

# Implementation of Localized Algorithm for Self-Organized Wireless Sensor Network

Arnab Kundu and Debadrita Roy

**Abstract:** Communication between two nodes is a big issue in now a days. To achieve that, wireless network plays a big role. With limited source of energy, memory & computation power wireless sensor networks can be composed by mass number of sensor nodes. Where the sensors' lifetime depend on energies. Like autonomous system in LAN, cluster can be defined for wireless sensor network, where cluster head plays the prime role for the energy conservation. So, the optimization of cluster head within cluster along with number of nodes is a big research issue. Here we have analyzed the advanced optimization algorithm for sensor network clustering theoretically. We have tried to test the proposed method as a clustering algorithm and compared it with other recent sensor network clustering algorithm named Algorithm for Cluster Establishment (ACE)

**Keywords:** ACE; Data sets; Localized algorithm; Migration; Spawning threshold function

## I. INTRODUCTION

The unique characteristics of wireless sensor networks and the nature of their applications motivate the development of new algorithmic solutions to provide services such as information processing, messages routing, fault-tolerance, localization, naming and addressing. Generally algorithms have been classified as either centralized or distributed. Centralized algorithms execute on a single node, which is impractical for large-scale highly constrained sensor networks. This approach would require all nodes to send their respective information (e.g., sensed data, or energy level) to a specified location (e.g., a sink node) and wait until this node executes the algorithm and sends back its output[1,6]. This approach is unsuitable for sensor networks because of its cost in terms of energy and delay. In addition, it suffers low fault-tolerance and low scalability and creates bottlenecks. In the case of distributed algorithms, different nodes participate in the computation which can reduce the execution time. To address this problem, a new class of algorithms called localized algorithms has been proposed for sensor networks. Localized algorithms are a special category of distributed algorithms that are designed specifically to enable efficient and reliable collaboration between sensor nodes by taking into account their specific constraints. For instance, a localized algorithm might limit the collaboration and information exchange to close neighbors of a node to save energy.[2,4,9,10].

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## II. ALGORITHM FOR CLUSTER ESTABLISHMENT (ACE)

Unlike other distributed clustering schemes, ACE employs an emergent algorithm. Emergent algorithms much like artificial neural networks evolve to optimal solution through a mix of local optimization steps. The main idea of ACE is to allow a node to assess its potential as a CH before becoming one and stepping down if it is not the best CH at the moment. The algorithm works in iterations that do not have to be synchronized at the individual nodes. Spawning new clusters and migration of existing ones are the two functional components of ACE. A node spawns of new cluster when it decides to become a CH. It broadcasts an invitation message to recruit its neighbors. Upon getting the invitation, a neighboring sensor joins the new cluster and becomes a follower of the new CH. At any moment, a node can be a follower of more than one cluster. However, the node can be a loyal follower, i.e. a member, of only a single cluster.[3,5]

### A. ACE Algorithm

```

procedure SCALE_ONE_ITERATION()
  if myTime > 3 × EXPECTED_ITERATION_LENGTH then
    if myState = CLUSTER-HEAD then
      return DONE
    else if myState = CLUSTERED then
      wait for my cluster-heads to terminate, then pick one as my cluster-head
      return DONE
    else if myState = UNCLUSTERED then
      pick a random clustered node to act as my proxy after it terminates
      wait for it to terminate, then return DONE
    end if
    else if myState = UNCLUSTERED
      and numLoyalFollowers() ≥ fmin(myTime) then
        myClusterID ← generate_New_Random_ID()
        locally_broadcast(RECRUIT, myID, myClusterID)
      else if myState = CLUSTER-HEAD then
        bestLeader ← myID
        bestFollowerCount ← numLoyalFollowers
        for all n where n is a potential new cluster-head do
          followerCount = Poll_For_Num_Loyal_Followers(n, myClusterID)
          if followerCount > bestFollowerCount then
            bestLeader ← n
            bestFollowerCount ← followerCount
          end if
        end for
        if bestLeader is not myID then
          send(bestLeader, PROMOTE, myClusterID)
          wait for bestLeader to broadcast its RECRUIT message
          locally_broadcast(ABDICATE, myID, myClusterID)
        end if
      end if
    end if
  end procedure

```

**B. Migration**

Migration is a process in which the best candidate for being CH is selected. Each CH periodically checks the ability of its neighbors for being a CH and decides to step down if one of these neighbors has more followers than it does. A node that has the largest number of followers and the least overlap with existing clusters will be considered as the best candidate for CH.

**C. Parameter Selection**

In the ACE protocol, an unclustered node will spawn a new cluster by declaring itself a cluster head whenever it finds that it can gain at least  $f_{min}$  loyal followers if it were to become a cluster head. The function  $f_{min}$  is called the *spawning threshold function* and is dependent on the time that has passed since the protocol was initiated for that node. In general,  $f_{min}$  should decrease as the algorithm proceeds. This causes fewer clusters to form near the beginning of the algorithm. Here, we used an exponentially decreasing function for  $f_{min}$ :

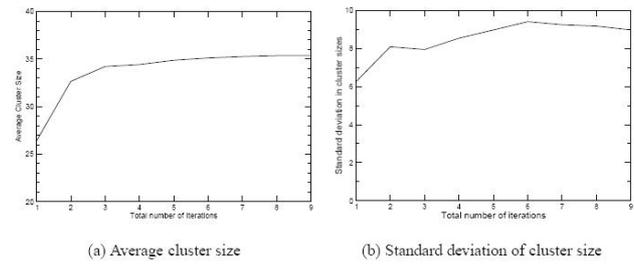
$$f_{min} = (e^{-k_1 \frac{t}{cI}} - k_2)d$$

In this formula,  $t$  is the time passed since the protocol began and  $cI$  is the duration of the protocol as described earlier.  $d$  is the estimated average degree (number of neighbours) of a node in the network.  $k_1$  and  $k_2$  are chosen constants that determine the shape of the exponential graph. In practice, it was empirically found that  $k_1 = 2:3$  and  $k_2 = 0:1$  have produced good results. In this case,  $f_{min}$  starts at  $0:9d$  at the beginning of the protocol and reduces to 0 by the final iteration. This ensures that any node left unclustered at the end of the protocol will declare itself a cluster head.[7]

An alternative parameter setting is  $k_1 = 2:3$  as before, but setting  $k_2 = 0$ . In this case the function starts near  $d$  when the protocol commences and reduces to  $0:1d$  at the end of the protocol. Since  $0:1d > 1$  if  $d > 10$ , it is possible that there will be a small number of nodes that will not be within one hop radius of any cluster-head at the end of the algorithm. It can be observed in simulation that the number of nodes not within one-hop radius of a cluster-head is, on average, less than 4% of the total number of nodes in low node deployment densities, and around 2% for moderate to high node deployment densities (20 or more neighbors per node). These nodes that are not within one hop radius of any cluster-head can simply pick a clustered neighbor to act as their bridge to the cluster head, thus becoming *two-hop followers* (because they take 2 hops to communicate with the cluster-head, instead of the usual 1 hop).

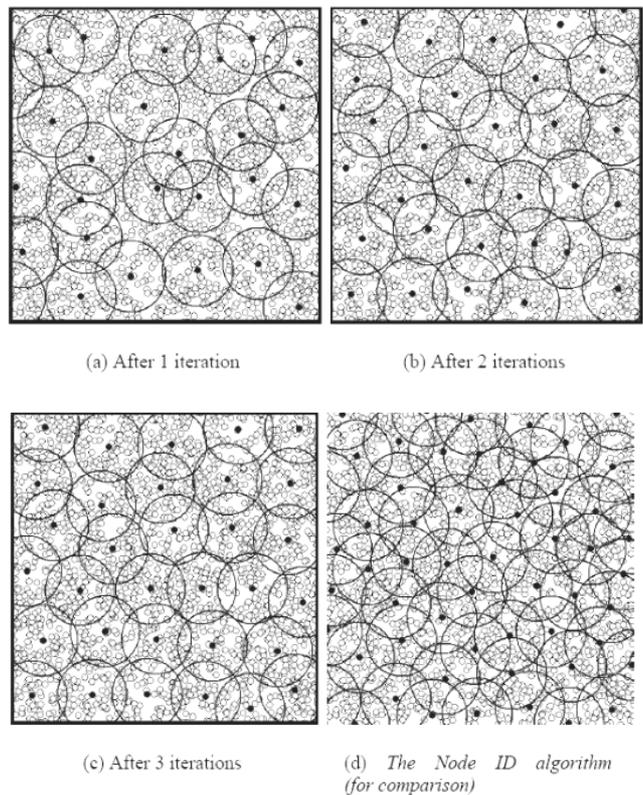
It remains to determine  $c$ , the number of iterations the algorithm should execute. Figure-1. reflects how the performance of ACE changes as it is given a longer number of iterations to operate. ACE was simulated in a 2D area with a uniform random distribution with an average deployment density  $d$  of 50 nodes per circle of one communication radius. Results for the simulation with  $k_1 = 2:3$  and  $k_2 = 0:1$  are shown (results for  $k_2 = 0$  have similar characteristics). It can be noted that increasing the number of iterations above 3

yielded only very slight improvements in average cluster size. In this simulation the total number of iterations did not significantly affect the standard deviation in cluster sizes, which was between 6 and 10 for all iterations > 1. [3,5]



**Fig.1. Performance of ACE at various maximum iterations,  $d=50, k_1=2.3, k_2=0.1$**

Figure-2. illustrates the ACE algorithm operating in simulation (with  $k_1 = 2:3$  and  $k_2 = 0$ ). The little circles represent nodes. Cluster-heads are highlighted in black, and their range is indicated with a large black circle (nodes within the circle are in that cluster). The clusters migrate away from each other in successive iterations to produce a highly efficient cover of the area. Clusters tend to center over areas where nodes are dense. The clusters overlap minimally, and when they do overlap, they tend to overlap in areas where nodes are sparse. Figure-2d. provides a qualitative visual comparison of the Node ID algorithm with ACE. It can be observed that ACE provides a packing with significantly less cluster overlap than Node ID.[8]



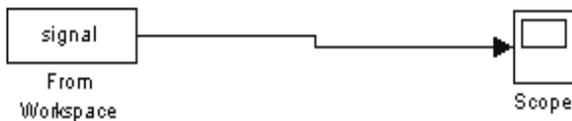
**Fig.2. The ACE algorithm(with  $k_1=2.3$  &  $k_2=0$ )**

### III. SIMULATED OUTPUTS & DISCUSSION

To implement the above algorithm we have used the Application of Program named “MATLAB – 7” as front end application and “Microsoft Excel – 2000” as a Backend application or database. In this section we have shown the outputs of various type of Sensor nodes for variable Lat-Long. Then we have evaluated the function *fmin* in the following program.

#### A. Sensor Node Outputs

Data Name : elcentro\_EW.dat  
 signal=load('elcentro\_EW.dat');  
 From Simulink Library Browser :



Output : 1.

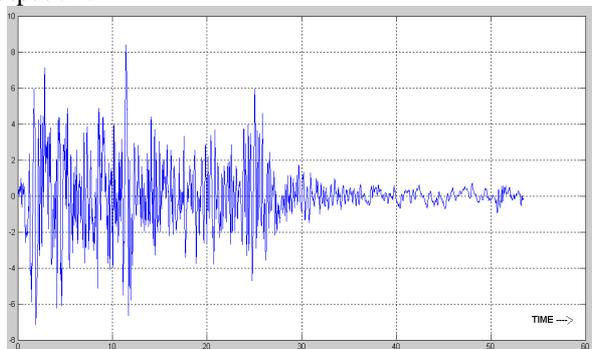
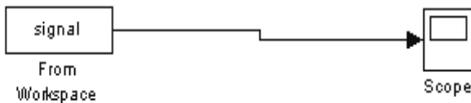


Fig.3. Vibration sensor output

Data Name : elcentro\_NS.dat  
 signal=load('elcentro\_NS.dat');  
 From Simulink Library Browser :



Output : 2.

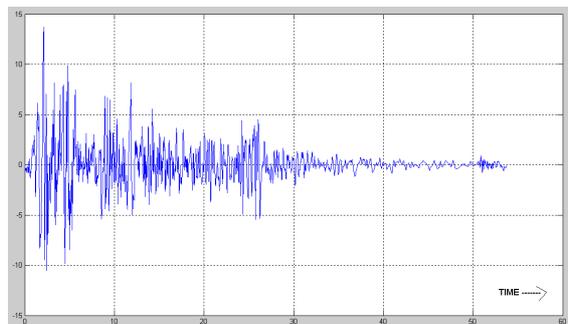


Fig.4. Vibration sensor output

Data Name: elcentro\_EW.dat  
 signal=load(' elcentro\_UP.dat');  
 From Simulink Library Browser :



Output :3.

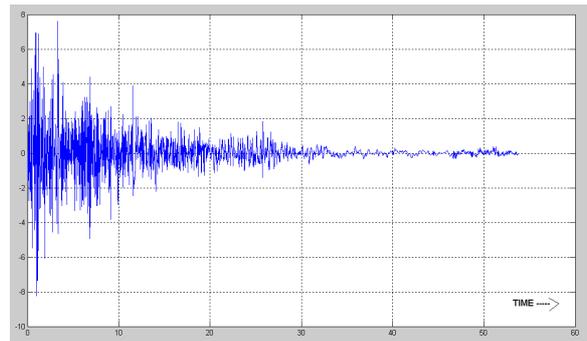


Fig.5. Vibration sensor output

#### B. Calculation of *fmin*

Considering,  $fmin = d$

- $k1 = a$
- $t/cI = b$
- $k2 = c$
- $d = h$

Output : 1.

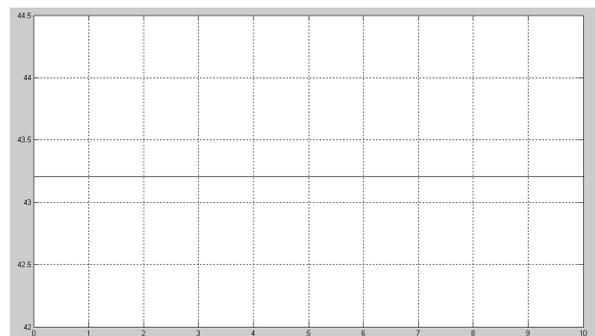
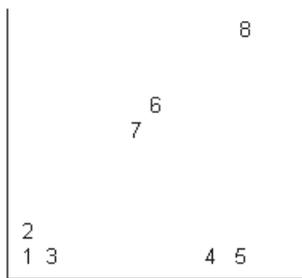


Fig.6. Performance of ACE at various maximum iterations,  $d=50, k1=2.3, k2=0.1$

Each object is represented by an ordered n-tuple of its attributes. Objects with similar characteristics must have individual attributes which are close in value. Since distance is a measure of "closeness", it can be used in clustering. For each cluster, there will be a single point which is designated as a hub. The placement of the hub within the cluster is determined by the algorithm. We will start with two-dimensional objects since they are the easiest to visualize. Suppose we have 8 points in the configuration below. Draw circles around the clusters that you see visually.

**TABLE I**

POINT	X	Y
1	1	1
2	1	2
3	2	1
4	8	1
5	9	1
6	6	6
7	5	5
8	9	9



Let's see what a computer model of clustering would give us.

This algorithm assumes a minimum of 2 clusters and a maximum of p clusters where p is the number of objects. The algorithm is as follows:

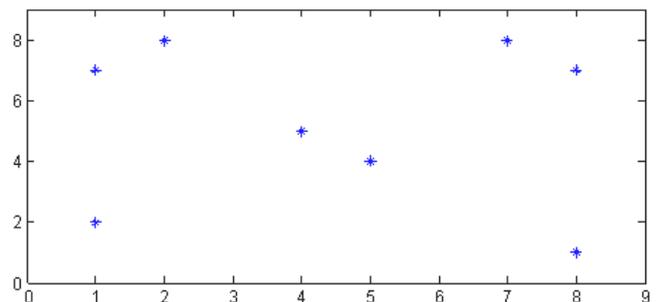
1. From the set of objects (data points), choose any one to be the first hub.
2. Find the point which is farthest from this hub, and make it the second hub.
3. For each remaining object, find the distance from the object to each hub and assign the object to the hub to which it is closer.
4. To determine whether it is necessary to find a third hub, do the following.
  - a. Find the distance between the two hubs and divide by two. Call this value R.
  - b. Calculate the distance from each object to its hub. If any distance is greater than R, it is necessary to find a third hub.
5. To find the third hub, find the object which is farthest from its hub (1 or 2). So, if Point 8 belongs to Hub 1, look at the distance from Point 8 to Hub 1, but not Point 8 to Hub 2.

6. Then, calculate the distance from each point to the new hub to see if that distance is less than the point's distance to its current hub. If so, reassign the point to Hub 3.
7. To determine the stop value, R, find the average distance between the hubs and divide by two. Since there are three hubs, there will be three distances to average together (H1 - H2, H1 - H3, and H2 - H3).
8. Once again, see if any distance from a point to its hub exceeds R.
9. Continue this process until all points are within R units of their hub or until all points are themselves hubs.

**TABLE III - data.dat (Data File)**

	1	2
1	4	5
2	1	7
3	1	2
4	2	8
5	8	1
6	7	8
7	8	7
8	5	4

Output: 1.



**Fig.7. Graphical plotting of datasheets (Top View)**

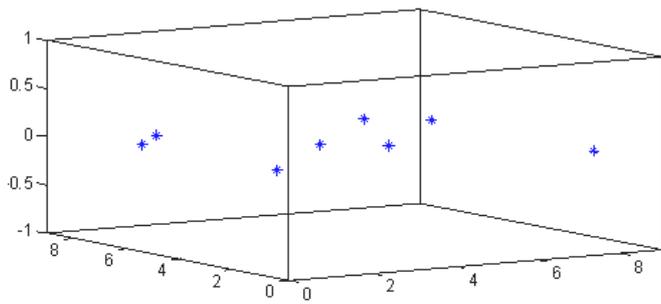


Fig.8. Graphical plotting of datasheets (3-D View)

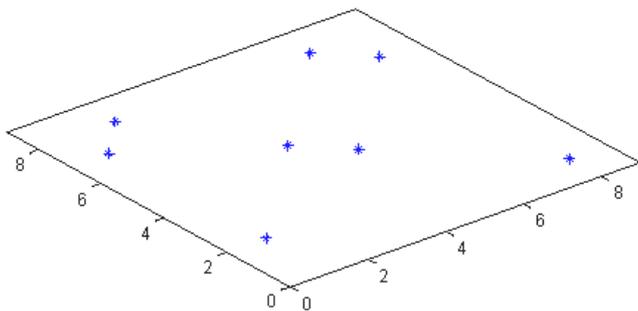


Fig.9. Graphical plotting of datasheets(Side top View)

Output: 2.

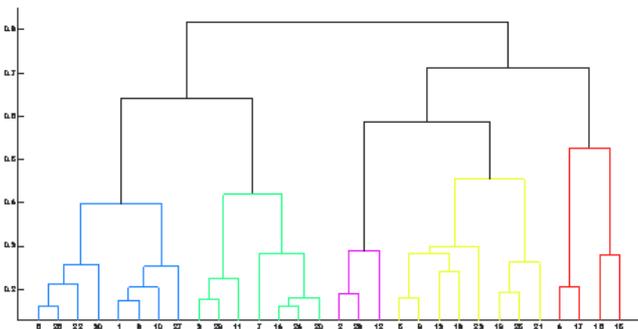


Fig.10. Clustering of Data Sets

Output: 3.

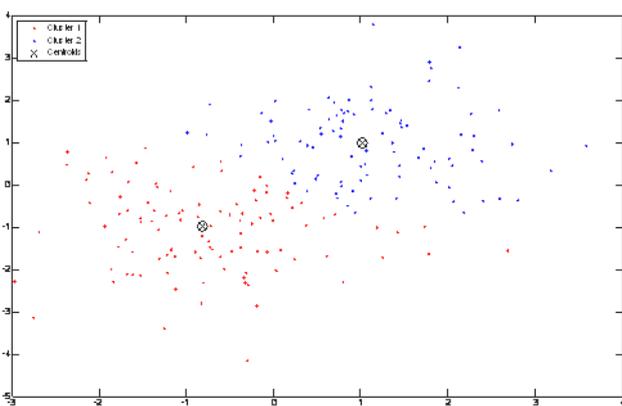


Fig.11. Clustering process

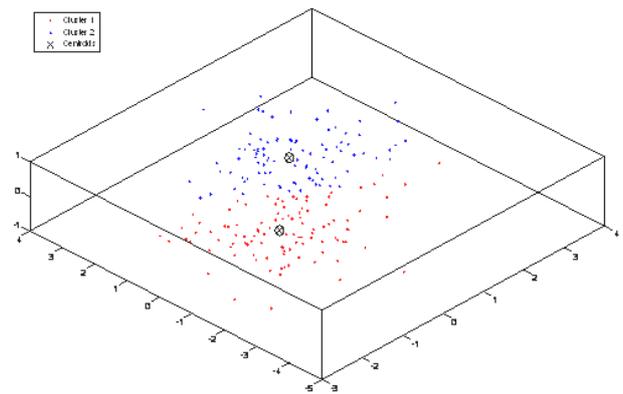


Fig.12. Clustered nodes

IV. CONCLUSION

Through the above theoretical analysis and simulated outputs we are trying to solve the network link between two Nodes of a cluster as well as measure the Stability Criteria of those (A.C.E., S.O.S., Optimization Algorithm) Clustering Algorithms. Clustering Algorithms are very important at present. There are huge scope of work. The basic concepts needed to understand the functionality of the algorithms. This paper was meant as an introduction to the algorithms and also the purpose was to create an efficient way to display the clustering of individual data elements. If something remains to be done a more throughout analysis of the behavior of the algorithms should be made. Specially cases where the samples are badly balanced or in some particular order should be generated and analyzed.

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