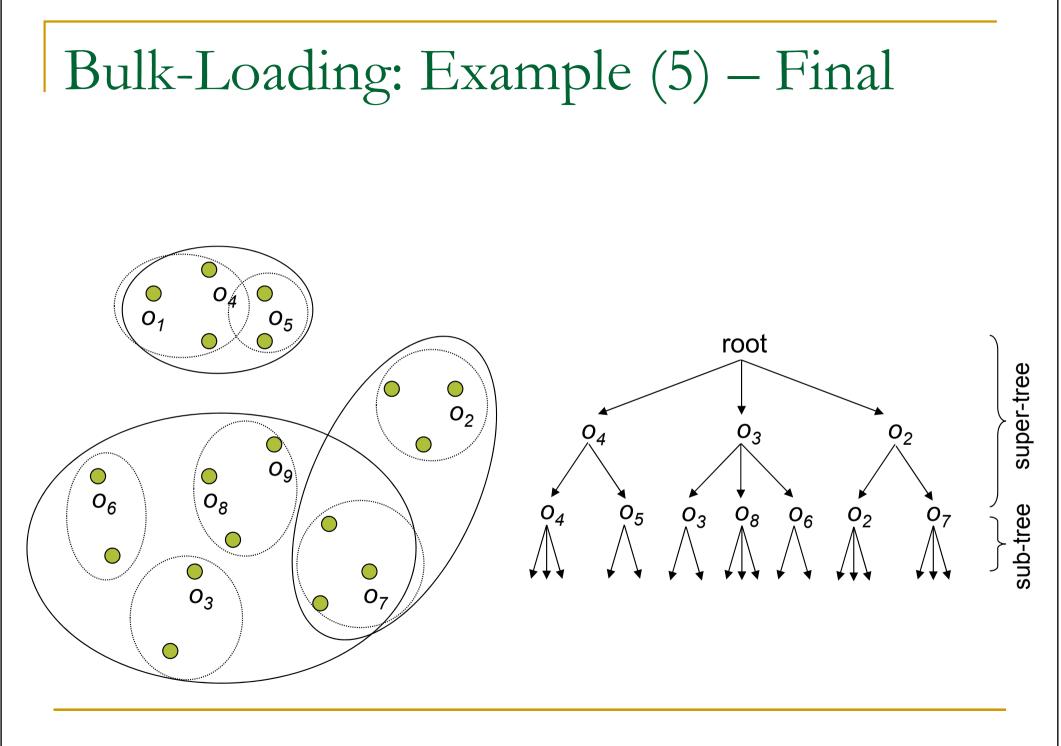
# Bulk-Loading: Example (3)

#### After elimination of under-filled nodes.



# Bulk-Loading: Example (4)

- Sub-trees rooted in o<sub>4</sub> and o<sub>3</sub> in the tree are deeper
- split them into new subtrees rooted in  $O'_4$ ,  $O_5$ ,  $O''_3$ ,  $O_8$ ,  $O_6$ ,  $O_7$
- add them into *P* and remove  $o_4, o_3$
- build the super-tree (two levels) over the final set of pivots P={o<sub>2</sub>, o'<sub>4</sub>, o<sub>5</sub>, o"<sub>3</sub>, o<sub>8</sub>, o<sub>6</sub>, o<sub>7</sub>} – from Sample (3)



# Bulk-Loading: Optimization

- Reduce the number of distance computations in the recursive calling of the algorithm
  - after initial phase, we have distances  $d(p_j, o_i)$  for all objects  $X = \{o_1, \dots, o_n\}$  and all pivots  $P = \{p_1, \dots, p_l\}$
  - Assume the recursive processing of  $P_1$
  - New set of pivots is picked  $\{p_{1,1}, ..., p_{1,l'}\}$
  - □ During clustering, we are assigning every object  $o \in P_1$  to its nearest pivot.
  - The distance  $d(p_{1,i}, o)$  can be lower-bounded:

 $|d(p_1,o) - d(p_1,p_{1,j})| \le d(p_{1,j},o)$ 

# Bulk-Loading: Optimization (cont.)

If this lower-bound is greater than the distance to the closest pivot  $p_{1,N}$  so far, i.e.,  $|d(p_1,o) - d(p_1,p_{1,j})| > d(p_{1,N},o)$ 

then the evaluation of  $d(p_{1,i}, o)$  can be avoided.

- Cuts costs by 11%
  - □ It uses pre-computed distances to a single pivot.
  - by 20% when pre-computed distances to multiple pivots are used.

# M-tree Family

- The M-tree
- Bulk-Loading Algorithm
- Multi-Way Insertion Algorithm
- The Slim Tree
- Slim-Down Algorithm
  - Generalized Slim-Down Algorithm
- Pivoting M-tree
- The M<sup>+</sup>-tree
- The M<sup>2</sup>-tree

# Multi-Way Insertion Algorithm

- another extension of M-tree insertion algorithm
- objective: *build more compact trees* 
  - reduce search costs (both I/O and CPU)
- for dynamic datasets (not necessarily given in advance)
- increase insertion costs slightly
- the original single-way insertion visits exactly one root-leaf branch
  - □ leaf with *no* or *minimum* increase of covering radius
  - not necessarily the most convenient

### Multi-Way Insertion: Principle

- when inserting an object  $o_N$
- run the *point query*  $R(o_N, 0)$
- for all visited leaves (they can store o<sub>N</sub> without radii enlargement): compute the distance between o<sub>N</sub> and the leaf's pivot
- choose the closest pivot (leaf)
- if no leaf visited run the single-way insertion

# Multi-Way Insertion: Analysis

### Insertion costs:

- 25% higher I/O costs (more nodes examined)
- higher CPU costs (more distances computed)

#### Search costs:

- 15% fewer disk accesses
- almost the same CPU costs for the range query
- 10% fewer distance computations for k-NN query

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### The Slim Tree

- extension of M-tree the same structure
  - speed up insertion and node splitting
  - improve storage utilization
- new node-selection heuristic for insertion
- new node-splitting algorithm
- special post-processing procedure
  - make the resulting trees more compact.

#### Slim Tree: Insertion

Starting at the root node, in each step:

- find a node that covers the incoming object
- if none, select the node whose pivot is *the nearest* 
  - M-tree would select the node whose covering radius requires the smallest expansion
- if several nodes qualify, select the one which occupies the minimum space
  - M-trees would choose the node with closest pivot

#### Slim Tree: Insertion Analysis

- fill insufficiently occupied nodes first
  - defer splitting, boost node utilization, and cut the tree size
- experimental results (the same *mM\_RAD\_2* splitting policy) show:
  - Iower I/O costs
  - nearly the same number of distance computations
  - this holds true for both the tree building procedure and the query execution

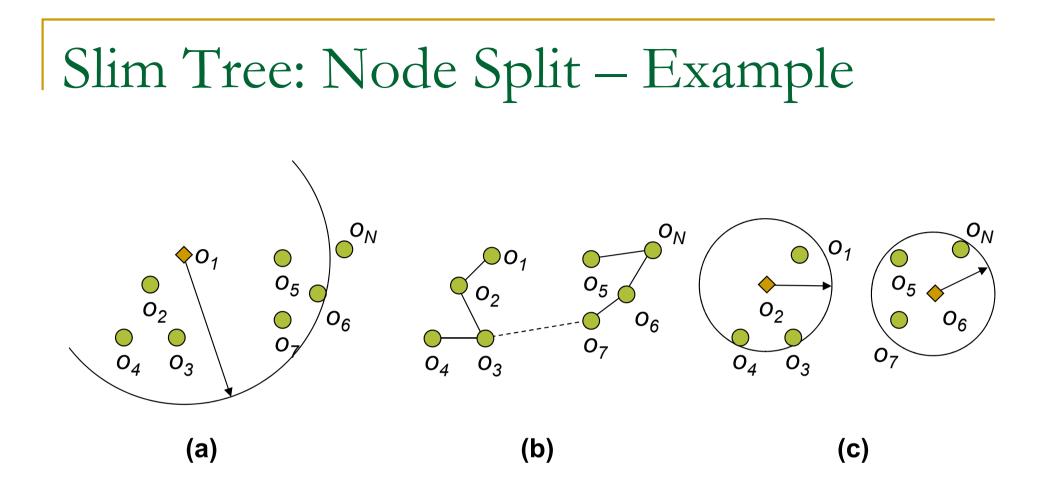
## Slim Tree: Node Split

- splitting of the overfilled nodes high costs
- mM\_RAD\_2 strategy is considered the best so far
  - Complexity  $O(n^3)$  using  $O(n^2)$  distance computations
- the Slim Tree splitting based on the *minimum* spanning tree (MST)
  - Complexity  $O(n^2 \log n)$  using  $O(n^2)$  distance computations
- the MST algorithm assumes a full graph
  - □ *n* objects
  - □ n(n-1) edges distances between objects

# Slim Tree: Node Split (cont.)

#### Splitting policy based on the MST:

- 1. build the *minimum* spanning tree on the full graph
- 2. delete the *longest edge*
- 3. the two resulting sub-graphs form the *new nodes*
- 4. choose the *pivot* for each node as the *object* whose distance to the *others* in the group is *the shortest*



- (a) the original Slim Tree node
- (b) the minimum spanning tree
- (c) the new two nodes

# Slim Tree: Node Split – Discussion

- does not guarantee the balanced split
- a possible variant (more balanced splits):
  - choose the most appropriate edge from among *the longer* edges in the MST
  - if no such edge is found (e.g., for a star-shaped dataset), accept the original unbalanced split
- experiments prove that:
  - tree building using the MST algorithm is at least forty times faster than the *mM\_RAD\_2* policy
  - query execution time is not significantly better

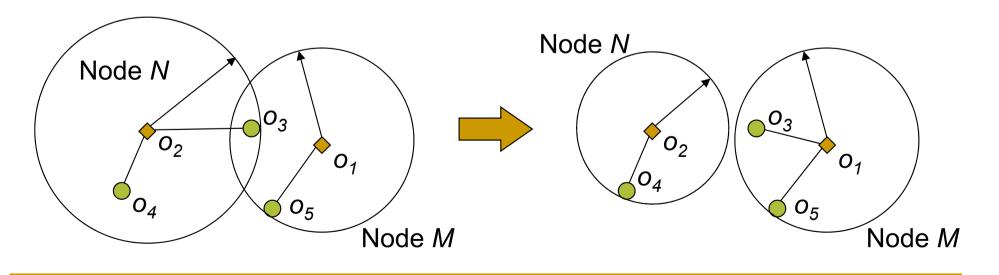
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### Slim-Down Algorithm

- post-processing procedure
- reduce the fat-factor of the tree
  - basic idea: reduce the overlap between nodes on one level
  - minimize number of nodes visited by a point query, e.g.,  $R(o_3, 0)$



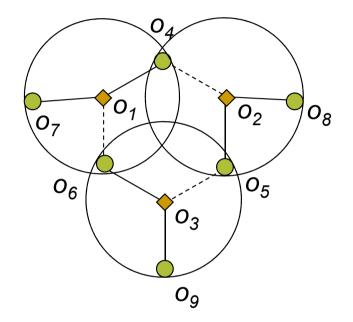
# Slim-Down Algorithm: The Principle

For each node *N* at the leaf level:

- 1. Find object *o* furthest from pivot of *N*
- 2. Search for a sibling node *M* that also covers *o*. If such a not fully occupied node exists, move *o* from *N* to *M* and update the covering radius of *N*.
- Steps 1 and 2 are applied to all nodes at the given level. If an object is relocated after a complete loop, the entire algorithm is executed again.
- Observe moving of  $o_3$  from N to M on previous slide.

### Slim-Down Algorithm: Discussion

- Prevent from infinite loop
  - cyclic moving of objects  $o_4, o_5, o_6$
- Limit the number of algorithm cycles



- Trials proved reducing of I/O costs of at least 10%
- The idea of dynamic object relocation can be also applied to defer splitting.
  - Move distant objects from a node instead of splitting it.

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## Generalized Slim-Down Algorithm

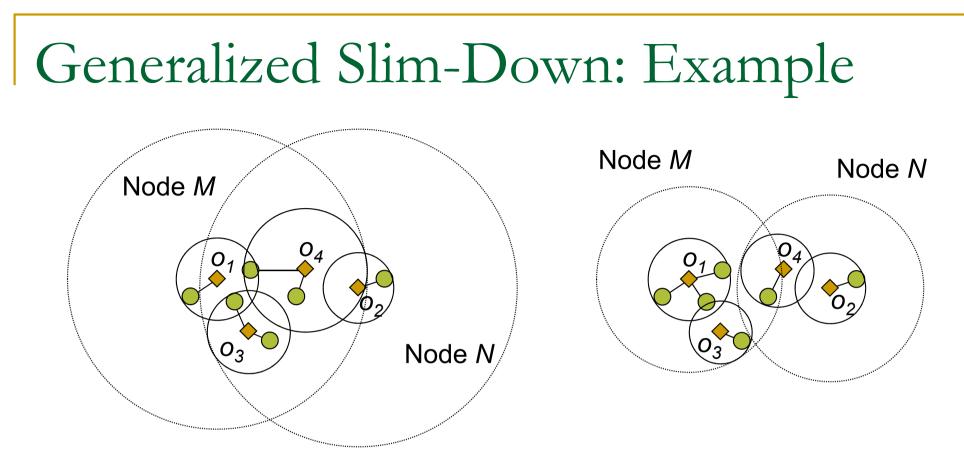
- generalization of Slim-down algorithm for non-leaf tree levels
- the covering radii r<sup>c</sup> must be taken into account before moving a non-leaf entry
- the generalized Slim-down starts from the leaf level
  - follow the original Slim-down algorithm for leaves
- ascend up the tree terminating in the root

## Generalized Slim-Down: The Principle

For each entry  $E = \langle p, r^c, ... \rangle$  at given non-leaf level:

pose range query R(p,r<sup>c</sup>),

- the query determines the set of nodes that *entirely* contain the query region,
- from this set, choose the node M whose parent pivot is closer to p than to p<sup>p</sup>,
- if such *M* exists, move the entry *E* from *N* to *M*,
- if possible, shrink the covering radius of *N*.



Leaf level:

• move two objects from  $o_3$  and  $o_4$  to  $o_1$  – shrink  $o_3$  and  $o_4$ 

#### Upper level:

- originally node M contains  $o_1, o_4$  and node N contains  $o_2, o_3$
- swap the nodes of  $o_3$  and  $o_4$

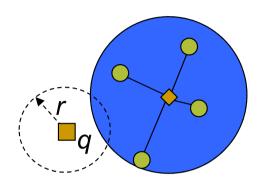
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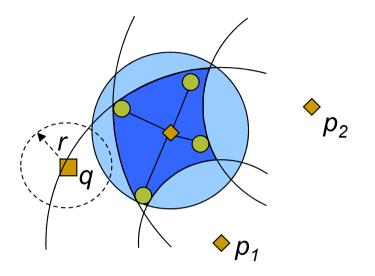
## Pivoting M-tree

- upgrade of the standard M-tree
- bound the region covered by nodes more tightly
  - define additional *ring regions* that restrict the ball regions
  - ring regions: pivot *p* and two radii  $r_{min}$ ,  $r_{max}$
  - □ such objects *o* that:  $r_{min} \le d(o,p) \le r_{max}$
- basic idea:
  - Select additional *pivots*
  - Every pivot defines two boundary values between which all node's objects lie.
  - Boundary values for each pivot are stored in every node.
     (see a motivation example on the next slide)

# PM-tree: Motivation Example



- original M-tree
- range query R(q,r) intersects the node region



- PM-tree (two pivots)
- this node not visited for query R(q,r)

#### PM-tree: Structure

- select additional set of pivots |P|=n<sub>p</sub>
- leaf node entry: (o,d(o,o<sup>p</sup>),PD)
  - □ PD array of  $n_{pd}$  pivot distances:  $PD[i]=d(p_i, o)$
  - Parameter  $n_{pd} < n_p$
- internal node entry: (p,r<sup>c</sup>,d(p,p<sup>p</sup>),ptr,HR)
  - *HR* array of  $n_{hr}$  intervals defining ring regions  $HR[j].min = min(\{d(o, p_j) | \forall o \in ptr\})$  $HR[j].max = max(\{d(o, p_j) | \forall o \in ptr\})$

□ parameter  $n_{hr} < n_p$ 

#### PM-tree: Insertion

- insertion of object  $o_N$
- the *HR* arrays of nodes visited during insertion must be updated by values  $d(o_N, p_i)$  for all  $i \le n_{hr}$
- the leaf node:
  - □ create array *PD* and fill it with values  $d(o_N, p_j)$ ,  $\forall j \le n_{pd}$
- values d(o<sub>N</sub>, p<sub>j</sub>) are computed only once and used several times – max(n<sub>hr</sub>, n<sub>pd</sub>) distance computations
- insertions may force node splits

# PM-tree: Node Split

- node splits require some maintenance
- leaf split:
  - □ set arrays *HR* of two new internal entries
  - set HR[i].min and HR[i].max as min/max of PD[j]
  - □ compute additional distances:  $d(p_j, o)$ ,  $\forall j (n_{pd} < j \le n_{hr})$ and take them into account
  - can be expensive if  $n_{hr} >> n_{pd}$
- internal node split:
  - creating two internal node entries with HR
  - set these HR arrays as union over all HR arrays of respective entries

# PM-tree: Range Query

#### Given R(q,r):

 $n_{hr}$ 

- evaluate distances  $d(q,p_i)$ ,  $\forall i (i \leq max(n_{hr},n_{pd}))$
- traverse the tree, internal node (p,r<sup>c</sup>,d(p,p<sup>p</sup>),ptr,HR) is visited if both the expressions hold:

$$d(q,p) \le r + r^{c}$$

$$\bigwedge_{i=1}^{n} (d(q, p_i) - r \le HR[i].\max \land d(q, p_i) + r \ge HR[i].\min)$$

leaf node entry test:

$$\bigwedge_{i=1}^{n_{pd}} (|d(q, p_i) - PD[i]| \le r)$$

M-tree: the first condition only

### PM-tree: Parameter Setting

#### general statements:

- existence of PD arrays in leaves reduce number of distance computations but increase the I/O cost
- □ the *HR* arrays reduce both CPU and I/O costs

#### experiments proof that:

- □  $n_{pd}$ =0 decreases I/O costs by 15% to 35% comparing to Mtree (for various values of  $n_{hr}$ )
- CPU cost reduced by about 30%
- □  $n_{pd} = n_{hr} / 4$  leads to the same I/O costs as for M-tree
- with this setting up to 10 times faster
- particular parameter setting depends on application

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#### The M<sup>+</sup>-tree

- modification of the M-tree
- restrict the application to  $L_p$  metrics (vector spaces)
- based on the concept of key dimension
- each node partitioned into two twin-nodes
  - partition according to a selected key dimension

## M<sup>+</sup>-tree: Principles

- in an *n*-dimensional vector space
- key dimension for a set of objects is the dimension along which the data objects are most spread
- for any dimension  $D_{key}$  and vectors  $(x_1, \dots, x_n), (y_1, \dots, y_n)$

$$|x_{D_{key}} - y_{D_{key}}| \le \sqrt{(x_1 - y_1)^2 + \dots + (x_n - y_n)^2}$$

- this holds also for other  $L_p$  metrics
- this fact is applied to prune the search space

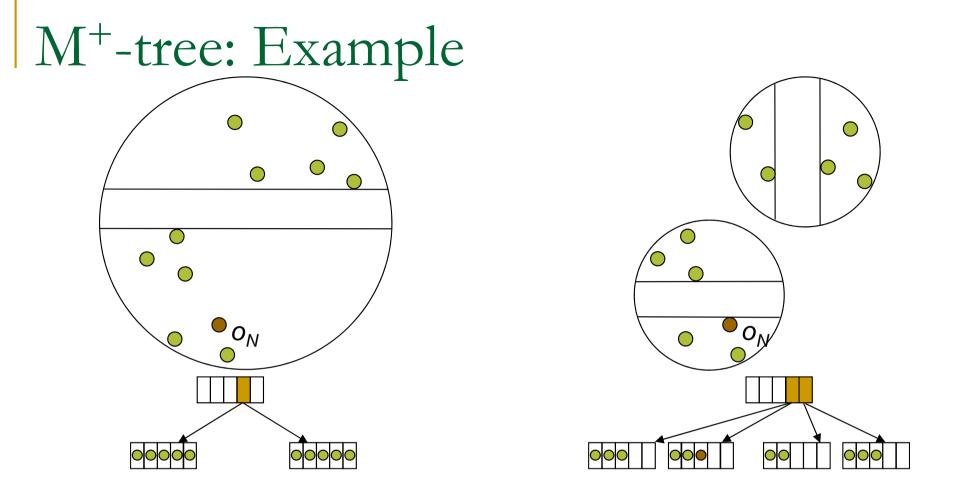
#### M<sup>+</sup>-tree: Structure

internal node is divided into two subsets

- according to a selected dimension
- leaving a gap between the two subsets
- the greater the gap the better filtering
- internal node entry:

 $\langle p, r^{c}, d(p, p^{p}), D_{key}, ptr_{left}, d_{lmax}, d_{rmin}, ptr_{right} \rangle$ 

- $\Box$   $D_{key}$  number of the key dimension
- $ptr_{left}$ ,  $ptr_{right}$  pointers to the left and right twin-nodes
- $d_{Imax} maximal$  key-dimension value of the left twin
- $\Box$   $d_{rmin}$  minimal key-dimension value of the right twin



splitting of an overfilled node:

- objects of both twins are considered as a single set
- apply standard *mM\_RAD\_2* strategy
- select the key dimension for each node separately

#### M<sup>+</sup>-tree: Performance

- slightly more efficient than M-tree
- better filtering for range queries with small radii
- practically the same for larger radii
- nearest neighbor queries:
  - □ a shorter *priority queue* only one of the twin-nodes
  - save some time for queue maintenance
- moderate performance improvements
- application restricted to vector datasets with L<sub>p</sub>

### M-tree Family

- The M-tree
- Bulk-Loading Algorithm
- Multi-Way Insertion Algorithm
- The Slim Tree
- Slim-Down Algorithm
  - Generalized Slim-Down Algorithm
- Pivoting M-tree
- The M<sup>+</sup>-tree
- The M<sup>2</sup>-tree

### The M<sup>2</sup>-tree

- generalization of M-tree
- able to process complex similarity queries
  - combined queries on several metrics at the same time
  - for instance: an image database with keyword-annotated objects and color histograms
  - query: Find images that contain a lion and the scenery around it like this.
- qualifying objects identified by a scoring function d<sub>f</sub>
  - combines the particular distances (according to several different measures)

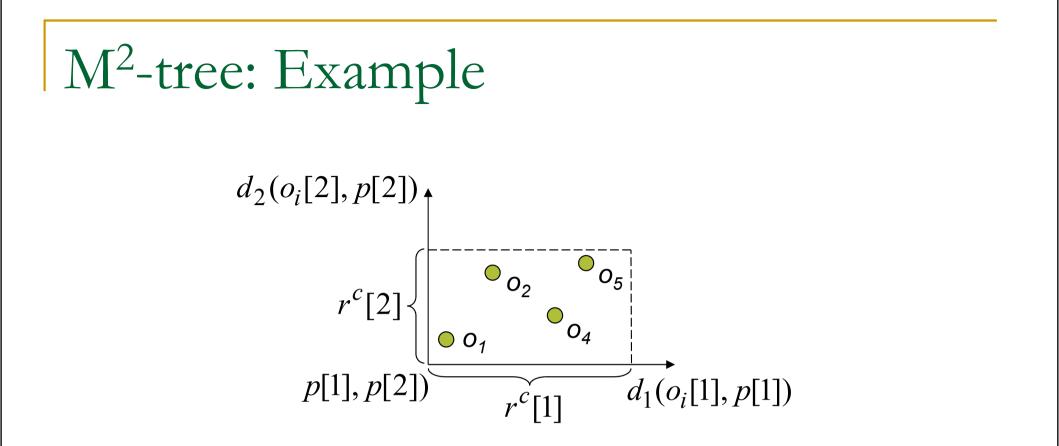
#### M<sup>2</sup>-tree: Structure

each object characterized by several *features* 

- e.g. *o[1],o[2]*
- respective distance measures may differ:  $d_1, d_2$
- leaf node: M-tree vs. M<sup>2</sup>-tree  $\langle o, d(o, p) \rangle$   $\langle o[1], d_1(o[1], p[1]), o[2], d_2(o[1], p[2]) \rangle$
- internal node: M-tree vs. M<sup>2</sup>-tree

 $\langle p, r^c, d(p, p^p), ptr \rangle$ 

 $\langle p[1], r^{c}[1], d_{1}(p[1], p^{p}[1]), p[2], r^{c}[2], d_{2}(p[2], p^{p}[2]), ptr \rangle$ 



- the space transformation according to particular features can be seen as an *n*-dimensional space
- the subtree region forms a hypercube

# M<sup>2</sup>-tree: Range Search

#### Given R(q,r):

- M-tree prunes a subtree if  $|d(q,p^p) d(p,p^p)| r^c > r$
- M<sup>2</sup>-tree: compute the lower bound for every feature  $\forall i, \min(|d_i(q[i], p^p[i]) - d_i(p[i], p^p[i])| - r^c[i], 0)$
- combine these bounds using the scoring function  $d_f$
- visit those entries for which the result is  $\leq r$
- analogous strategy for nearest neighbor queries

#### M<sup>2</sup>-tree: Performance

- running k-NN queries
- image database mentioned in the example
- M<sup>2</sup>-tree compared with sequential scan
  - the same I/O costs
  - reduced number of distance computations
- M<sup>2</sup>-tree compared with Fagin's  $\mathcal{A}_0$  (two M-trees)
  - M<sup>2</sup>-tree saves about 30% of I/Os
  - about 20% of distance computations
  - $\mathcal{A}_0$  have higher I/O cost than the sequential scan

Centralized Index Structures for Large Databases

1. M-tree family

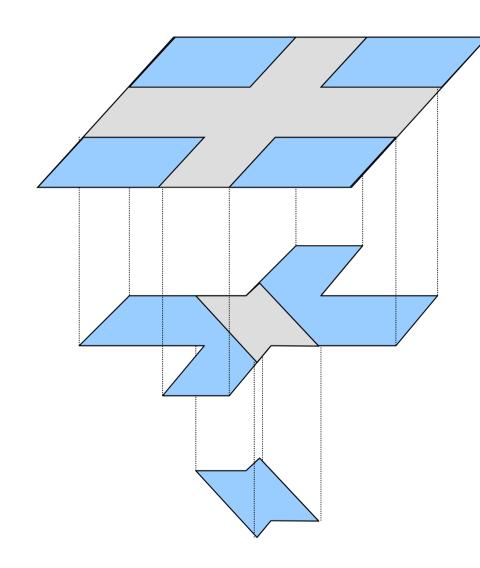
#### 2. hash-based metric indexing

- Distance Index (D-index)
- Extended D-Index (eD-index)
- 3. performance trials

## Distance Index (D-index)

- Hybrid structure
  - combines pivot-filtering and partitioning.
- Multilevel structure based on hashing
   one ρ-split function per level.
- The first level splits the whole data set.
- Next level partitions the exclusion zone of the previous level.
- The exclusion zone of the last level forms the exclusion bucket of the whole structure.

### D-index: Structure



4 separable buckets at the first level



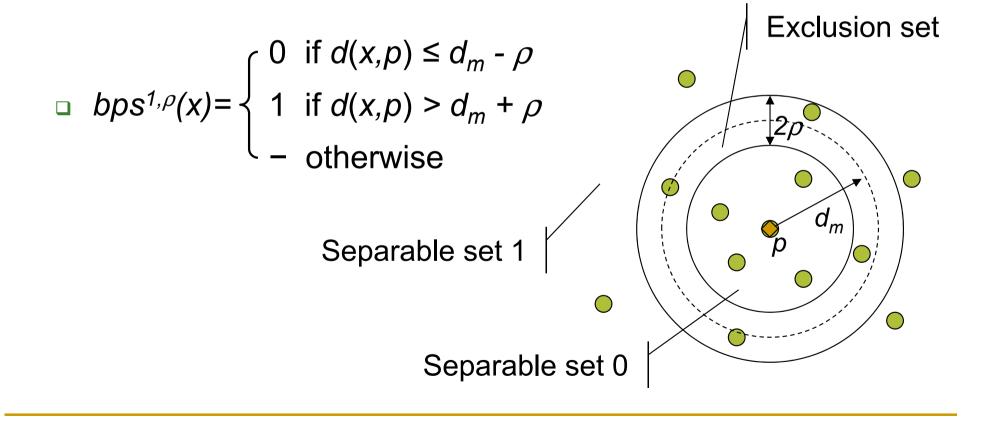
2 separable buckets at the second level



exclusion bucket of the whole structure

## D-index: Partitioning

Based on excluded middle partitioning
 ball partitioning variant is used.



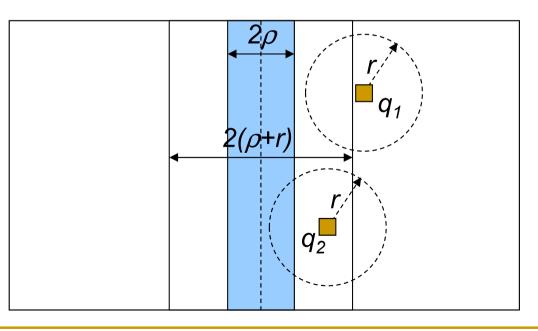
# D-index: Binary $\rho$ -Split Function

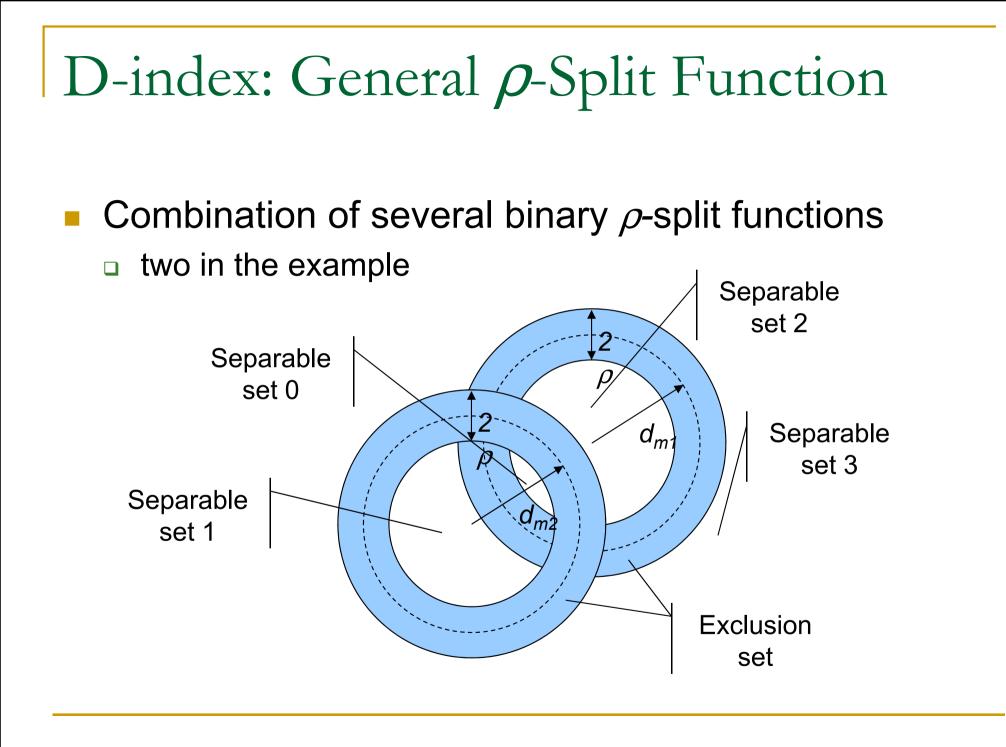
- Binary mapping:  $bps^{1,\rho}: \mathcal{D} \rightarrow \{0, 1, \neg\}$ 
  - □  $\rho$ -split function,  $\rho \ge 0$
  - $\Box$  also called the first order  $\rho$ -split function
- Separable property (up to  $2\rho$ ):  $\forall x, y \in \mathcal{D}, bps^{1,\rho}(x) = 0 \text{ and } bps^{1,\rho}(y) = 1 \Rightarrow d(x,y) > 2\rho$ 
  - No objects closer than 2p can be found in both the separable sets.
  - Symmetry property:  $\forall x, y \in \mathcal{D}, \rho_2 \ge \rho_1$ ,

 $bps^{1,\rho_2}(x) \neq -$ ,  $bps^{1,\rho_1}(y) = - \implies d(x,y) > \rho_2 - \rho_1$ 

# D-index: Symmetry Property

- We want to test whether a query intersects the exclusion set or not.





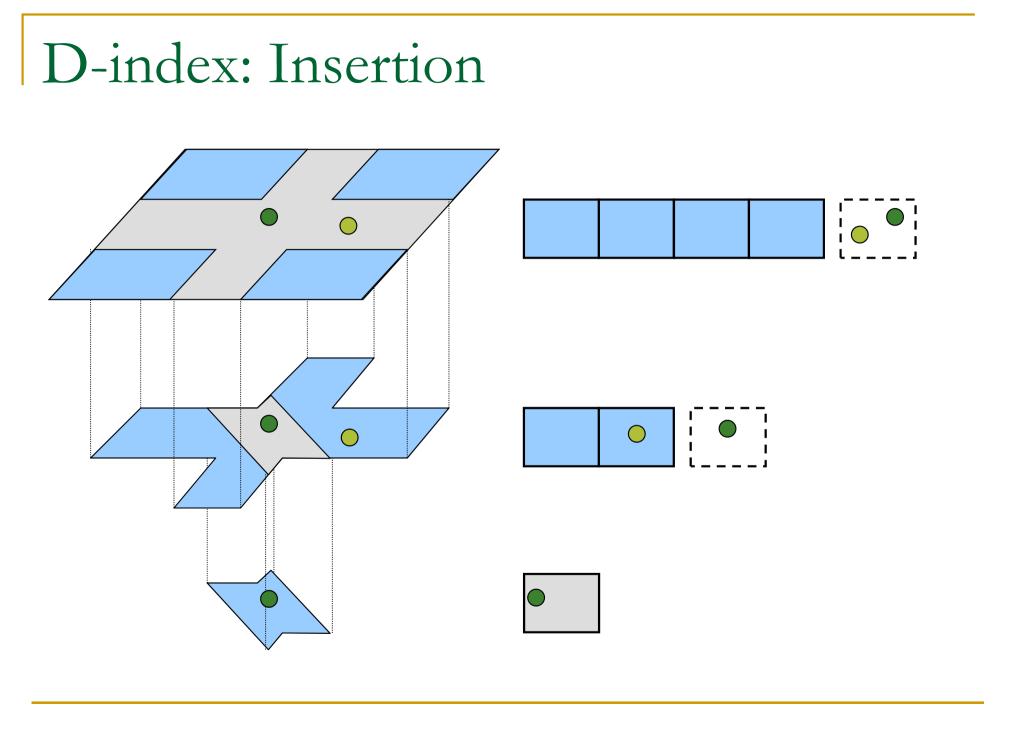
# D-index: General *p*-Split Function

A combination of *n* first order *ρ*-split functions:

□  $bps^{n,\rho}: \mathcal{D} \rightarrow \{0..2^{n}-1, -\}$ 

$$bps^{n,\rho}(x) = \begin{cases} - & \text{if } \exists i, bps_i^{1,\rho}(x) = - \\ b & \text{all } bps_i^{1,\rho}(x) \text{ form a binary number } b \end{cases}$$

Separable & symmetry properties hold
 resulting sets are also separable up to 2ρ.



## D-index: Insertion Algorithm

- Dindex $^{\rho}(X, m_1, m_2, ..., m_h)$ 
  - h number of levels,
  - $n_i$  number of binary functions combined on level *i*.
- Algorithm insert the object  $o_N$ :

```
for i=1 to h do
```

```
if bps^{m_i,\rho}(o_N) \neq - then
```

```
o_N \rightarrow bucket with the index bps<sup>m<sub>i</sub>, \rho</sup>(o_N).
```

exit

end if

end do

 $o_N \rightarrow global exclusion bucket.$ 

# D-index: Insertion Algorithm (cont.)

The new object is inserted with one bucket access.

Requires  $\sum_{i=1}^{j} m_i$  distance computations
 assuming  $o_N$  was inserted in a bucket on the level *j*.

# D-index: Range Query

Dindex<sup>p</sup>(X, m<sub>1</sub>, m<sub>2</sub>, ..., m<sub>h</sub>)

• h – number of levels,

•  $m_i$  – number of binary functions combined on level *i*.

#### Given a query R(q,r) with $r \leq \rho$ :

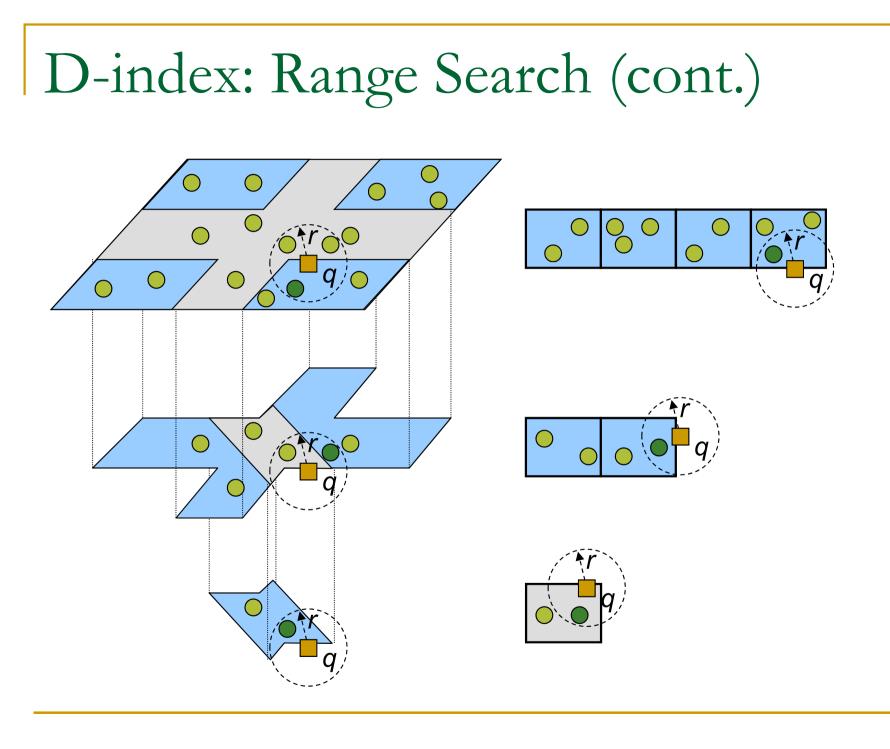
**for** *i*=1 **to** *h* **do** 

search in the bucket with the index  $bps^{m_i, 0}(q)$ .

end do

search in the global exclusion bucket.

□ Objects *o*,  $d(q,o) \le r$ , are reported on the output.



# D-index: Range Query (cont.)

- The call  $bps^{m_i,0}(q)$  always returns a value between 0 and  $2^{m_i}$ -1.
- Exactly one bucket per level is accessed if  $r \leq \rho$ 
  - h+1 bucket access.
- Reducing the number of bucket accesses:
  - the query region is in the exclusion set  $\Rightarrow$  proceed the next level directly,
  - the query region is in a separable set  $\Rightarrow$  terminate the search.

# D-index: Advanced Range Query

```
for i = 1 to h
      if bps^{m_i,\rho+r}(q) \neq - then (exclusively in the separable bucket)
           search in the bucket with the index bps^{m_i,\rho+r}(q).
                                           (search terminates)
          exit
      end if
      if r \leq \rho then
                                          (the search radius up to \rho)
           if bps^{m_{i,\rho}-r}(q) \neq - then (not exclusively in the exclusion zone)
               search in the bucket with the index bps^{m_i,\rho-r}(q).
          end if
                                           (the search radius greater than \rho)
      else
           let \{i_1, ..., i_n\} = G(bps^{m_{i,r-\rho}}(q))
           search in the buckets with the indexes i_1, \ldots, i_n.
      end if
end for
search in the global exclusion bucket.
```

#### D-index: Advanced Range Query (cont.)

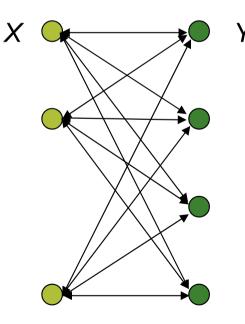
- The advanced algorithm is not limited to  $r \leq \rho$ .
- All tests for avoiding some bucket accesses are based on manipulation of parameters of split functions (i.e. *ρ*).
- The function *G*() returns a set of bucket indexes:
  - all minuses (-) in the split functions' results are substituted by all combinations of ones and zeros,
  - □ e.g. *bps*<sup>3,ρ</sup>(q)='1--'
  - $G(bps^{3,\rho}(q)) = \{100, 101, 110, 111\}$

### D-index: Features

- supports disk storage
- insertion needs one bucket access
  - distance computations vary from  $m_1$  up to  $\sum_{i=1..h} m_i$
- h+1 bucket accesses at maximum
  - $\hfill \hfill \hfill$
- exact match (R(q,0))
  - successful one bucket access
  - unsuccessful typically no bucket is accessed

# Similarity Join Query

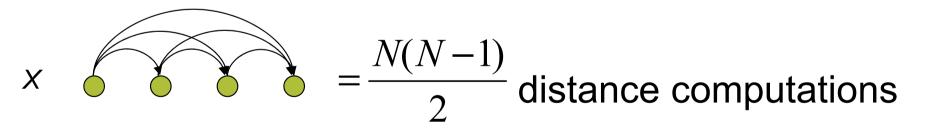
The similarity join can be evaluated by a simple algorithm which computes |X||Y| distances between all the pairs of objects.



= NM distance computations

# Similarity Self Join Query

- The similarity self join examines all pairs of objects of a set X, which is |X||X| distance computations.
- Due to the symmetry property, d(x,y) = d(y,x), we can reduce the costs.



This is called the nested loops algorithm (NL).

# Similarity Self Join Query (cont.)

#### Specialized algorithms

- usually built on top of a commercial DB system, or
- tailored to specific needs of application.
- D-index provides a very efficient algorithm for range queries:
  - a self join query can be evaluated using

Range Join Algorithm (RJ):

for each o in dataset X do

range\_query(o, μ)

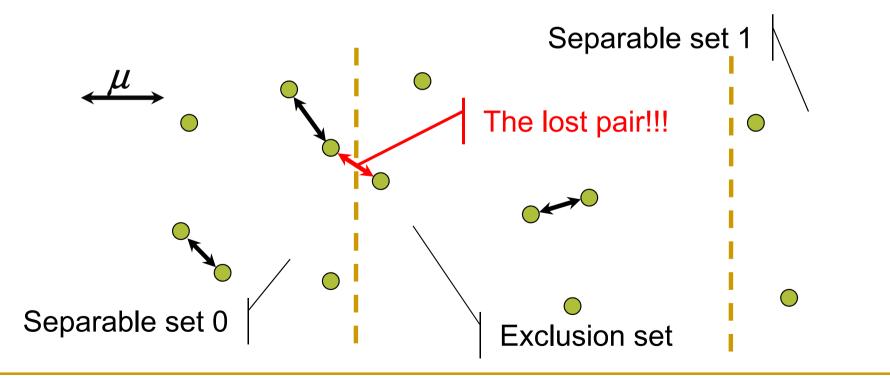
end do

# Extended D-index (eD-index)

- A variant of D-index which provides a specialized algorithm for similarity joins.
- Application independent general solution.
- Split functions manage replication.
- D-index's algorithms for range & k-NN queries are only slightly modified.

# eD-index: Similarity Self Join Query

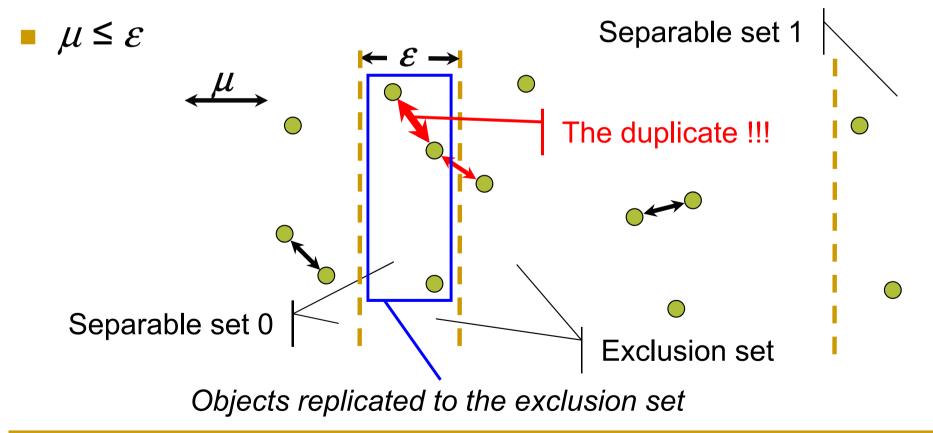
- Similarity self join is elaborated independently in each bucket.
- The result set is a union of answers of all sub-queries.

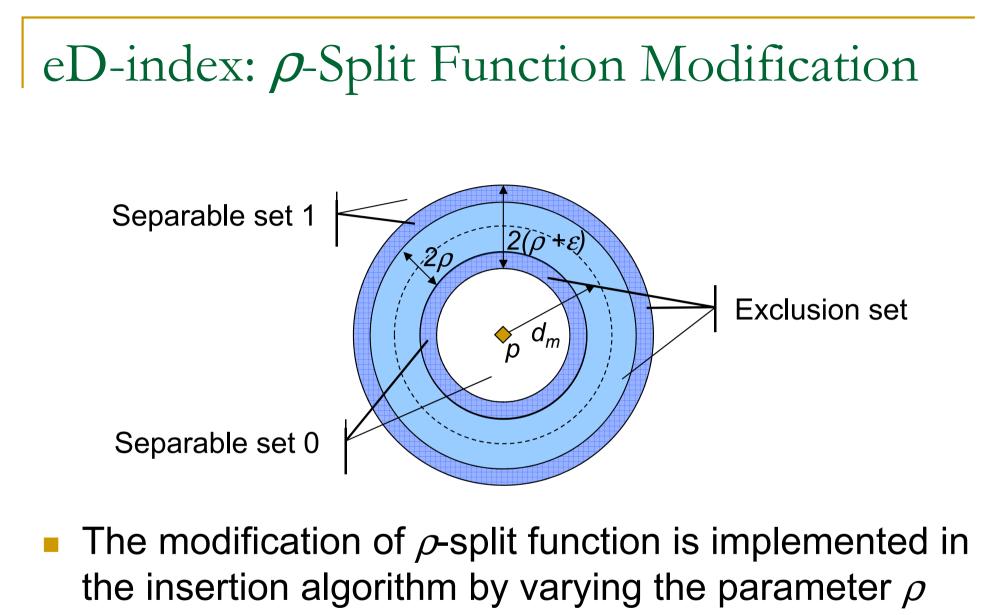


# eD-index: Overloading Principle

Lost pairs are handled by replications

 $\Box$  areas of width  $\varepsilon$  are replicated in the exclusion set.



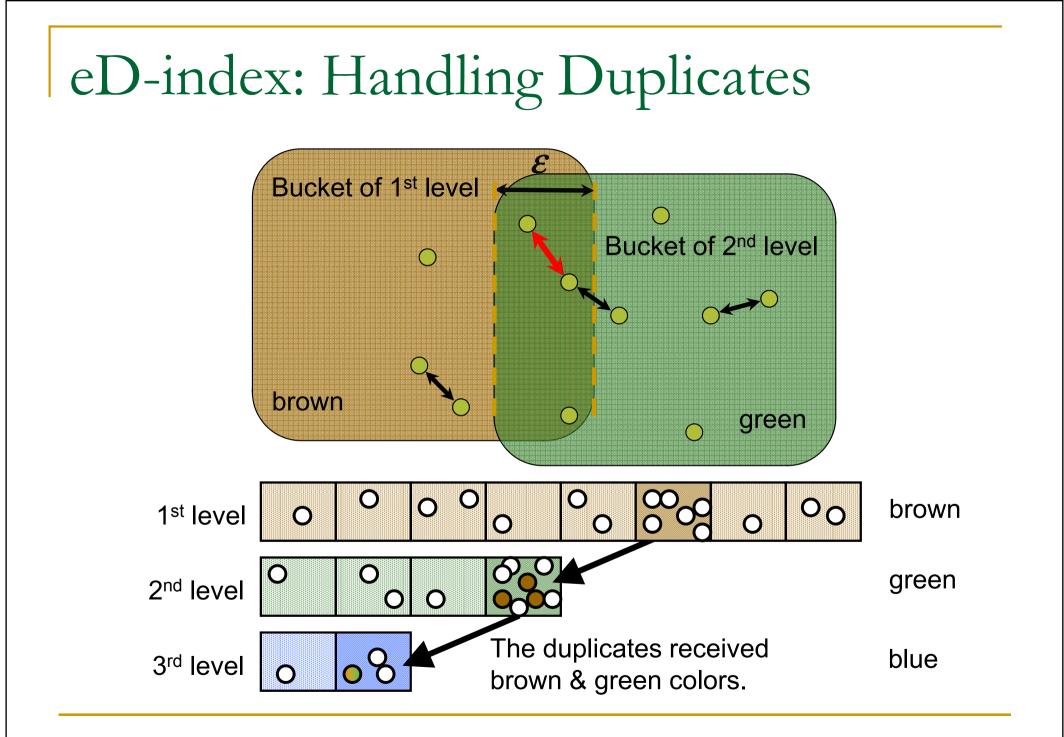


the original stop condition in the D-index's algorithm is changed.

## eD-index: Insertion Algorithm

- $eDindex^{\rho,\varepsilon}(X, m_1, m_2, ..., m_h)$
- Algorithm insert the object o<sub>N</sub>:
   for *i*=1 to *h* do

   if bps<sup>m<sub>i</sub>,ρ</sup>(o<sub>N</sub>) ≠ '-' then
   o<sub>N</sub> → bucket with the index bps<sup>m<sub>i</sub>,ρ</sup>(o<sub>N</sub>).
   if bps<sup>m<sub>i</sub>,ρ+ε</sup>(o<sub>N</sub>) ≠ '-' then (not in the overloading area)
   exit
   end if
   end if
   end if
  - $o_N \rightarrow global exclusion bucket.$



# eD-index: Overloading Join Algorithm

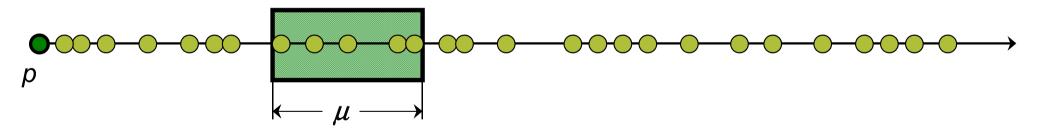
Given similarity self-join query  $SJ(\mu)$ :

- Execute the query in every separable bucket on every level
  - and in the global exclusion bucket.
- In the bucket, apply sliding window algorithm.
- The query's result is formed by concatenation of all sub-results.

# eD-index: Sliding Window

Use the triangle inequality

- to avoid checking all pairs of objects in the bucket.
- Order all objects on distances to one pivot.
- The sliding window is then moved over all objects.
   only pairs of objects in the window are examined.



- Due to the triangle inequality, the pair of objects outside the window cannot qualify:
  - $\Box \quad d(x,y) \geq d(x,p) d(y,p) > \mu$

# eD-index: Sliding Window (cont.)

#### The algorithm also employs

- the pivot filtering and
- the eD-index's coloring technique.

#### Given a pair of objects $o_1, o_2$ :

- if a color is shared, this pair must have been reported on the level having this color – the pair is ignored without distance computation, else
- □ if  $d(o_1, o_2) \le \mu$ , it is an original qualifying pair.

#### eD-index: Limitations

- Similarity self-join queries only
  - the query selectivity must satisfy:  $\mu \leq \varepsilon$ .
  - □ it is not very restrictive since we usually look for close pairs.
- The parameters  $\rho$  and  $\varepsilon$  depend on each other.
  - $\Box \quad \mathcal{E} \leq 2\rho$
  - If  $\varepsilon > 2\rho$ , the overloading zone is wider than the exclusion zone.
    - because we do not replicate objects between separable sets only between a separable set and the exclusion zone,
    - some qualifying pairs might be missed.

Centralized Index Structures for Large Databases

- 1. M-tree family
- 2. hash-based metric indexing
- **3.** performance trials

### Performance Trials

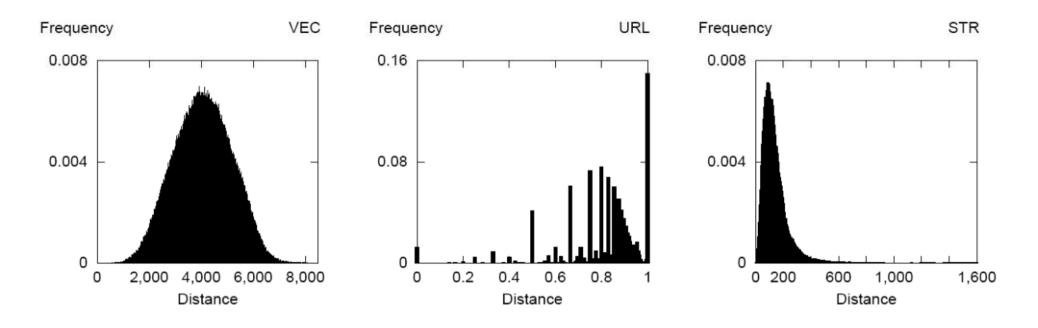
- experiments on M-tree and D-index
- three sets of experiments:
  - comparison of M-tree (tree-based approach) vs. D-index (hash-based approach)
  - 2. processing different **types of queries**
  - scalability of the centralized indexes growing the size of indexed dataset

#### Datasets and Distance Measures

trials performed on three datasets:

- VEC: 45-dimensional vectors of image color features compared by the *quadratic distance* measure
- URL: sets of URL addresses; the distance measure is based on the similarity of sets (*Jaccard's coefficient*)
- STR: sentences of a Czech language corpus compared using an *edit distance*

### Datasets: Distance Distribution



distribution of distances within the datasets:

- VEC: practically normal distance distribution
- URL: discrete distribution
- STR: skewed distribution

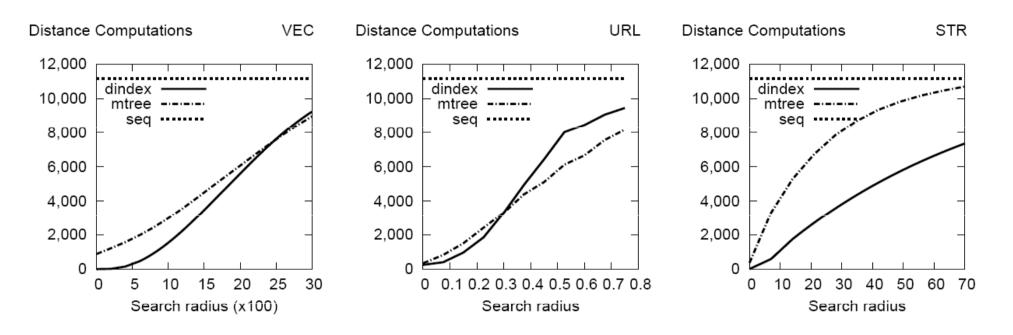
## Trials: Measurements & Settings

- CPU costs: number of distance computations
- I/O costs: number of block reads
  - The same size of disk blocks
- Query objects follow the dataset distribution
- Average values over 50 queries:
  - Different query objects
  - The same selectivity
    - Radius or number of nearest neighbors

## Comparison of Indexes

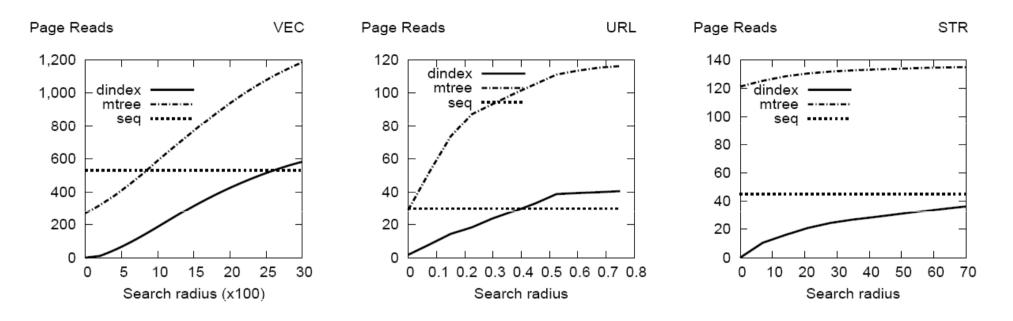
- Comparing performance of
  - □ M-tree a tree-based approach
  - D-index hash-based approach
  - sequential scan (baseline)
- Dataset of 11,100 objects
- Range queries increasing radius
  - maximal selectivity about 20% of the dataset

## Comparison: CPU Costs



- generally, D-index outperforms M-tree for smaller radii
- D-index: pivot-based filtering depends on data distribution and query size
- M-tree outperforms D-index for discrete distribution
  - pivot selection is more difficult for discrete distributions

# Comparison: I/O Costs

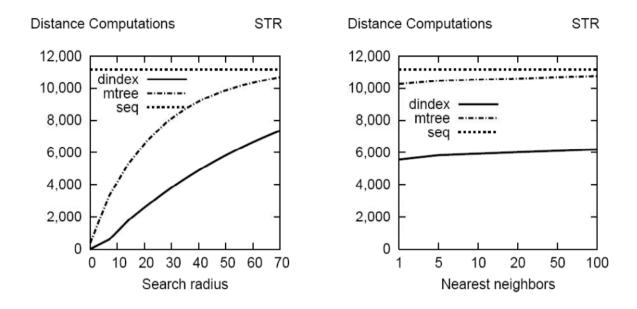


- M-tree needs twice the disk space to stored data than SEQ
- □ inefficient if the *distance function* is easy to compute
- D-index more efficient
- a query with *r*=0: D-index accesses only one page (important, e.g., for deletion)

# Different Query Types

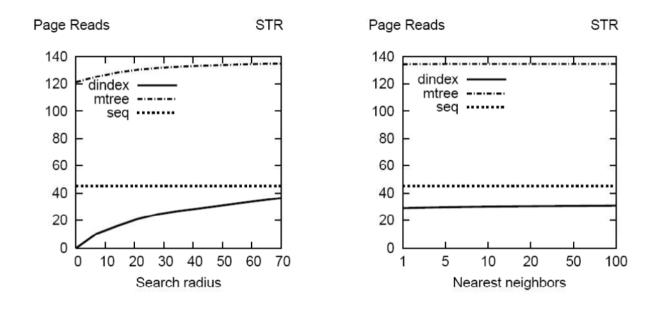
- comparing processing performance of different types of queries
  - range query
  - nearest neighbor query
  - similarity self join
- M-tree, D-index, sequential scan

# Range vs. k-NN: CPU Costs



- nearest neighbor query:
  - similar trends for M-tree and D-index
  - □ the D-index advantage of small radii processing decreases
  - expensive even for small k similar costs for both 1 and 100
  - D-index still twice as fast as M-tree

## Range vs. k-NN: I/O Costs



- nearest neighbor query:
  - similar trends for I/O costs as for CPU costs
  - D-index four times faster than M-tree

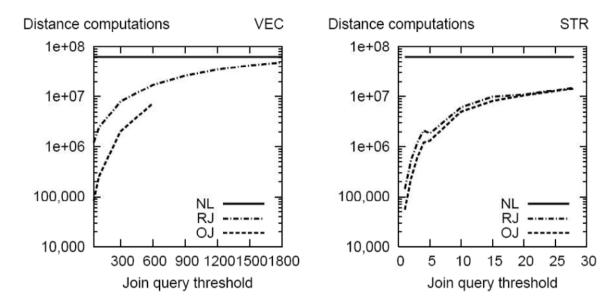
# Similarity Self Join: Settings

- $J(X,X,\mu)$  very demanding operation
- three algorithms to compare:
  - NL: nested loops naive approach
  - □ RJ: range join based on D-index
  - OJ: overloading join eD-index
    - for  $\mu$ :  $2\mu \le \rho$ , i.e.  $\mu \le 600$  for vectors
- datasets of about 11,000 objects
- selectivity retrieving up to 1,000,000 pairs (for high values of μ)

# Similarity Self Join: Complexity

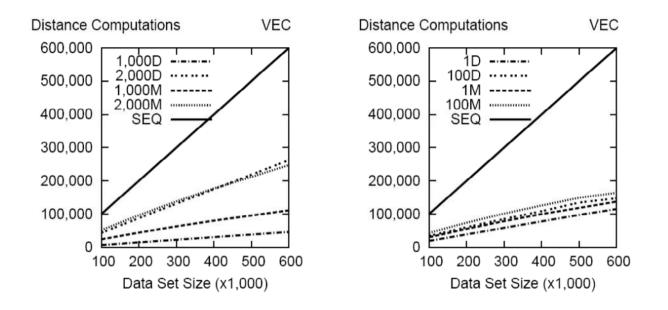
- Quadratic complexity
  - prohibitive for large DB
  - example: 50,000 sentences
  - a range query:
    - sequential scan takes about 16 seconds
  - □ a self join query:
    - nested loops algorithm takes 25,000 times more
    - about 4 days and 15 hours!

# Similarity Join: Results



- RJ and OJ costs increase rapidly (logarithmic scale)
- OJ outperforms RJ twice (STR) and 7 times for VEC:
  - high distances between VEC objects
  - high pruning effectiveness of pivot-based filtering for smaller  $\mu$

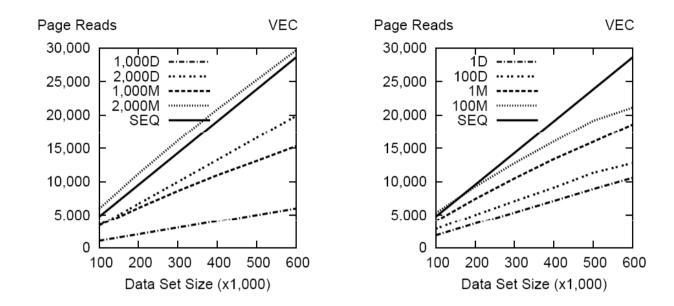
### Scalability: CPU Costs



□ range query: r = 1,000; 2,000 □ k-NN query: k = 1; 100

- Iabels: radius or k + D (D-index), M (M-tree), SEQ
- data: from 100,000 to 600,000 objects
- M-tree and D-index are faster (D-index slightly better)linear trends

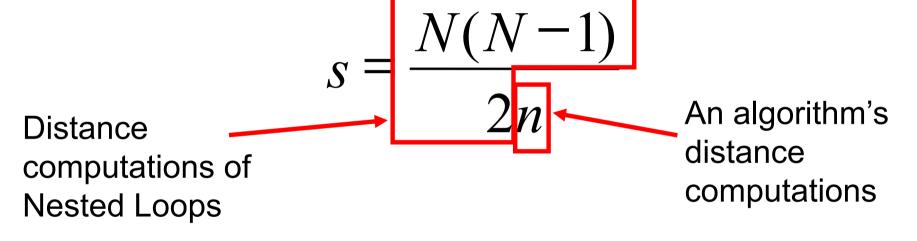
# Scalability: I/O Costs



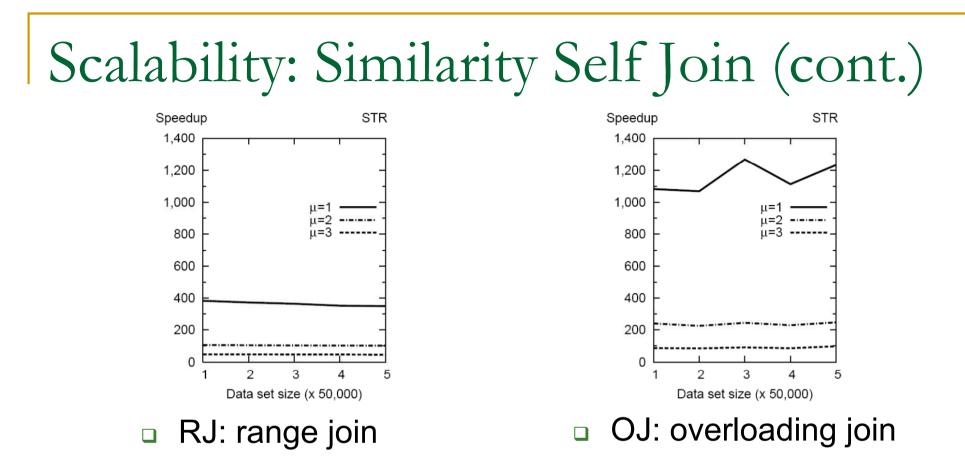
- the same trends as for CPU costs
- D-index more efficient than M-tree
- *exact match* contrast:
  - □ M-tree: 6,000 block reads + 20,000 d. c. for 600,000 objects
  - D-index: read 1 block + 18 d. c. regardless of the data size

# Scalability: Similarity Self Join

We use the speedup s as the performance measure:



Speedup measures how many times is a specific algorithm faster than NL.



#### STR dataset: from 50,000 to 250,000 sentences

- constant speedup
  - □ E.g. a join query on 100,000 objects takes 10 minutes.
  - The same join query on 200,000 objects takes 40 minutes.
- OJ at least twice faster than RJ

# Scalability Experiments: Conclusions

- similarity search is expensive
- the scalability of centralized indexes is linear
- cannot be applied to huge data archives
   become inefficient after a certain point

Possible solutions:

- sacrifice some precision: approximate techniques
- use more storage & computational power:
   distributed data structures

