

Lecture outline

- Nearest-neighbor search in low dimensions
 - kd-trees
- Nearest-neighbor search in high dimensions
 - LSH
- Applications to data mining

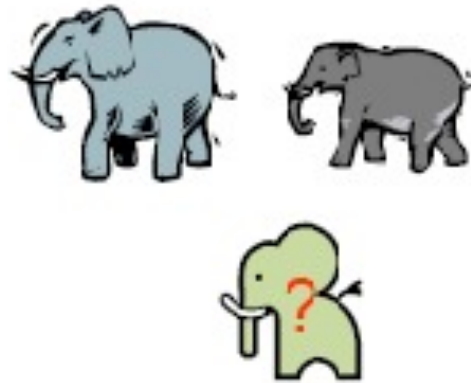
Definition

- Given: a set X of n points in \mathbb{R}^d
- Nearest neighbor: for any query point $q \in \mathbb{R}^d$ return the point $x \in X$ minimizing $D(x, q)$
- **Intuition:** Find the point in X that is the closest to q

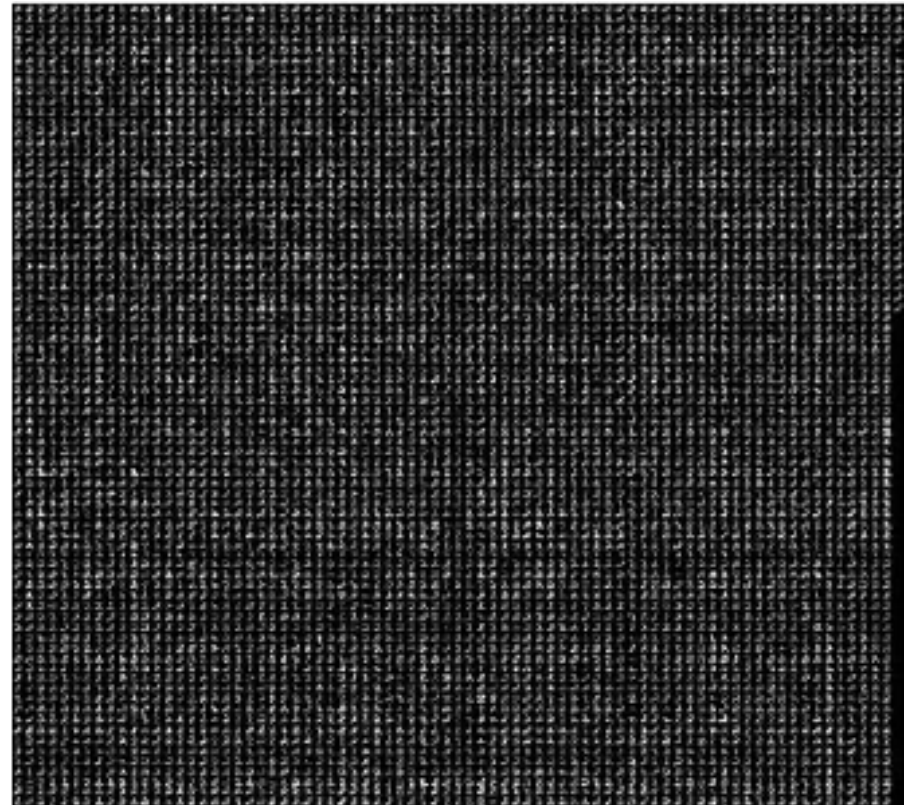
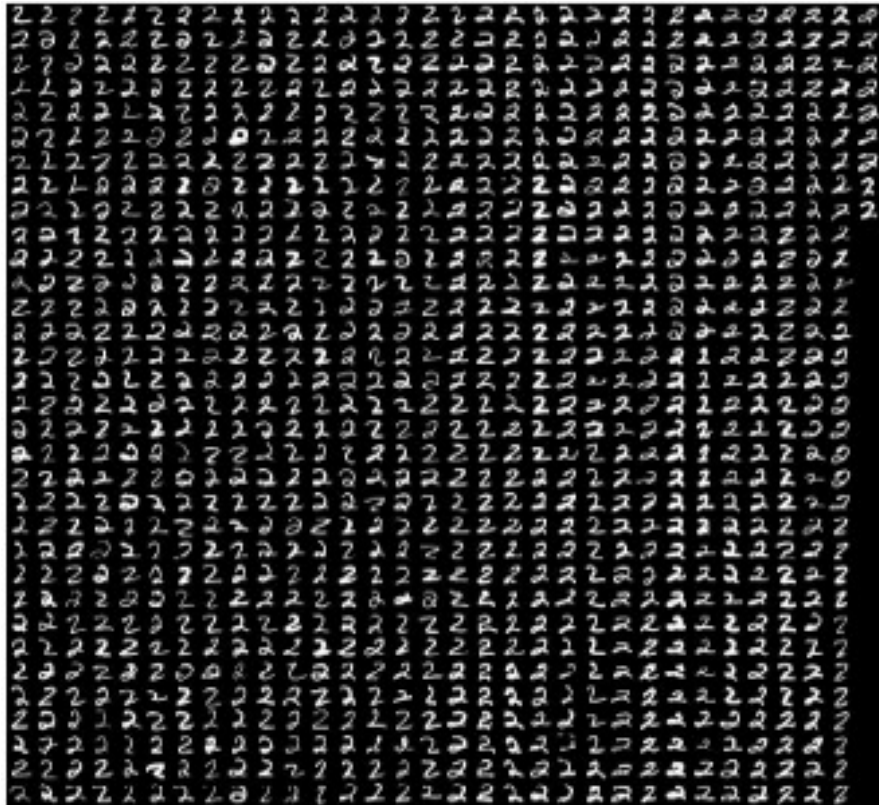
Motivation

- **Learning:** Nearest neighbor rule
- **Databases:** Retrieval
- **Data mining:** Clustering
- Donald Knuth in vol.3 of **The Art of Computer Programming** called it the post-office problem, referring to the application of assigning a resident to the **nearest-post office**

Nearest-neighbor rule



MNIST dataset “2”



Methods for computing NN

- **Linear scan:** $O(nd)$ time
- This is pretty much all what is known for exact algorithms with theoretical guarantees
- In practice:
 - **kd-trees** work “well” in “low-medium” dimensions

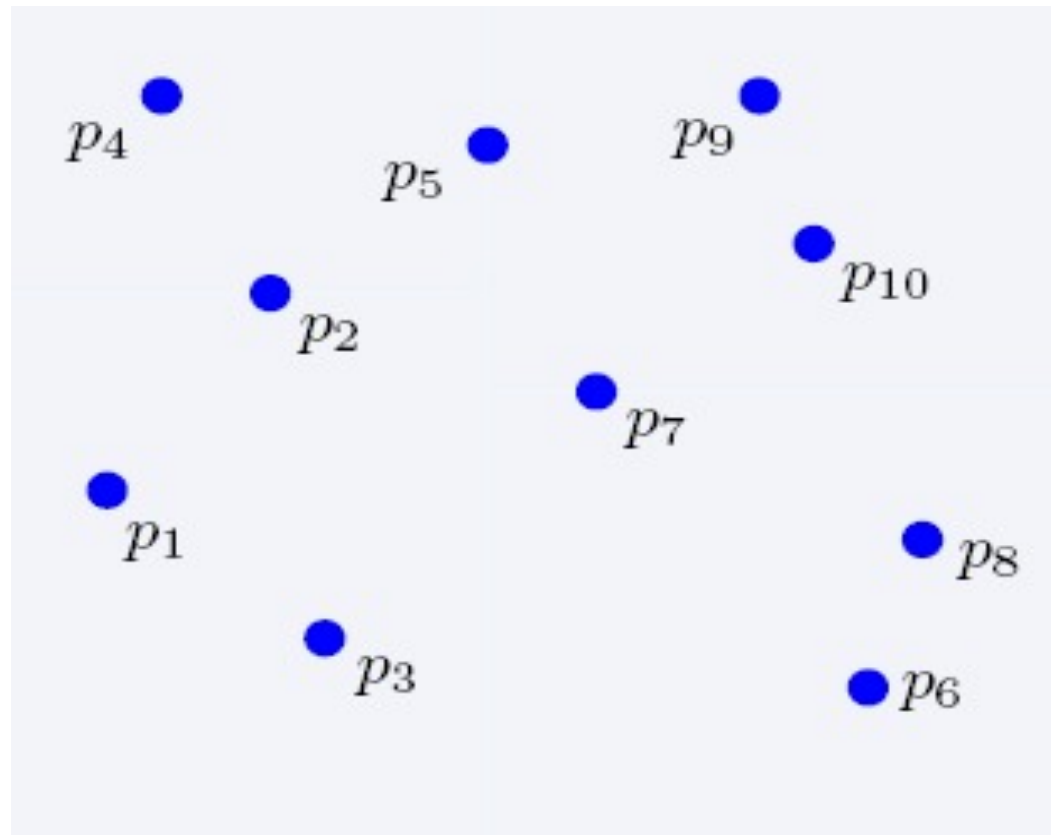
2-dimensional kd-trees

- A data structure to support range queries in \mathbf{R}^2
 - Not the most efficient solution in theory
 - Everyone uses it in practice
- Preprocessing time: $\mathbf{O(n \log n)}$
- Space complexity: $\mathbf{O(n)}$
- Query time: $\mathbf{O(n^{1/2} + k)}$

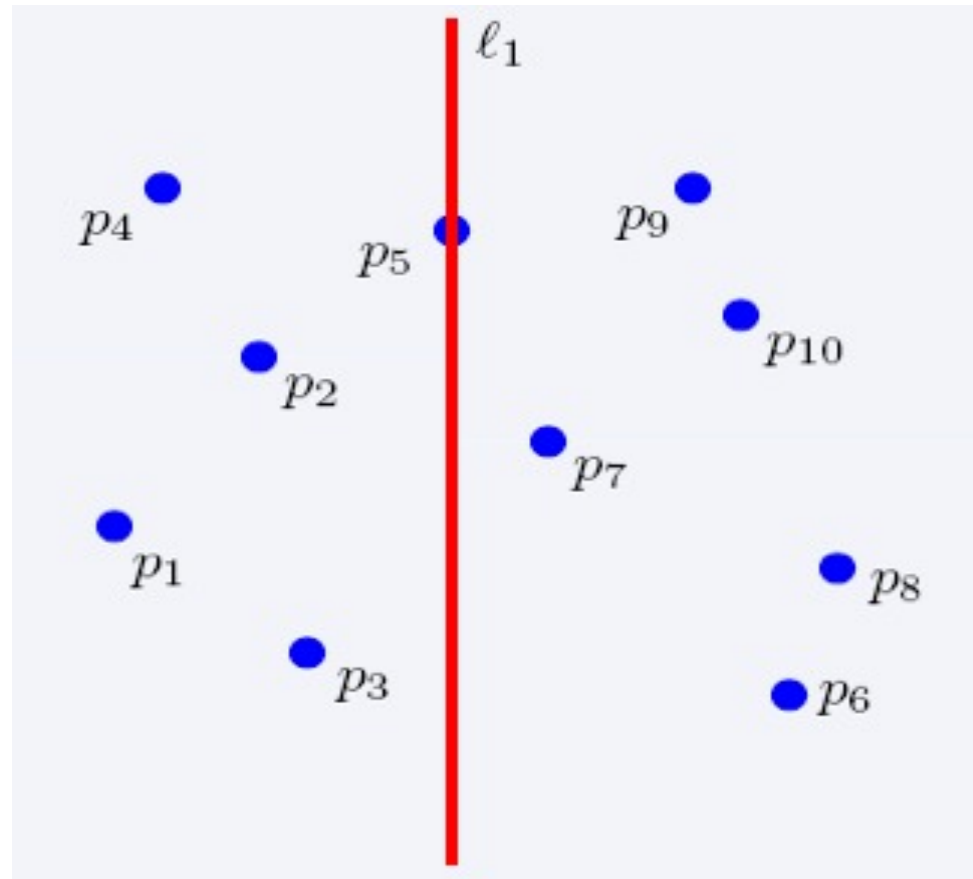
2-dimensional kd-trees

- Algorithm:
 - Choose **x** or **y** coordinate (alternate)
 - Choose the median of the coordinate; this defines a horizontal or vertical line
 - Recurse on both sides
- We get a binary tree:
 - Size **$O(n)$**
 - Depth **$O(\log n)$**
 - Construction time **$O(n \log n)$**

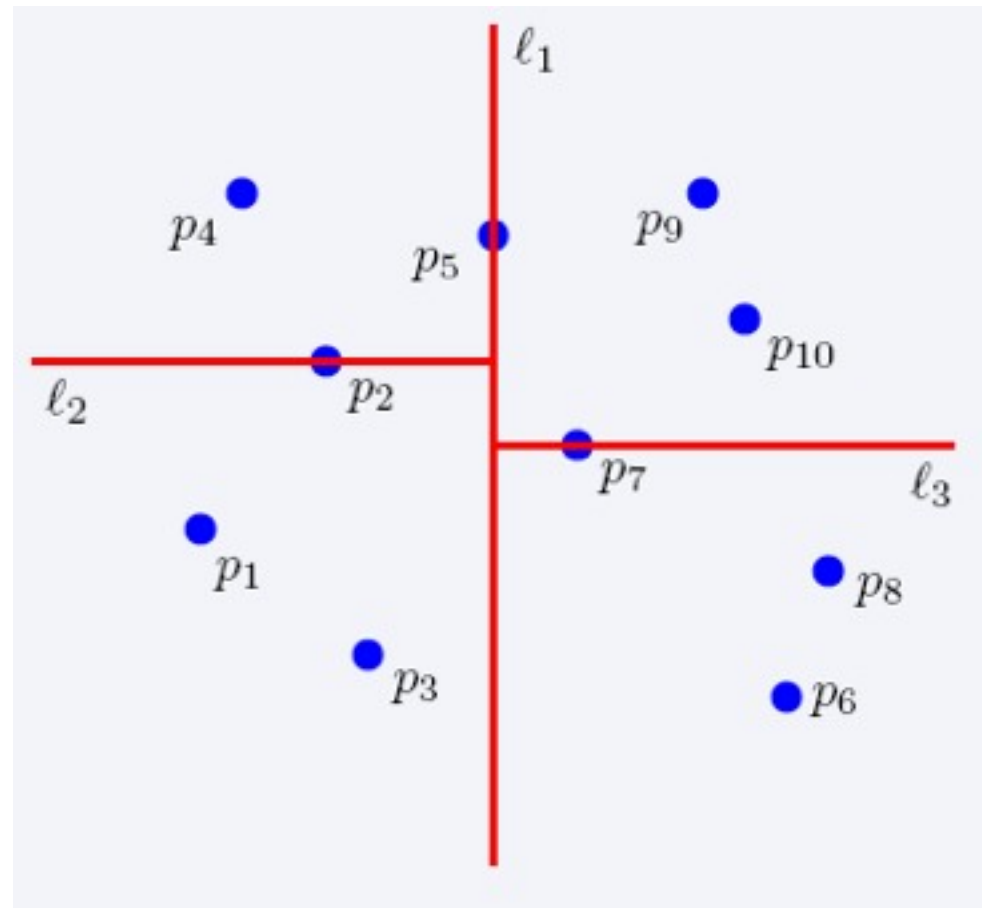
Construction of kd-trees



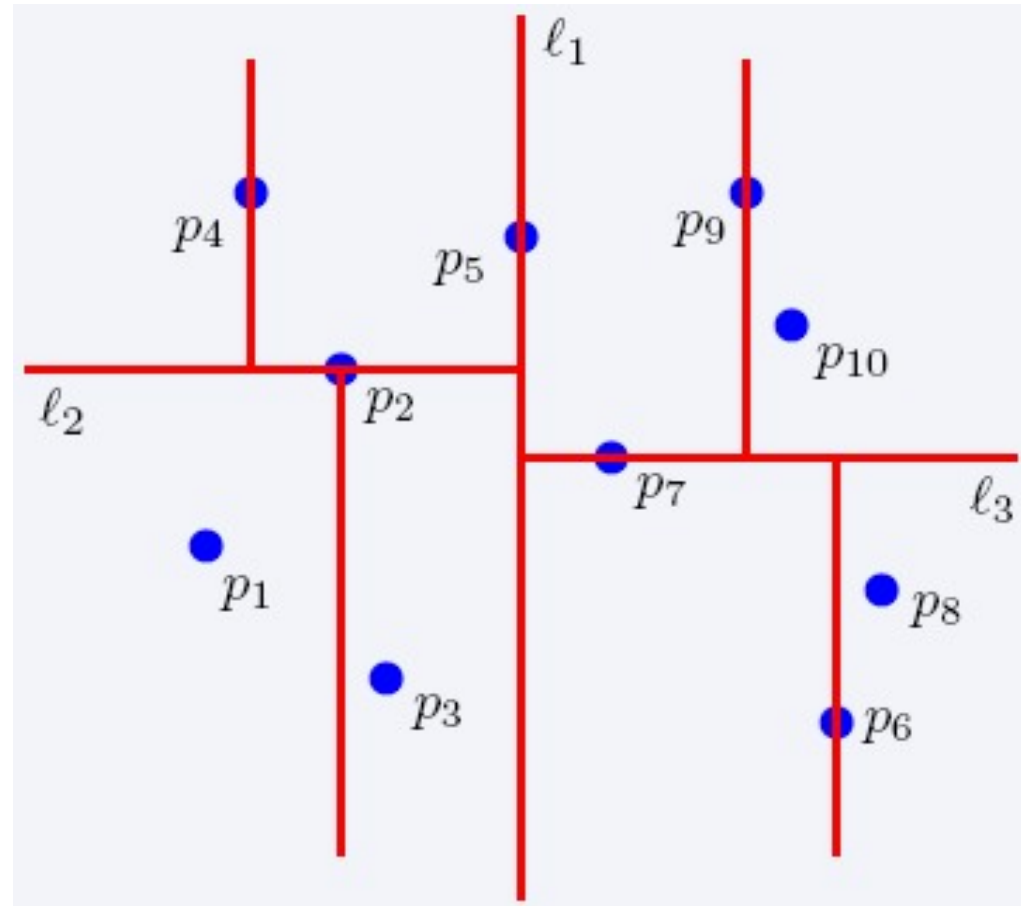
Construction of kd-trees



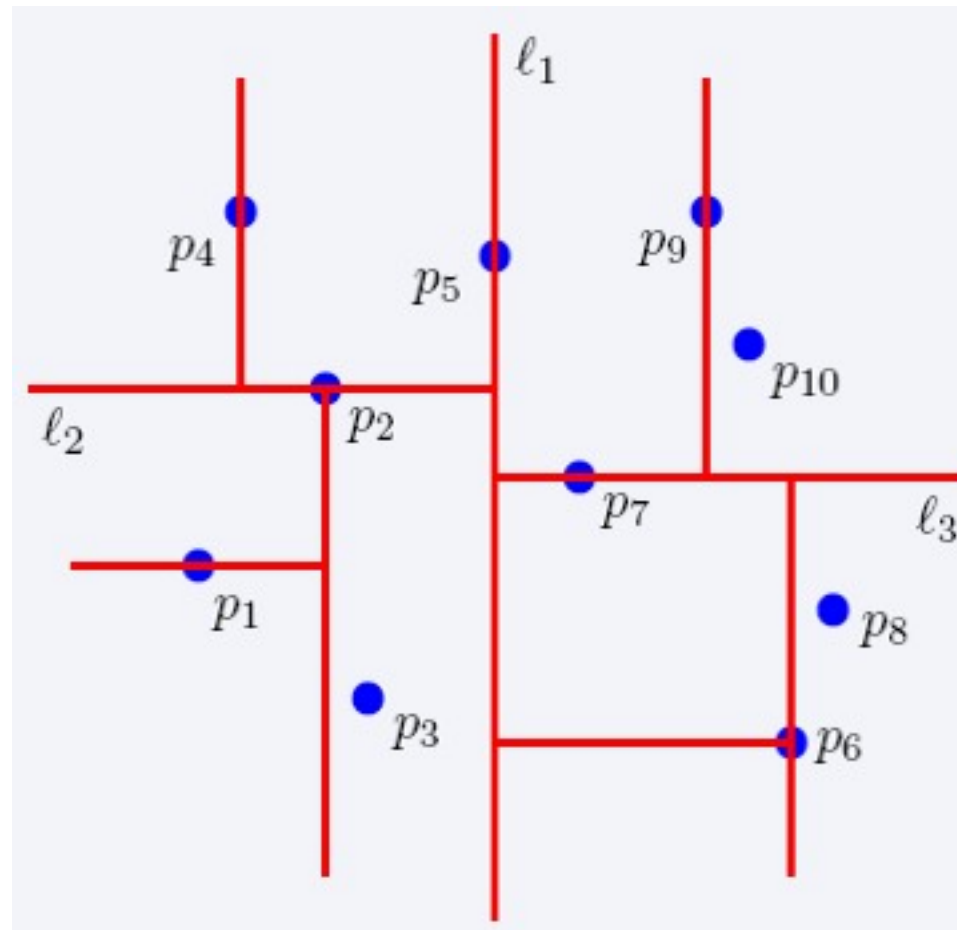
Construction of kd-trees



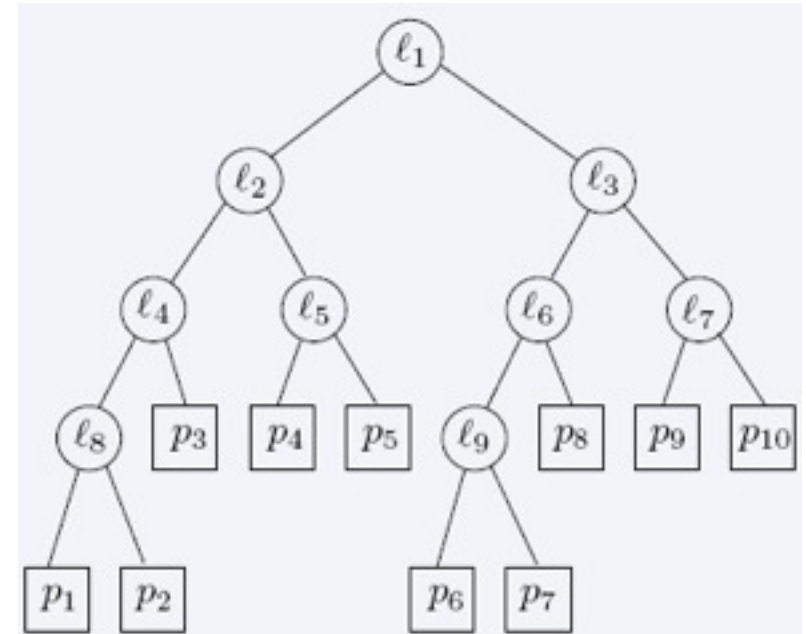
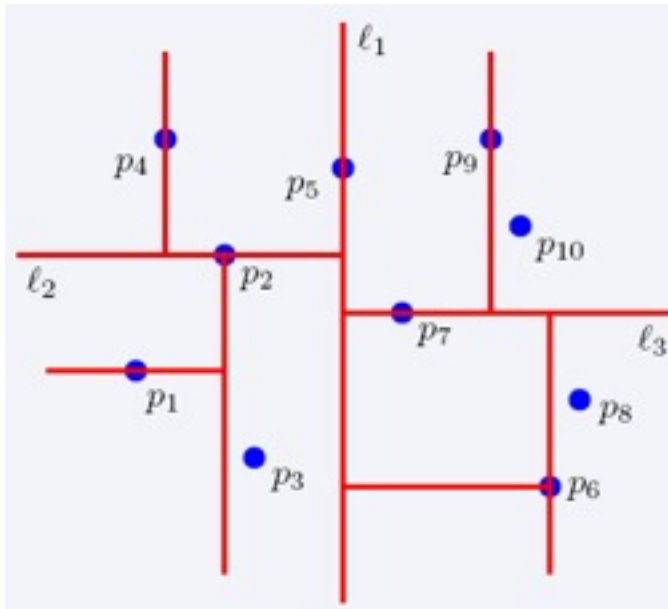
Construction of kd-trees



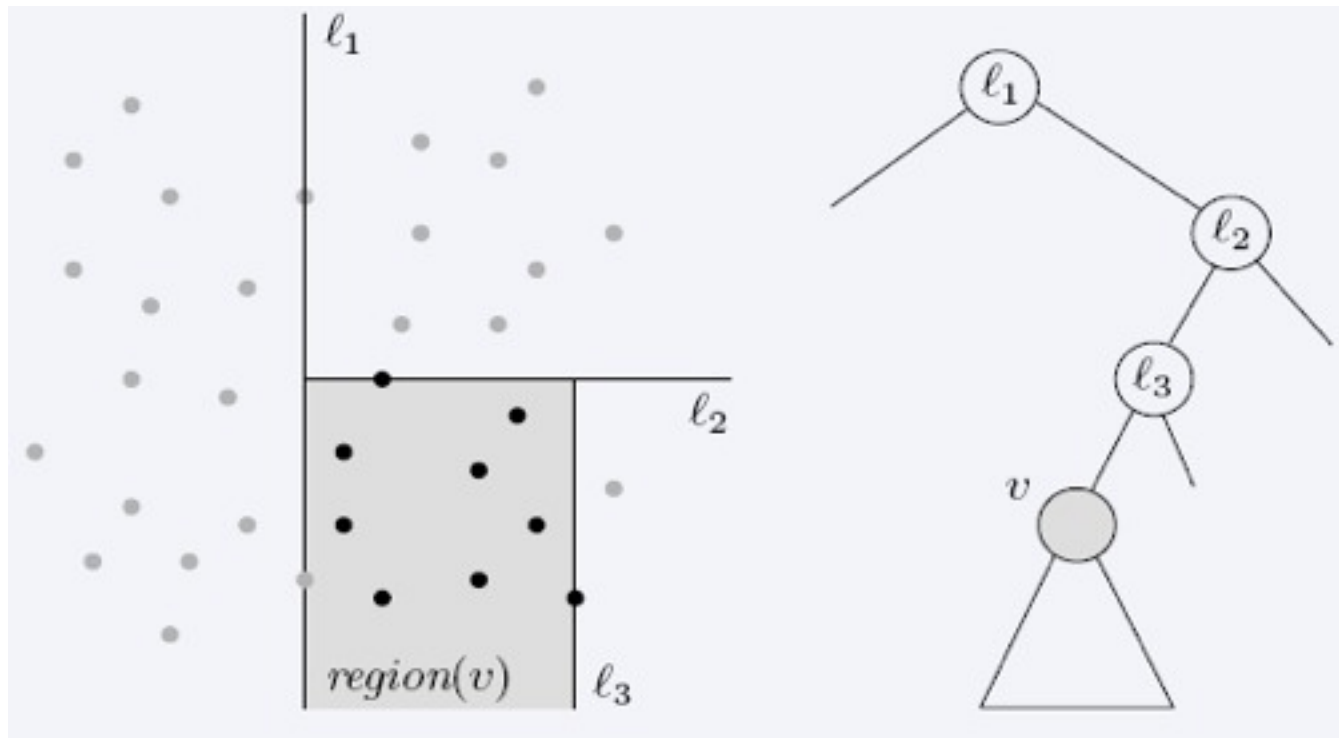
Construction of kd-trees



The complete kd-tree



Region of node v



Region(v) : the subtree rooted at v stores the points in black dots

Searching in kd-trees

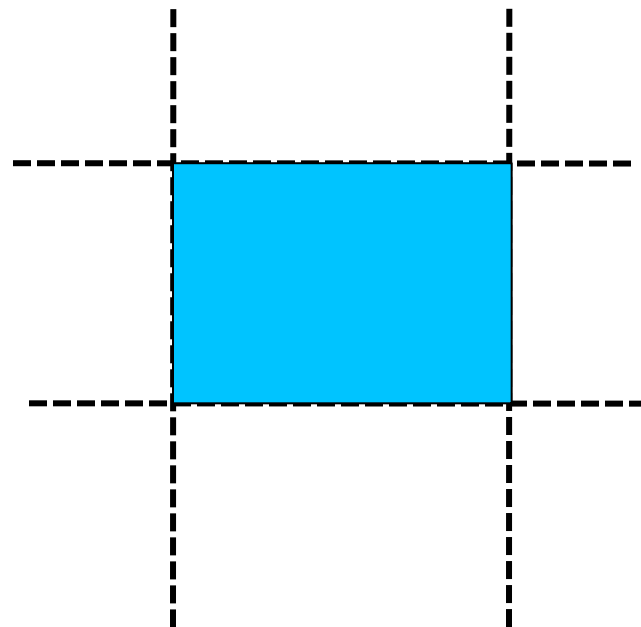
- Range-searching in **2-d**
 - Given a set of **n** points, build a data structure that for any query rectangle **R** reports all point in **R**

kd-tree: range queries

- Recursive procedure starting from $v = \text{root}$
- **Search** (v, R)
 - If v is a leaf, then report the point stored in v if it lies in R
 - Otherwise, if $\text{Reg}(v)$ is contained in R , report all points in the **subtree**(v)
 - Otherwise:
 - If $\text{Reg}(\text{left}(v))$ intersects R , then **Search**($\text{left}(v), R$)
 - If $\text{Reg}(\text{right}(v))$ intersects R , then

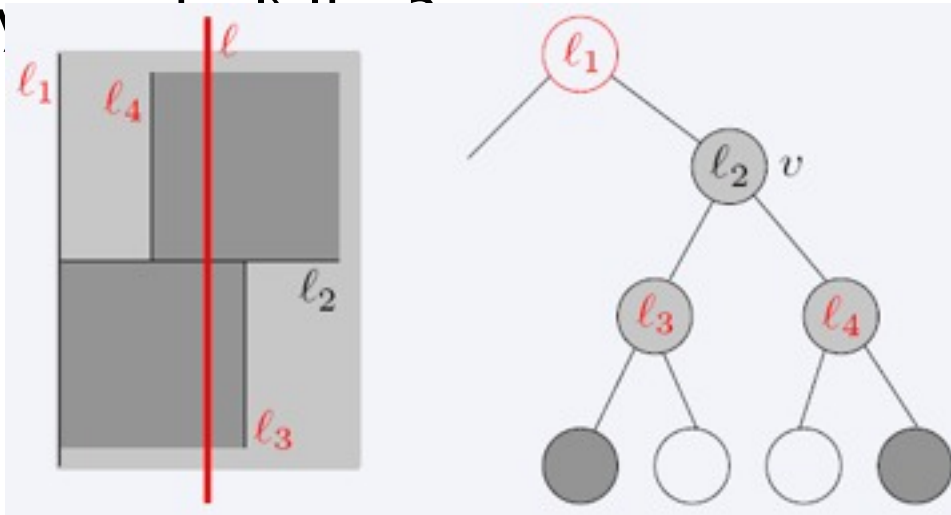
Query time analysis

- We will show that **Search** takes at most $O(n^{1/2} + P)$ time, where P is the number of reported points
 - The total time needed to report all points in all sub-trees is $O(P)$
 - We just need to bound the number of nodes v such that **region(v)** intersects R but is not contained in R (i.e., boundary of R intersects the boundary of **region(v)**)
 - **gross overestimation**: bound the number of **region(v)** which are crossed by any of the **4** horizontal/vertical lines



Query time (Cont'd)

- $Q(n)$: max number of regions in an n -point kd-tree intersecting a (say) vertical line ℓ



- If ℓ intersects $\text{region}(v)$ (due to vertical line splitting), then after two levels it intersects **2** regions (due to 2 vertical splitting lines)
- The number of regions intersecting ℓ is $Q(n) = 2 + 2Q(n/4) \rightarrow Q(n) = (n^{1/2})$

d-dimensional kd-trees

- A data structure to support range queries in \mathbb{R}^d
- Preprocessing time: $O(n \log n)$
- Space complexity: $O(n)$
- Query time: $O(n^{1-1/d} + k)$

Construction of the d -dimensional kd-trees

- The construction algorithm is similar as in $2-d$
- At the root we split the set of points into two subsets of same size by a hyperplane vertical to x_1 -axis
- At the children of the root, the partition is based on the second coordinate: x_2 -coordinate
- At depth d , we start all over again by partitioning on the first coordinate
- The recursion stops until there is only one

Locality-sensitive hashing (LSH)

- **Idea:** Construct hash functions $h: \mathbb{R}^d \rightarrow \mathcal{U}$ such that for any pair of points p, q :
 - If $D(p, q) \leq r$, then $\Pr[h(p) = h(q)]$ is high
 - If $D(p, q) \geq cr$, then $\Pr[h(p) = h(q)]$ is small
- Then, we can solve the “approximate NN” problem by hashing
- LSH is a general framework; for a given D we need to find the right h

Approximate Nearest Neighbor

- Given a set of points X in \mathbb{R}^d and query point $q \in \mathbb{R}^d$
 c -Approximate r -Nearest Neighbor search returns:
 - Returns $p \in P, D(p, q) \leq r$
 - Returns NO if there is no $p' \in X, D(p', q) \leq cr$

Locality-Sensitive Hashing (LSH)

- A family \mathbf{H} of functions $\mathbf{h}: \mathbf{R}^d \rightarrow \mathbf{U}$ is called (P_1, P_2, r, cr) -sensitive if for any p, q :
 - if $\mathbf{D}(p, q) \leq r$, then $\Pr[\mathbf{h}(p) = \mathbf{h}(q)] \geq P_1$
 - If $\mathbf{D}(p, q) \geq cr$, then $\Pr[\mathbf{h}(p) = \mathbf{h}(q)] \leq P_2$
- $P_1 > P_2$
- Example: **Hamming** distance
 - LSH functions: $\mathbf{h}(p) = p_i$, i.e., the i -th bit of p
 - Probabilities: $\Pr[\mathbf{h}(p) = \mathbf{h}(q)] = 1 - \mathbf{D}(p, q) / d$

Algorithm -- preprocessing

- $g(p) = \langle h_1(p), h_2(p), \dots, h_k(p) \rangle$
- Preprocessing
 - Select g_1, g_2, \dots, g_L
 - For all $p \in X$ hash p to buckets $g_1(p), \dots, g_L(p)$
 - Since the number of possible buckets might be large we only **maintain the non empty ones**
- Running time?

Algorithm -- query

- Query q :
 - Retrieve the points from buckets $g_1(q), g_2(q), \dots, g_L(q)$ and let points retrieved be x_1, \dots, x_L
 - If $D(x_i, q) \leq r$ report it
 - Otherwise report that there does not exist such a NN
 - Answer the query based on the retrieved points
 - Time $O(dL)$

Applications of LSH in data mining

- Numerous.....

Applications

- Find pages with similar sets of words (for clustering or classification)
- Find users in Netflix data that watch similar movies
- Find movies with similar sets of users
- Find images of related things

How would you do it?

- Finding very similar items might be computationally demanding task
- We can relax our requirement to finding **somewhat similar** items

Running example: comparing documents

- Documents have common text, but no common topic
- Easy special cases:
 - Identical documents
 - Fully contained documents (letter by letter)
- General case:
 - Many small pieces of one document appear out of order in another. What do we do then?

Finding similar documents

- Given a collection of documents, find pairs of documents that have lots of text in common
 - Identify mirror sites or web pages
 - Plagiarism
 - Similar news articles

Key steps

- **Shingling:** convert documents (news articles, emails, etc) to sets
- **LSH:** convert large sets to **small signatures**, while preserving the similarity
- Compare the signatures instead of the actual documents

Shingles

- A **k-shingle** (or **k-gram**) is a sequence of **k** characters that appears in a document
- If doc = abcab and $k=3$, then 2-singles: {ab, bc, ca}
- Represent a document by a set of **k-shingles**

Assumption

- Documents that have similar sets of **k**-shingles are similar: same text appears in the two documents; the position of the text does not matter
- What should be the value of **k**?
 - What would large or small **k** mean?

Data model: sets

- Data points are represented as sets (i.e., sets of shingles)
- Similar data points have large intersections in their sets
 - Think of documents and shingles
 - Customers and products
 - Users and movies

Similarity measures for sets

- Now we have a set representation of the data
- Jaccard coefficient
- **A, B** sets (subsets of some, large, universe **U**)

$$\text{sim}(A, B) = \frac{|A \cap B|}{|A \cup B|}$$

Find similar objects using the Jaccard similarity

- Naïve method?
- Problems with the naïve method?
 - There are too many objects
 - Each object consists of too many sets

Speeding up the naïve method

- Represent every object by a signature (summary of the object)
- Examine pairs of signatures rather than pairs of objects
- Find all similar pairs of signatures
- **Check point:** check that objects with similar signatures are actually similar

Still problems

- Comparing large number of signatures with each other may take too much time (although it takes less space)
- The method can produce pairs of objects that might not be similar (false positives). The check point needs to be enforced

Creating signatures

- For object x , signature of x ($\text{sign}(x)$) is much smaller (in space) than x
- For objects x, y it should hold that $\text{sim}(x,y)$ is almost the same as $\text{sim}(\text{sign}(x),\text{sign}(y))$

Intuition behind Jaccard similarity

- Consider two objects: x, y

	x	y
a	1	1
b	1	0
c	0	1
d	0	0

- a : # of rows of form same as a
- $\text{sim}(x, y) = a / (a + b + c)$

A type of signatures -- minhashes

- Randomly **permute** the rows
- **$h(x)$** : first row (in permuted data) in which column **x** has an **1**

	x	y
a	1	1
b	1	0
c	0	1
d	0	0

- Use several (e.g., 100) independent hash functions to design a signature

	x	y
a	0	1
b	0	0
c	1	1
d	1	0

“Surprising” property

- The probability (over all permutations of rows) that $h(x)=h(y)$ is the same as $\text{sim}(x,y)$
- Both of them are $a/(a+b+c)$
- So?
 - **The similarity of signatures is the fraction of the hash functions on which they agree**

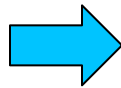
Minhash algorithm

- Pick **k** (e.g., 100) permutations of the rows
- Think of **sign(x)** as a new vector
- Let **sign(x)[i]**: in the **i**-th permutation, the index of the **first row that has 1** for object **x**

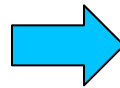
Example of minhash signatures

- Input matrix

	x1	x2	x3	x4
1	1	0	1	0
2	1	0	0	1
3	0	1	0	1
4	0	1	0	1
5	0	1	0	1
6	1	0	1	0
7	1	0	1	0



1
3
7
6
2
5
4



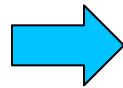
	x1	x2	x3	x4
1	1	0	1	0
3	0	1	0	1
7	1	0	1	0
6	1	0	1	0
2	1	0	0	1
5	0	1	0	1
4	0	1	0	1

1	2	1	2
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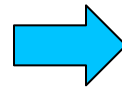
Example of minhash signatures

- Input matrix

	x1	x2	x3	x4
1	1	0	1	0
2	1	0	0	1
3	0	1	0	1
4	0	1	0	1
5	0	1	0	1
6	1	0	1	0
7	1	0	1	0



4
2
1
3
6
7
5



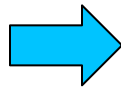
	x1	x2	x3	x4
4	0	1	0	1
2	1	0	0	1
1	1	0	1	0
3	0	1	0	1
6	1	0	1	0
7	1	0	1	0
5	0	1	0	1

2	1	3	1
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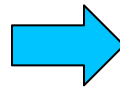
Example of minhash signatures

- Input matrix

	x1	x2	x3	x4
1	1	0	1	0
2	1	0	0	1
3	0	1	0	1
4	0	1	0	1
5	0	1	0	1
6	1	0	1	0
7	1	0	1	0



3
4
7
6
1
2
5



	x1	x2	x3	x4
3	0	1	0	1
4	0	1	0	1
7	1	0	1	0
6	1	0	1	0
1	1	0	1	0
2	1	0	0	1
5	0	1	0	1

3	1	3	1
---	---	---	---

Example of minhash signatures

- Input matrix

	x1	x2	x3	x4
1	1	0	1	0
2	1	0	0	1
3	0	1	0	1
4	0	1	0	1
5	0	1	0	1
6	1	0	1	0
7	1	0	1	0

\approx

	x1	x2	x3	x4
1	2	1	2	
2	1	3	1	
3	1	3	1	

	actua	signs
(x1,x2)	0	0
(x1,x3)	0.75	2/3
(x1,x4)	1/7	0
(x2,x3)	0	0
(x2,x4)	0.75	1
(x3,x4)	0	0

Is it now feasible?

- Assume a billion rows
- Hard to pick a random permutation of 1...billion
- **Even representing a random permutation requires 1 billion entries!!!**
- How about accessing rows in permuted order?
- ☹️

Being more practical

- Approximating row permutations: pick $k=100$ (?) hash functions (h_1, \dots, h_k)

for each row r

for each column c

if c has 1 in row r

for each hash function h_i **do**

if $h_i(r)$ is a smaller value than $M(i,c)$

then

$M(i,c) = h_i(r);$

Being more practical

- Approximating row permutations: pick $k=100$ (?) hash functions (h_1, \dots, h_k)

for each row r

for each column

if c has 1 in row

for each hash

if $h_i(r)$ is a smaller value than $M(i,c)$

then

$$M(i,c) = h_i(r);$$

$M(i,c)$ will become the smallest value of $h_i(r)$ for which column c has 1 in row r ; i.e., $h_i(r)$ gives order of rows for i -th

Example of minhash signatures

- Input matrix

	x1	x2
1	1	0
2	0	1
3	1	1
4	1	0
5	0	1

	x1	x2
1	0	1
2	2	0

$$h(r) = r + 1 \pmod{5}$$

$$g(r) = 2r + 1 \pmod{5}$$