Accelerating Exact k-means Algorithms with Geometric Reasoning

Close reading of a paper

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Accelerating Exact k-means Algorithms with Geometric Reasoning

Dan Pelleg and Andrew Moore School of Computer Science, Carnegie Mellon University Pittsburgh, PA 15213, (dpelleg,awm)@cs.cmu.edu

Abstract

We present new algorithms for the k-means clustering problem. They use the kd-tree data structure to reduce the large number of nearest-neighbor queries issued by the traditional algorithm. Sufficient statistics are stored in the nodes of the kd-tree. Then, an analysis of the geometry of the current cluster centers results in great reduction of the work needed to update the centers. Our algorithms behave exactly as the traditional k-means algorithm. Proofs of correctness are included. The kd-tree can also be used to initialize the k-means starting centers efficiently. Our algorithms can be easily extended to provide fast ways of computing the error of a given cluster assignment, regardless of the method in which those clusters were obtained. We also show how to use them in a setting which allows approximate clustering results, with the benefit of running faster.

We have implemented and tested our algorithms on both real and simulated data. Results show a speedup factor of up to 170 on real astrophysical data, and superiority over the naive algorithm on simulated data in up to 5 dimensions. Our algorithms scale well with respect to the number of points and number of centers, allowing for clustering with tens of thousands of centers.

This paper:

Review: k-means

k-means algorithm

The standard algorithm alternates between two steps until convergence: Given k initial center points,

- 1 Assign each point to its closest center
- 2 Update each center to the centroid of the points assigned to it, where the centroid is the arithmetic mean.

```
k-means analysis
                              dist(\vec{u}, \vec{v}) = \|\vec{u} - \vec{v}\| = \sqrt{(u_1 - v_1)^2 + (u_2 - v_2)^2 + \dots}
def dist(p1: tuple, p2: tuple) -> float:
        Squared distance between two points. """
    return sum((p1[i] - p2[i])*(p1[i] - p2[i]) for i in range(len(p1)))
 I: Number of iterations (hard to know a priori)
 For each iteration:
                                                 For each center:
     For each point:
                                                      Compute dist(point, center)
                                                 Pick minimum distance center
          Find its nearest center -
                                                 = O(KD)
```

= O(INKD), analyze per iteration -> O(NKD)

k-means analysis, cont.

- Updating centers to centroid
- Assume we have point -> nearest center from last slide

For each center, go through its points and average.

Note: points may be weighted (duplicates)

```
E[S] = \frac{1}{\|S\|} \sum_{\vec{p} \in S} \vec{p}
```

```
[(0, 1), (5, 1), (2, 3), (0, 1), (2, 3)] \rightarrow [((0, 1), 2), ((2, 3), 2), ((5, 1), 1)]
```

def mix(pairs: list) -> list:
 """ Takes a list of vector, weight pairs and returns a final centroid. """
 D, denom = len(pairs[0][0]), sum(list(zip(*pairs))[1])
 return tuple(sum(u[d]*w for u, w in pairs)/denom for d in range(D))

= $D(S_1 + S_2 + ... + S_n) = O(DN)$ -> time dominated by assignment

Consider the following...

N = 60,000 pixels

K = 256 (8-bit image)

I = 111 iterations



NKI = 60,000*256*111 > 1.7 billion

 $1.7*10^{9}/1$ MHz = 1700 seconds/60 = 30 minutes!

Can we do better? Yes, kd-trees!

kd-tree, review

- BST insert: compare value against root's value
 - If less, recur on the left subtree
 - \circ If greater, on the right subtree.
- kd-tree: similar, but node holds *point* (vector), not scalar
 - Need some way of comparing
 - Pick an arbitrary *cutting dimension* to compare
 - Cycle through cutting dimensions
- Note: *x*_{cd} = *k* defines a *hyperplane* (high dimensional plane)



Building a kd-tree

- See previous lecture
- O(d² n log n) "pre-sort"
 - O(d n log n) with careful representation

- If all points are known in advance, balanced kd-tree can be built in O(n log n) time
 - Recall: sort the points by x and y coordinates
 - Always split on the median point so each split divides remaining points nearly in half.
 - Time dominated by the initial sorting.

Silly <u>CMU</u>, pre-sort is O(d n log n)!

- O(n log n) "median" if you read too much Introduction to Algorithms
 - Irrespective of dimensionality!
 - High constant factor, however...
- tl;dr fast

mrkd-tree

- "Multi-resolution kd-tree"
- Additional data at each node
 - Bounding box
 - Number of points
 - Centroid
- Can we just annotate a regular kd-tree?



Using the kd-tree

- Can be used to speed up nearest neighbor queries
- $\sim O(\log n + 2^d)$
 - log *n* from tree search
 - \circ 2^d from nodes "near" the target point
- Build kd-tree on centers and use for queries
- Is this really the best approach?

Comparison

kd-tree on centers

- Few points, $K \le N$
 - In practice, K < 10
 - Can maybe save a *K* factor vs log *K*
- Dynamic, centers change every iteration
 - Need to build a new kd-tree
- Easily convert existing algorithm

kd-tree on points

- Many points
 - In practice, *N* > 100,000
 - Can save time as it's based on *N*
 - Static, points are the same
 - Can build once and recycle
 - Requires a bit of adjustment

Clearly, we should build a kd-tree on the *points*, not the centers! How?

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This is where the paper comes in...

Ownership

- Let C be the list of centers, len(C) = K
- Center *c* "owns" a hyper-rectangle *h* (kd-node and descendants)
 - Every point in *h* is closer to *c* than to any other center
 - "Every point" is really "every point"; not just the ones in the kd-tree
 - Obviously, all points assigned to *c*
- If we can show *c* owns *h*, no need to check points in *h*
- How to show ownership?

Point to Hyper-rectangle



```
def distbb(p: tuple, bb: list) -> float:
    """ Squared distance between a point and a bounding box. """
    # three cases, use x if x is in the box, otherwise one of the bounds
    bbp = tuple(box[0] if x < box[0] else (box[1] if x > box[1] else x)
        for x, box in zip(p, bb))
    return dist(p, bbp)
```

- Find the closest point to *p* in *bb*, *q*
- o dist(*p*, *q*)
- 1. *p*[*d*] < *bb*[*d*][0] -> Take *bb*[*d*][0]
- 2. *p*[*d*] > *bb*[*d*][1] -> Take *bb*[*d*][1]
- 3. $bb[d][0] \le p[d] \le bb[d][1] \implies Take p[d]$



Condition #1

- *c* must have the minimum distance to *h*
 - If not true, contradiction!
 - Suppose other center c' with dist(c', h) < dist(c, h)
 - Note that dist(*c*, *h*) *constructive*, finds closest point
 - Therefore there's a point closer to *c*' than *c*
 - Contradicts definition of owner
- Must also be unique
 - Can't have two points with minimum distance
 - Neither owner

```
dists = [distbb(c, t.bb) for c in centers]
```

- i = min(range(len(centers)), key=lambda i: dists[i])
- c, dis = centers[i], dists[i]
- if dists.count(dis) == 1: # c is uniquely the smallest
 # check c is a true owner

Condition #2

- Condition #1 narrows down the owner candidates to either 1 or 0
- If there isn't an owner, we're going to recur on the children
 - Left and right hyper-rectangles
- Let's say we have a candidate c_1
- Is c_1 a true owner?
 - Check against all other centers, call one of them c_2
 - Is every point in *h* closer to c_1 than c_2 ? c_1 "dominates" c_2

```
if dists.count(dis) == 1: # c is uniquely the smallest
    new_centers = centers[:i] + centers[i + 1:]
    owner = True
    for j in range(len(new_centers)):
        if not dominate(c, new_centers[j], t.bb):
            owner = False
            break
```

Domination

- Line of points equidistant to c_1 and $c_2 L$
- If there's a point p on the wrong side of L,
 c₁ doesn't dominate c₂
- Optimization problem:
 - Objective: Be as far in the half-space as possible (not the same as being close to c_{2} !)
 - Constraint: Contained in h
- How to measure objective?
- Look at $v = c_2 c_1$
- *v* dot *p*
- Linear programming!



Linear Programming

- Linear objective and linear constraints
- Yes! Objective v dot p, v is a constant
 - $\circ \quad = v_1 p_1 + v_2 p_2 + \ldots + v_D p_D$, linear function of p
- Constraints of the form
 - $\circ \quad h[d][0] \le p[d] \le h[d][1] \text{ for } d \text{ between } 0 \dots D 1$
- Algorithms for linear programming irrelevant
 - Mathematical ideas relevant

Solving Linear Programs

- Feasible value: satisfies constraints
- Region of feasible values: *simplex*
- Optimal value must be at a corner of the simplex
- *Proof.* Start at arbitrary point and follow the gradient (vector) grad grad until we reach the exterior. From here, move along edges. All edges are orthogonal to the gradient or not.
 (1) If all are, then our function value can't change, current value is optimal. Move along an edge to a corner.
 (2) If all aren't, move along an edge in the right direction. (I do Must eventually run out of edges, at which point (1) applies.

(moving in direction of gradient increases function value)

(I don't know enough math to justify this)

Back to Domination

- Gradient is simply $v = c_2 c_1$, direction towards c_2 's side of the line
- Simplex is the hyper-rectangle *h*
- Optimal value occurs at corners of simplex = corners of *h*
- Can efficiently compute best point *p*
 - Take max or min depending on which is in the direction of $c_2 c_1$

```
def dominate(c1: tuple, c2: tuple, bb: list) -> bool:
    """ Does c1 dominate c2 with respect to bb?
    (every point in bb is closer to c1 than it is to c2). """
    p = tuple(box[0] if x2 < x1 else box[1] for x1, x2, box in zip(c1, c2, bb))
    return dist(p, c1) < dist(p, c2)</pre>
```

Example



Putting it Together

Update(h, C):

- 1. If h is a leaf: base case
 - (a) For each data point in h, find the closest center to it and update that center's counters.
 - (b) Return.
- 2. Compute d(c, h) for all centers c. If there exists one condition #1 center c with shortest distance:

If for all other centers c', c dominates c' with respect to h (so we have established $c = \operatorname{owner}(h)$):

- (a) Update c's counters using the data in h. (b) Return.
- 3. Call Update (h_l, C) . recursion

4. Call Update (h_r, C) .

"For each data point in h" when h is a leaf and leaves of a kd-tree contain exactly one point? 「 (ツ) /

condition #2

This is where the algorithm saves time!

Putting it Together, Code

def __update(self, t: kdNode, centers: list, d: dict) -> None: Update(h, C): 111111 Updates each point to its closest center. """ l. r = t.child1. If h is a leaf: if l is None and r is None: # leaf (a) For each data point in h, find the closest center to it and update that center's counters. d[closest(centers, t.point)].append((t.centroid, t.num)) return (b) Return. dists = [distbb(c, t.bb) for c in centers] 2. Compute d(c, h) for all centers c. If there exists one i = min(range(len(centers)), key=lambda i: dists[i]) c, dis = centers[i], dists[i] center c with shortest distance: if dists.count(dis) == 1: # c is uniquely the smallest new centers = centers[:i] + centers[i + 1:] owner = TrueIf for all other centers c', c dominates c' with refor j in range(len(new centers)): spect to h (so we have established $c = \operatorname{owner}(h)$): if not dominate(c, new centers[j], t.bb): owner = False(a) Update c's counters using the data in h. break *# c dominates all other centers* if owner: (b) Return. d[c].append((t.centroid, t.num)) return # recur on children 3. Call Update (h_l, C) . if l is not None: self. update(l, centers, d, cont) 4. Call Update (h_r, C) . if r is not None: self. update(r, centers, d, cont)

Finishing Touches

```
def update(self, centers: list) -> list:
    """ Wrapper over the recursive helper function __update. """
    d = {c: [] for c in centers}
    self.__update(self, centers, d, cont)
    return [mix(pairs) for c, pairs in d.items()]
```

One Small Adjustment...

- Paper's kd-tree stores points at leaves
- In our kd-tree, intermediate nodes also have points
- Simply add its point with brute force before recursion

since our tree has points on intermediate nodes, handle
d[closest(centers, t.point)].append((t.point, self.weight[t.point]))
recur on children

Blacklisting

- Suppose c_1 is an owner candidate
- c_1 dominates c_2 but isn't an owner overall
- Normally, recur and have to check again
- But c_2 can't be an owner anymore!
 - If we recur on h', every point in h' is in h
 - Thus, c_1 still dominates c_2
 - \circ c_2 can't be an owner
- Saves time in multiple ways!
 - Checking ownership
 - If at a leaf, less centers to check

Blacklisting, code

if dists.count(dis) == 1: # c is uniquely the smallest
 poss, new_centers = [c], centers[:i] + centers[i + 1:]
 for j in range(len(new_centers)):
 if not dominate(c, new_centers[j], t.bb):
 poss.append(new_centers[j])
 if len(poss) == 1: # c dominates all other centers
 d[c].append((t.centroid, t.num))
 return
 centers = poss # blacklist dominated centers

Approximate Pruning

- Speed up the algorithm by cutting off recursion
- Heuristic for when this is "safe"

$$n \cdot \sum_{j=1}^{M} \left(\frac{\operatorname{width}(h)_j}{\operatorname{width}(U)_j} \right)^2 \le d$$

- Few points, small rectangle, early on = less error
- Divide points between current centers

h: hyper-rectangle
n: number of points
U: hyper-rectangle
of all points
M: number of dimensions
i: iteration number
d: parameter,
recommended 0.8

Pruning, Code
$$n \cdot \sum_{j=1}^{M} \left(\frac{\operatorname{width}(h)_j}{\operatorname{width}(U)_j} \right)^2 \leq d^i$$

if APPROX and self.__prune(t):

for c in centers: # split points evenly amongst centers
d[c].append((t.centroid, t.num/len(centers)))

return

since our tree has points on intermediate nodes, handle
d[closest(centers, t.point)].append((t.point, self.weight[t.point]))
recur on children
if l is not None: self.__update(l, centers, d, cont)
if r is not None: self. update(r, centers, d, cont)

Paper is done! How's the results?

Benchmarks, Dataset

hanekawa.png: https://stephen-huan.github.io/assets/images/hanekawa.png

Size: 570 x 517

Pixels: 294690

Distinct pixel count: 59809

Most common pixel: (250, 228, 213) => 19377



Benchmarks, Result

testcases/hanekawa.png, K = 8
k-means++ initialization scheme

	РуРу	Python
kd-tree	5.220s	6.420s
naive	8.379s	17.164s

Benchmarks, Results

```
testcases/hanekawa.png, K = 256, PyPy

naive (run on all pixels): 35m 22s073

naive (distinct pixels): 8m 38s343

kd-tree median (n log n): 3m 52s540

kd-tree pre-sort (dn log n): 3m 48s248

kd-tree approximate (d=0.8): 2m 18s735 ← Slower!

Implementation mistake? ~\_("")_/"

Worse centers -> more iterations

Overhead in calculating prune
```

Extensions

Continue Parameter

- Trying to show *c* is an owner
- If it isn't, should we break?
 - Don't waste time comparing against other centers
 - But continuing could save time if we removed them with blacklisting
- Usually faster to continue than not continue

```
o 4m 0s716 without continue vs 26s802 with continue
for j in range(len(new_centers)):
    if not dominate(c, new_centers[j], t.bb):
        poss.append(new_centers[j])
        if not cont: # whether to continue
        poss += new_centers[j + 1:]
        break
```

kd-tree for Leaf Nodes

- O(0) if point is pruned
- *O*(*K*) if it isn't (leaf or intermediate)
- Can we speed the worst case up?
- Build a kd-tree on the centers again!
- ~ O(log *K*)
- *K* needs to be large in practice; blacklisting removes many centers
 - Average 1-2 centers for K = 256
 - Not *O*(*K*), actually *O*(# of centers)
- Use kd-tree if len(centers) > $F(\log_2 K + 2^d)$, brute force otherwise
 - \circ *F* = 4, see appendix for empirical justification

kd-tree for Leaf Nodes, Code

```
def __close(self, centers: list) -> None:
    """ Set up paramaters for kd-tree closest search. """
    self.t = kdTree(centers)
    # cost roughly 0(log n + 2^d)
    self.cost = F*((len(centers) - 1).bit_length() + (1 << len(centers[0])))</pre>
```

```
def close(self, centers: list, p: tuple) -> tuple:
    """ Finds the closest point with a mix of brute force and kd-tree. """
    if KDTREE and len(centers) >= self.cost:
        return self.t.closest(p)
    return closest(centers, p)
```

"Pseudo"-owner

- Original paper restrictions seem excessive
 - Does *c really* need to be the unique minimum?
 - Does *c really* need to completely dominate?
- No, can relax the conditions
 - Allowed to arbitrarily choose between equidistant centers
 - Maintains same quality and sometimes the same centers
- Reduces number of candidates and improves pruning

```
dists = [distbb(c, t.bb) for c in centers]
dists = [distbb(c, t.bb) for c in centers]
                                                                   i = min(range(len(centers)), key=lambda i: dists[i])
i = min(range(len(centers)), key=lambda i: dists[i])
                                                                   c, dis = centers[i], dists[i]
c, dis = centers[i], dists[i]
                                                                   if PSUEDO or dists.count(dis) == 1:
if dists.count(dis) == 1: # c is uniquely the smallest
                                                                        . . .
                                                                def dominate(c1: tuple, c2: tuple, bb: list) -> bool:
def dominate(c1: tuple, c2: tuple, bb: list) -> bool:
                                                                    """ Does c1 dominate c2 with respect to bb?
    """ Does c1 dominate c2 with respect to bb?
    (every point in bb is closer to c1 than it is to c2). """
                                                                    (every point in bb is closer to c1 than it is to c2). """
                                                                    p = tuple(box[0] if x2 < x1 else box[1]
    p = tuple(box[0] \text{ if } x2 < x1 \text{ else } box[1]
              for x1, x2, box in zip(c1, c2, bb))
                                                                              for x1, x2, box in zip(c1, c2, bb))
    d1, d2 = dist(p, c1), dist(p, c2)
                                                                    d1, d2 = dist(p, c1), dist(p, c2)
    return d1 < d2
                                                                    return d1 < d2 or (PSUEDO and d1 == d2)
```

Benchmarks, Results



k-means++

k-means++, Review

- Initialization strategy
- For more information, see the previous lecture

k-means++ algorithm

- **1** Pick the first center point at random
- 2 From there, pick the next center point by sampling the probability distribution where a point \vec{p} is picked with weighting dist $(\vec{p}, \vec{c})^2$, where \vec{c} is the closest center
- **3** Repeat until k centers are picked
- 4 Run k-means as usual

Speedup

- We can apply the kd-tree algorithm
- Simply change the counters
- Leaf/internal node:

```
def __get_children(self, t: kdNode, children: list) -> list:
    """ Recursively gets the children of a kd-tree. """
    l, r = t.child
    children.append(t.point)
    if l is not None: self.__get_children(l, children)
    if r is not None: self.__get_children(r, children)
    return children
```

```
def get_children(self, t: kdNode) -> list:
    """ Wrapper function. """
    return self.__get_children(t, [])
```

```
ids[t.point] = closest(centers, t.point)
```

• Pruned point:

```
for point in self.get_children(t):
    ids[point] = c
return
```

Speedup, Analysis

- Might not seem like we're doing much
- Need to touch every point still
- However, save O(K) distance checking if pruned
 - Save O(average # of centers) after blacklisting for leaves
- Could be up to K times faster
- Can we do better?

Caching IDs

- Need to iterate over each point when pruned
 - Original algorithm can simply skip over
- Centers aren't changing, only added
- Cache IDs if the new center doesn't change ownership
 - If kd-node points to *c* and *c* owns *h*, then *c* must have owned *h* in the past
 - If *c* owned *h* and *c* dominates the new center, all points already point to *c*
 - No need to do anything

```
if len(poss) == 1:  # c dominates all other centers
    if ids.get(t.point, None) != c:  # changed
        for point in self.get_children(t): # slow but necessary
        ids[point] = c
        return
```

Caching IDs, Analysis

- Expected time savings?
- Assume points are split evenly among centers
 - $K = 1 \rightarrow$ need to update all points
 - $K = 2 \rightarrow$ need to update half of the points
 - etc.
- $N + \frac{1}{2}N + \frac{1}{3}N + \dots + \frac{1}{K}N = N(1 + \frac{1}{2} + \frac{1}{3} + \dots + \frac{1}{K})$
- = $N H_{K} = O(N \log K)$
- Very optimistic
- In practice:
 - 108,248 for K = 8 vs *NK* = 480,000
 - 4,794,349 for K = 256 vs *NK* = 15,360,000

Probability Distribution

- Still maintaining a list of distances
- Need to sample from this PMF
- = O(N), O(NK) over K centers
- Can we speed this up?

BIT CMF

- Maintain the *cumulative mass function* (CMF) instead
- CMF is just the prefix sum of the PMF
- Only update when a point changes
- Can efficiently update with <u>segment trees</u> or <u>binary indexed trees</u> (BITs)
 - My BIT and segtree implementation <u>here</u>
- Sample by binary searching on the CMF
 - Why does this work? See: <u>my analysis of Mudae</u> (page 9) and the <u>previous lecture</u>

BIT CMF, Code

def __update_point(self, ids: dict, point: tuple, center: tuple) -> None:
 """ Updates the probability distribution of a point. """
 ids[point] = center
 i, w = self.point_id[point], self.weight[point]
 self.cmf.update(i, w*dist(point, center) - self.cmf.range(i, i))

```
def sample(self) -> tuple:
    """ Samples a point according to k-means++. """
    p = random.random()
    # sum of all values
    denom = self.cmf.query(len(self.points) - 1)
    l, r = 0, len(self.points)
    while l < r:
        m = (l + r)>>1
        if self.cmf.query(m)/denom <= p:
            l, r = m + 1, r
        else:
            l, r = l, m
        return self.points[l]
```

BIT CMF, Analysis

- Sample runs in $O(\log^2 N)$
 - \circ log *N* queries from the binary search, each query is log *N*
 - = $O(K \log^2 N)$ over K centers
- The total amount of work for updating IDs is harder
- Optimistically, O(N log K) total updates
 - $\circ = O(N \log N \log K) \text{ overall}$
- In practice, closer to $O(NK) \rightarrow O(N \log N K)$
- Naive is $O(NK^2)$
 - *K* iterations of finding the closest center for *N* points

Benchmarks

testcases/hanekawa.png, K = 256, PyPy
k-means++ naive: 311.600s, 111 iterations
k-means++ kd-tree: 122.594s, --k-means++ geometric reasoning: 75.269s, --k-means++ cache ids: 53.181s, --k-means++ BIT CMF: 44.143s, --random.sample(): 0.001s, 108

k-means++, Judgement

- *k*-means++ has strong theoretical properties
- On "real-world" datasets, irrelevant
 - *K* iterations over *N* is just way too much!
 - Can't parallelize, each iteration depends on previous
- Need some way of speeding up...

Conclusion

- Build a kd-tree on the points, not the centers
- Works well in practice and in theory
- Many avenues for improvement



256 color image generated in less than 30 seconds

References

- Accelerating Exact k-means Algorithms with Geometric Reasoning
- <u>k-means, kd-Trees, and Median of Medians</u> (CV club lecture)
- CMU <u>kd-tree</u>, <u>kd-tree continued</u>

Appendix

Convergence

- How do we test if *k*-means has converged?
- Centers hasn't changed
- Sort new center list, check if equal to previous
- Not necessarily $O(D K \log K)$ (although this is a fairly trivial cost)
- Some sorting algorithms proportional to displacement, e.g. bubble sort
- Can run in nearly linear time if the centers don't change that much

Determining F

• Naive:

- \circ T₁(x) = mx + b
- \circ 0 at x = 0, 6 at x = 100
- o m = 6/100, b = 0

• kd-tree:

- $T_2(x) = a \ln x + b$
- 2 at 20, 3 at 100
- a = 1/(ln(100) ln(20)), b = 2 a ln 20
- K = 256, $T_2(K) = 3.58$, $T_1^{-1}(K) = 59.7$
 - $\circ \quad \log_2 K + 2^d = 8 + 8 = 16$
 - Therefore $F(\log_2 K + 2^d) = 16F \approx 64$
 - *F* = 4

