

## a2/kmeans/omp\_reduction\_kmeans.c

```
1  #include <stdio.h>
2  #include <stdlib.h>
3  #include "kmeans.h"
4  /*
5   * TODO: include openmp header file
6   */
7  #include <omp.h>
8
9  // square of Euclid distance between two multi-dimensional points
10 inline static double euclid_dist_2(int numdims, /* no. dimensions */
11                                     double *coord1, /* [numdims] */
12                                     double *coord2) /* [numdims] */
13 {
14     int i;
15     double ans = 0.0;
16
17     for (i = 0; i < numdims; i++)
18         ans += (coord1[i] - coord2[i]) * (coord1[i] - coord2[i]);
19
20     return ans;
21 }
22
23 inline static int find_nearest_cluster(int numClusters, /* no. clusters */
24                                       int numCoords, /* no. coordinates */
25                                       double *object, /* [numCoords] */
26                                       double *clusters) /* [numClusters][numCoords] */
27 {
28     int index, i;
29     double dist, min_dist;
30
31     // find the cluster id that has min distance to object
32     index = 0;
33     min_dist = euclid_dist_2(numCoords, object, clusters);
34
35     for (i = 1; i < numClusters; i++)
36     {
37         dist = euclid_dist_2(numCoords, object, &clusters[i * numCoords]);
38         // no need square root
39         if (dist < min_dist)
40         { // find the min and its array index
41             min_dist = dist;
42             index = i;
43         }
44     }
45     return index;
46 }
47
48 void kmeans(double *objects, /* in: [numObjs][numCoords] */
49             int numCoords, /* no. coordinates */
50             int numObjs, /* no. objects */
51             int numClusters, /* no. clusters */

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52     double threshold, /* minimum fraction of objects that change membership */
53     long loop_threshold, /* maximum number of iterations */
54     int *membership, /* out: [numObjs] */
55     double *clusters) /* out: [numClusters][numCoords] */
56 {
57     int i, j, k;
58     int index, loop = 0;
59     double timing = 0;
60
61     double delta; // fraction of objects whose clusters change in each loop
62     int *newClusterSize; // [numClusters]: no. objects assigned in each new cluster
63     double *newClusters; // [numClusters][numCoords]
64     int nthreads; // no. threads
65
66     nthreads = omp_get_max_threads();
67     printf("OpenMP Kmeans - Reduction\t(number of threads: %d)\n", nthreads);
68
69     // initialize membership
70     for (i = 0; i < numObjs; i++)
71         membership[i] = -1;
72
73     // initialize newClusterSize and newClusters to all 0
74     newClusterSize = (typeof(newClusterSize))calloc(numClusters, sizeof(*newClusterSize));
75     newClusters = (typeof(newClusters))calloc(numClusters * numCoords,
76     sizeof(*newClusters));
77
78     // Each thread calculates new centers using a private space. After that, thread 0 does
79     // an array reduction on them.
80     int *local_newClusterSize[nthreads]; // [nthreads][numClusters]
81     double *local_newClusters[nthreads]; // [nthreads][numClusters][numCoords]
82
83     /*
84     * Hint for false-sharing
85     * This is noticed when numCoords is low (and neighboring local_newClusters exist
86     close to each other).
87     * Allocate local cluster data with a "first-touch" policy.
88     */
89     // Initialize local (per-thread) arrays (and later collect result on global arrays)
90     for (k = 0; k < nthreads; k++)
91     {
92         local_newClusterSize[k] = (typeof(*local_newClusterSize))calloc(numClusters,
93         sizeof(**local_newClusterSize));
94         local_newClusters[k] = (typeof(*local_newClusters))calloc(numClusters * numCoords,
95         sizeof(**local_newClusters));
96     }
97
98     timing = wtime();
99     do
100     {
101         // before each loop, set cluster data to 0
102         for (i = 0; i < numClusters; i++)
103         {
104             for (j = 0; j < numCoords; j++)
105                 newClusters[i * numCoords + j] = 0.0;

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101         newClusterSize[i] = 0;
102     }
103
104     // reset delta before each iteration; it will be updated via reduction in the
parallel region
105     delta = 0.0;
106
107     /*
108      * TODO: Initiliaze local cluster data to zero (separate for each thread)
109      *
110      * We now use an OpenMP parallel region where:
111      * - Each thread zeroes its own local_newClusterSize/local_newClusters.
112      * - The object loop is distributed with 'omp for' and 'reduction(+ : delta)'.
113      * - A single thread reduces the per-thread local arrays into the shared arrays.
114      */
115 #pragma omp parallel private(i, j, k, index)
116 {
117     int tid = omp_get_thread_num();
118     int T = omp_get_num_threads(); // actual number of threads in this team
119
120     /* per-thread zeroing (first-touch initialization of local cluster data) */
121     for (i = 0; i < numClusters; i++)
122         local_newClusterSize[tid][i] = 0;
123     for (i = 0; i < numClusters * numCoords; i++)
124         local_newClusters[tid][i] = 0.0;
125
126     // Distribute objects across threads and compute per-thread contributions.
127     // delta is accumulated using a reduction to avoid atomics on a shared
variable.
128 #pragma omp for reduction(+ : delta)
129     for (i = 0; i < numObjs; i++)
130     {
131         // find the array index of nearest cluster center
132         index = find_nearest_cluster(numClusters, numCoords,
&objects[i * numCoords], clusters);
133
134         // if membership changes, increase delta by 1
135         if (membership[i] != index)
136             delta += 1.0;
137
138         // assign the membership to object i
139         membership[i] = index;
140
141         // update new cluster centers : sum of all objects located within (average
will be performed later)
142         /*
143          * TODO: Collect cluster data in local arrays (local to each thread)
144          * Replace global arrays with local per-thread
145          */
146         local_newClusterSize[tid][index]++;
147         for (j = 0; j < numCoords; j++)
148             local_newClusters[tid][index * numCoords + j] += objects[i * numCoords
+ j];
149     }
150 }

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151
152     /*
153     * TODO: Reduction of cluster data from local arrays to shared.
154     *     This operation will be performed by one thread
155     *
156     * Here we use 'omp single' so that exactly one thread accumulates
157     * all per-thread local arrays into the shared newClusterSize/newClusters.
158     */
159 #pragma omp single
160 {
161     for (k = 0; k < T; k++)    // only sum over the threads actually in this
team
162     {
163         int *srcS = local_newClusterSize[k];
164         double *srcC = local_newClusters[k];
165         if (!srcS || !srcC)
166             continue;
167         for (i = 0; i < numClusters; i++)
168         {
169             newClusterSize[i] += srcS[i];
170             for (j = 0; j < numCoords; j++)
171                 newClusters[i * numCoords + j] += srcC[i * numCoords + j];
172         }
173     }
174     } /* implicit barrier after single */
175 }    /* end parallel region */
176
177 // average the sum and replace old cluster centers with newClusters
178 for (i = 0; i < numClusters; i++)
179 {
180     if (newClusterSize[i] > 0)
181     {
182         for (j = 0; j < numCoords; j++)
183         {
184             clusters[i * numCoords + j] = newClusters[i * numCoords + j] /
newClusterSize[i];
185         }
186     }
187 }
188
189 // Get fraction of objects whose membership changed during this loop. This is used
as a convergence criterion.
190 delta /= numObjs;
191
192 loop++;
193 printf("\r\tcompleted loop %d", loop);
194 fflush(stdout);
195 } while (delta > threshold && loop < loop_threshold);
196 timing = wtime() - timing;
197 printf("\n nloops = %3d (total = %7.4fs) (per loop = %7.4fs)\n", loop, timing, timing
/ loop);
198
199 for (k = 0; k < nthreads; k++)
200 {

```

```
201         free(local_newClusterSize[k]);
202         free(local_newClusters[k]);
203     }
204     free(newClusters);
205     free(newClusterSize);
206 }
207
208
```