

K-Means Algorithm

Implementation of sequential and parallel version using OpenMp

Marco Trambusti - 7135350

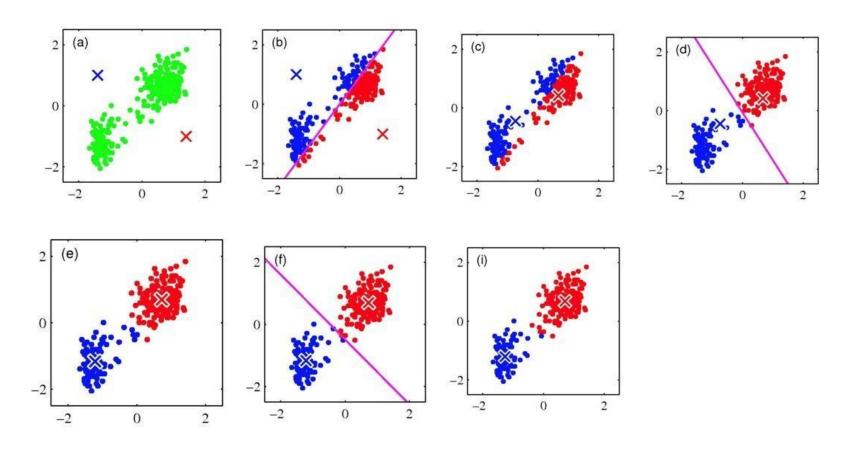
Introduction

Purpose of the Assignement

- Implement the K-Means algorithm in C++ in both sequential and parallelized versions using OpenMP.
- Dataset: 300147 values related to age and total spending of supermarket customers.
- Performance comparison on a machine with an Intel Core i7-6700K processor (4 cores).

K-Means Algorithm

- **1. Initialization**: Random selection of K initial cluster centers.
- 2. Cluster Assignment: Each data point is assigned to the nearest centroid.
- **3. Centroid Update:** Centroids are recalculated as the mean of the points assigned to each cluster.
- **4. Iteration:** The assignment and update steps are repeated until the centroids no longer change significantly or a maximum number of iterations is reached.



Esempio di K-Means



Algorithm Parameters

- Number of clusters (K)
- Initialization of centroids
- Choice of distance metric (squared Euclidean distance)
- Maximum number of iterations

Sequential Implementation

It was chosen C++ for the sequential implementation and OpenMP for the parallel implementation.

In this program, 2D points points are readed from a csv file and clusters are initialized randomly;

The points in the dataset are represented using a struct **Point**, composed by:

- x, y coordinates (double)
- cluster (int);
- minDist (double)
- distance(Point p) (method)

Initially, the point does not belong to any cluster, and was arbitrarily set to -1. As a result, minDist was set to the maximum possible value.



Sequential Implementation

Kmeans method: Cluster assigment part

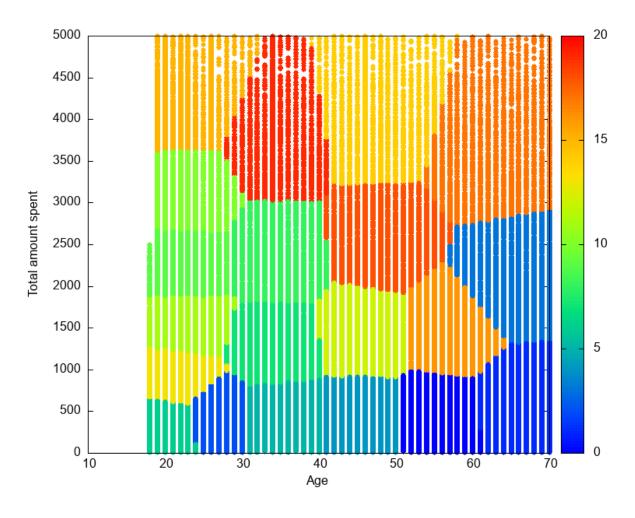
```
1 void kmeans(std::vector<Point>& points, std::vector<Point> centroids,
      int k, int epochs) {
      for (int epoch = 0; epoch < epochs; ++epoch) {</pre>
           int nPoints[k] = {0};
           double sumX[k] = \{0.0\};
5
           double sumY[k] = \{0.0\};
           for (auto& point : points) {
               for (int i = 0; i < k; ++i) {
                   double dist = point.distance(centroids[i]);
9
                   if (dist < point.minDist) {</pre>
10
                        point.minDist = dist;
11
                        point.cluster = i;
12
13
14
               //append data to centroids
15
               int clusterId = point.cluster;
16
               nPoints[clusterId]++;
17
               sumX[clusterId] += point.x;
               sumY[clusterId] += point.y;
19
               //reset distance
20
               point.minDist = std::numeric_limits<double>::max();
23
```



Sequential Implementation

Kmeans method: Centroid Update Code

```
for (int i = 0; i < k; ++i) {
   if (nPoints[i] > 0) {
      double newX = sumX[i] / nPoints[i];
      double newY = sumY[i] / nPoints[i];
      centroids[i].x = newX;
      centroids[i].y = newY;
}
```



Example of results plotted using GnuPlotted

Parallel Implementation

Cluster Assignment

- Cluster assignment step is independent for each data point,
 - ideal candidate for parallelization.
- Using OpenMP, it is possible to distribute the calculation parallelizing the outer loop.
- Updating the sum of coordinates and the number of points relative to the cluster, adds complexity due to data sharing,
 - Optimized using reductions



Parallel Implementation

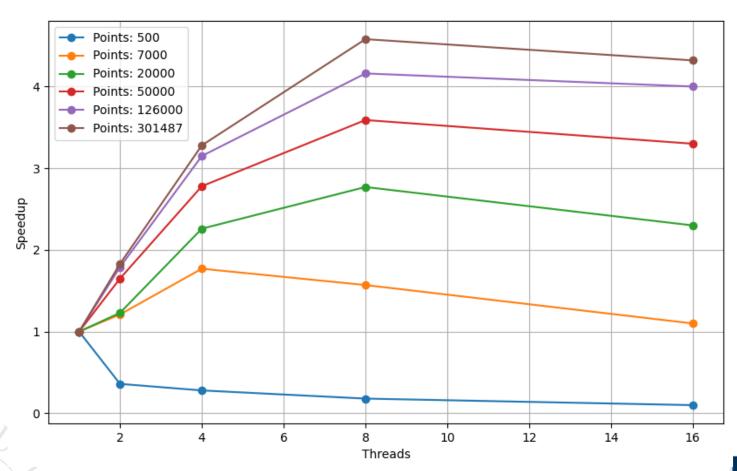
Centroids Update

```
1 // Update centroids
2 #pragma omp for
3 for (int i = 0; i < k; ++i) {
4    if (nPoints[i] > 0) {
5         ...
6    }
7 }
```



Results and Performance Evaluation

Speedup: Ratio of sequential to parallel execution time.



Conclusions

- 1. Parallelizing the K-Means algorithm proved effective, especially with larger datasets.
- 2. OpenMP simplified implementation and performance optimization.
- 3. The larger the number of points, the more effectively threads can be utilized up to a certain point.