

## In-Network Outlier Detection in Wireless Sensor Networks

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**Abstract** To address the problem of unsupervised outlier detection in wireless sensor networks, we develop an approach that (1) is flexible with respect to the outlier definition, (2) computes the result in-network to reduce both bandwidth and energy consumption, (3) uses only single-hop communication, thus permitting very simple node failure detection and message reliability assurance mechanisms (e.g., carrier-sense), and (4) seamlessly accommodates dynamic updates to data. We examine performance by simulation, using real sensor data streams. Our results demonstrate that our approach is accurate and imposes reasonable communication and power consumption demands.

**Keywords** Outlier detection · Wireless sensor networks

### 1 Introduction

Outlier detection, an essential step preceding most any data analysis routine, is used either to suppress or amplify outliers. Suppressing outliers (also known as data cleansing) improves the robustness of data analysis, for instance, in clustering [31], time series

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analysis [9] or text categorization [60]. Amplifying outliers helps find rare patterns in domains such as fraud analysis [47], intrusion detection, and Web purchase analysis.

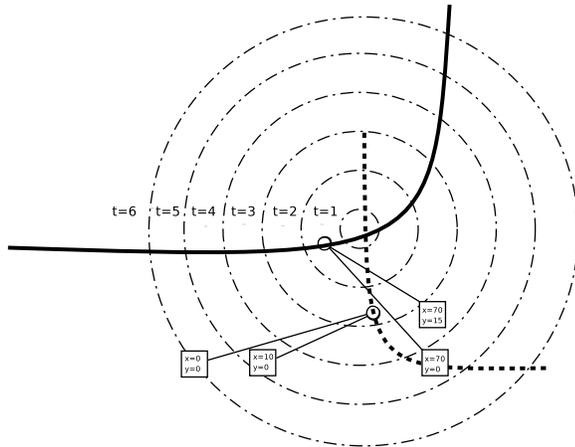
Several factors make wireless sensor networks (WSNs) especially prone to outliers. First, they collect their data from the real world using imperfect sensing devices. Second, they are battery powered and thus their performance tends to deteriorate as power dwindles. Third, since these networks may include a large number of sensors, the chance of error accumulates. Finally, when used for security and military purposes, sensors are especially prone to manipulation by adversaries. Hence, it is clear that outlier detection should be an inseparable part of any data processing routine in WSNs.

Simply put, outliers are observations whose probability of occurrence is extremely small. Since the actual distribution of the data is usually unknown, direct computation of probabilities is difficult. Because the problem is fundamental, a huge variety of outlier detection methods have been developed. In this paper we focus on non-parametric, unsupervised methods. A simplistic implementation of these methods would require collecting all data at one central node and executing the outlier detection algorithm there. Such a centralized solution has several disadvantages in WSNs [34]. The two most important are (i) the energy overhead incurred by sending the data across the WSN and (ii) the time delays incurred between data collection and processing. Both disadvantages increase proportionally to the average distance a data item needs to travel. They are also exacerbated by the dense flows arising in the areas close to the central node into which data from the entire WSN, by necessity of a centralized algorithm, converge. The delays increase in such areas because of likely packet collisions, so either the colliding packets need to be retransmitted, or bandwidth sharing needs to be imposed. Additionally, the nodes in such areas transmit far more than the average number of messages, and therefore expend their energy faster than other nodes.

We developed a technique for the computation of outliers in WSNs. This technique (1) is flexible with respect to the outlier definition, (2) computes the result in-network to reduce both bandwidth and energy consumption (see, e.g., [30]), (3) uses only single-hop communication, thus permitting very simple node failure detection and message reliability assurance mechanisms (e.g., carrier-sense), and (4) seamlessly accommodates dynamic updates to data. In addition to these essential features, the algorithm presented here has two highly desirable properties: it is generic – suitable for many outlier detection heuristics, and its communication load is proportional to the outcome (i.e., the number of outliers reported).

We exemplify the benefits of our algorithm by means of two different outlier detection heuristics and 53 sensors, which we simulate using the SENSE network simulator [20] on real data streams. Our results show that the algorithms both converge to an accurate result with reasonable communication load and power consumption. In most tested cases, our algorithms outperform the equivalent centralized algorithm.

The rest of the paper is organized as follows. In the next section, we provide a motivating example, in which a network of acoustic sensors attempts to locate the possible source of a sound. In Section 3, we discuss related work, including prior publications on outlier detection, wireless sensor networks, and distributed data mining. We introduce preliminaries in Section 4 and use them in Section 5 and 6 to describe the global and semi-global distributed outlier detection algorithms. A the detailed proof of the correctness of the former is given in the Appendix A. In Section 7 we present the performance evaluation of our algorithms and our conclusions in Section 8.



**Fig. 1** The expansion of a sound over time and the possible source location as computed by two different pairs of sensors according to the time difference of arrival. The origin of the sound lies in the intersection of the two hyperbolas.

## 2 Motivating Application

The importance of efficient outlier detection in wireless sensor networks is best understood in the context of popular applications of those systems. Consider, for instance, the acoustic source localization problem. In this problem, a set of synchronized sensors all register the arrival of a certain sound at a certain time. Given the distance of two sensors from one another and the time difference of arrival (TDOA) of the sound, the possible locations of the source vis-a-vis the two sensors can be deduced. Theoretically, given data from several sensors, the possible relative locations (each a hyperbola in the plane) can be intersected and the location of the source pinpointed (see, for example [3, 61] and Fig. 1).

In practice, the problem is much more complex. First, the real terrain in which the problem occurs is rarely flat, or an unobstructed three-dimensional space. Secondly, echoes and multiple concurrent sounds may add many possible hyperbolas from which the relevant ones have to be selected. Last, and perhaps most importantly, the method is sensitive to errors in sensor synchronization and positioning which may result from unsuccessful initialization, degradation with time, or power depletion. All these factors amount to a multiplicity of possible hyperbolas, only a few of which intersect at the correct location of the source.

A similar principle of localization applies in a broader setting, known as *binary sensing object location* [65], in which a single sensor detects an object's whereabouts to some accuracy and the neighboring nodes then cooperate to decrease measurement inaccuracy. Note that the modality of the signal and of the sensor (e.g., acoustic, seismic, visual, electromagnetic) is irrelevant. Also notice any detection, be it true or false, of object presence within a sensor's range will trigger a tracking algorithm or even the entire tracking service [21]. Such algorithm or service is usually associated with significant costs, which WSN applications often try to reduce or avoid by using, e.g., Maximum Likelihood [59]. Unfortunately, such error reduction methods often require

centralization of the raw data which, in itself, can be a costly process; especially when the sensor outputs a stream of measurements.

It is therefore crucial to be able to perform data cleaning in the network prior to any decision protocol. With the method suggested in this paper, sensors can constantly and efficiently prune away data suspected of being false. Only then, and only if the remaining data seems to require further analysis, would the more complex and costly procedure for source localization be executed. In this way, much energy can be saved and system lifetime extended. In the rest of this paper we avoid the discussion of any specific WSN application, which deserve thorough discussion in their own right (see, e.g., [12] for source localization and tracking). Instead, we focus on the detection and elimination of outliers.

### 3 Related work

#### 3.1 Outlier detection

Outlier detection is a long studied problem in data analysis, and its treatment has been thoroughly described in several studies [8, 33]. An outlier is an observation which appears to deviate markedly from other members of the sample in which it occurs. Outlier detection methods can be divided into two kinds: model based, and density based. Model based methods assume that the data follows a model which, excluding the specifics of its parameters, is known, so they focus on evaluating the model's parameters and singling out the outliers which obstruct their true value. Some methods assume a statistical model. For instance, the Tietjen-Moore test [64] assumes the data follows the normal distribution, with unknown mean and variance. However, many popular methods assume far more abstract models: e.g., that the non-outlier data can be modeled using a neural network [32, 45] or a decision tree [37].

Density based outlier detection methods evaluate the probability of an observation using the density of data near it. The advantage of this second approach is that no assumption of a global model for the data source is required but only a metric on the data domain<sup>1</sup>. Examples of such metrics include: distance to  $k^{th}$  nearest neighbor [11], [55]; average distance to the  $k$  nearest neighbors [6], [11]; inverse of the number of neighbors within a fixed distance  $\alpha$  [42]. A more complicated approach is to utilize both the local properties of the data (as do the metrics above) and its global properties, e.g., by denoting outliers the  $n$  most sparse data points [29]. Alternatively, methods such as LOF [18] use the ratio of the distance from an observation to its nearest neighbors to the average of their own distances to their nearest neighbors. The algorithms described in this work are suitable, as is, to all density based methods except LOF. Generalization to LOF, while apparently possible, is left for further research.

#### 3.2 Outlier detection in wireless sensor networks

WSNs combine the ability to sense, compute, and coordinate their activities with the ability to communicate results to the outside world. They have revolutionized data collection in all kinds of environments. Yet, the design and deployment of these networks

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<sup>1</sup> Often, methods which fall short of even the triangle inequality are sufficient.

creates unique research and engineering challenges due to their intended large size (up to thousands of sensor nodes), their often random and hazardous deployment, obstacles to their communication, their limited power supply, and their high failure rate. These limitations must be taken into account when developing WSN software. In [28], Estrin et al. introduce scalable coordination as an important software component. A survey on the state-of-the-art for WSNs is given in [5], and another survey [4] focuses on challenges arising from military, health care, ecology, and security applications.

Energy-efficiency is a cardinal WSN requirement and much research has focused on meeting it [7], [22]. One common strategy for achieving energy-efficiency is by minimizing communication using topology-control algorithms that dictate the active/sleep cycles of sensor nodes. Examples include ESCORT [16], ASCENT [19], STEM [56], and GAF [70]. While the focus of this paper is on WSN outlier detection, the challenge is the same. Hence, while we do not propose a topology-control algorithm, we do aim to design an energy-efficient algorithm by minimizing the required communication overhead.

Other research efforts have also focused on developing a framework for distributed outlier detection in WSNs. Zhuang et al. [72] use a weighted moving average approach to smooth noise from the data stream arriving at each sensor. Sensors also use data from neighboring sensors (spatial smoothing) to reduce data propagation to the sink by not sending observed data whose value remains within the established spatio-temporal trend. Unlike our approach, theirs does not seek to detect outliers.

Sheng et al. [58] developed a framework for the discovery of k-nearest-neighbor based outliers: points whose distance to their k-nn exceeds a fixed threshold or the top n points with respect to the distance to their k-nns. Each sensor maintains a histogram-type summary of pertinent information over a sliding window of its data points. The sink node collects these summaries and queries the network for any additional information needed to correctly determine the outliers over the whole network. The use of summaries allows less communication than a naive, centralized approach. Their approach differs from ours in several ways. First, they only detect outliers over one-dimensional data and difficulty of building compact, multi-dimensional histograms will hinder any extension beyond that. Second, they only consider the two k-nn based outlier definitions described above, while our approach encompasses these and more. Thirdly, their approach only applies in settings where spatial proximity is unimportant while our approach can, if needed, to accommodate spatial proximity ("semi-local" outlier detection).

Subramaniam et al. [63] require the sensors to maintain a tree communication topology and compute outliers using an estimate of the underlying probability distribution from which the data arises. Such an estimate is computed by each sensor maintaining a random sample of its data observations. Our approach differs in at least four ways. First, ours does not make any assumptions about the communication topology (e.g., that it is a tree), save that it is connected. Second, ours computes outliers with respect to all of the data observations at each sensor, not a sample. Third, ours can smoothly take into account spatial proximity among the sensors ("semi-local" outliers) while Subramaniam et. al do not focus on this task. Fourth, our approach is designed to smoothly adjust to changes in the underlying network topology while theirs requires that the underlying communication tree be reestablished by other means before the algorithm can resume operation.

Janakiram et al. [36] developed a framework based on a Bayesian Belief Network (BBN) that has been constructed over the WSN (and distributed to each sensor).

Using this, each sensor can estimate the likelihood of an observed tuple and, therefore, detect outliers, yet it is not clear to what extent the BBN construction phase can be carried out in-network. Moreover, the authors do not discuss the problem of updating the BBN given network/data change. In contrast, our processing is entirely in-network and smoothly adjusts to changes in data/network.

Zhuang and Chen [71] use a wavelet based technique for correcting large isolated spikes from single sensor data streams. A dynamic time warping (DTW) distance-based technique is also used to identify more steady intervals of erroneous sensor data by comparing the data streams of spatially close sensors assumed to produce similar streams. To reduce energy consumption, anomalous data streams are not transmitted to the base station. Our method is similar in that it is in-network. However, Zhuang and Chen’s use of DTW is tightly integrated with a minimum hop count routing algorithm, which makes the approach more restrictive than ours.

Rajasegarar et al. [54] describe an approach that is based on distributed non-parametric anomaly detection and requires sensors to maintain a tree communication network topology. Here each sensor clusters its sampled measurements using a fixed-width clustering algorithm, then extracts statistics of the clusters (i.e., the centroid and number of contained data vectors) and then sends them its parent node. The parent uses its children’s cluster statistics to form a merged cluster. The parent then transmits that cluster to its own parent. This process continues recursively until the base station receives all clusters, after which it will perform anomaly detection to identify all outliers. While this approach supports energy-efficiency by distributing the clustering operation throughout the network, anomaly detection is only performed at the base station. Our approach differs in that it distributes the anomaly detection process itself throughout the network, quickly enabling nodes to identify outliers and autonomously make further data processing decisions. Nor does our approach rely on the use and maintenance of a routing tree; hence it smoothly adjusts to changes in the underlying network topology.

Adam et al. [1] address the issue of accounting for spatially neighboring peers when detecting outliers in sensor networks. However, they assume the sensor datasets are centralized and the outlier processing is carried out at the central processing node. They do not consider the problem of carrying out the outlier detection *in-network* as we do.

Palpanas et al. [51] propose a technique for distributed deviation detection using a network hierarchy of low and high capacity sensors that are differentiated with respect to processing power and communication range. Here, low capacity sensors aim to detect local outliers while high capacity sensors detect more spatially dispersed outliers using an aggregation of low capacity sensors’ data. Kernel density estimators are used to model the distribution of data values reported by sensors and distance-based detection techniques are used for identifying outliers. The authors present no formal evaluation of the proposed technique. Our approach differs in that it does not rely on a hierarchy of device capabilities.

Radivojac et al. [53] address the process of sensors learning data distributions from class-imbalanced data. Here, sensors send data points to a central base station which generates a classification model from class-imbalanced data (i.e., having abundant negative samples and few positives). the total cost of detection and classification (e.g., costs of transmitting false positives and false negatives). In contrast, our framework operates in-network.

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This paper is an extension of our preliminary work appearing in conference proceedings [17]. In the extended version we provide complete correctness proofs for the global outlier detection algorithm along with improved experimental analysis. We have also added a localized outlier detection algorithm and experimental analysis of it.

### 3.3 Distributed data mining

Distributed Data Mining (DDM) has recently emerged as an important area of research. DDM is concerned with analysis of data in distributed environments, while paying careful attention to computation, communication, storage, and human-computer interaction. Detailed surveys of Distributed Data Mining algorithms and techniques have been presented in [38], [39], [40]. Some of the common data-analysis tasks include association rule mining, clustering, classification, kernel density estimation, and so on.

Recently, researchers have started to consider data analysis and data mining in large-scale dynamic networks with the goal of developing techniques that are highly asynchronous, scalable, and robust to network changes. Efficient data analysis algorithms often rely on efficient primitives, so researchers have developed several different approaches to computing basic operations (e.g., average, sum, max, or random sampling) on dynamic networks. Mehyar et al. [48] have developed an asynchronous, deterministic technique for computing an average over a large, dynamic network. Boyd et al. [15] and Kempe et al. [41] have investigated gossip based randomized algorithms. Jelasity and Eiben [43] develop the “newscast model.” Bawa et al. [10] have developed an approach in which similar primitives are evaluated to within an error margin. Wolff et al. [67] develop a local algorithm for majority voting. Datta and Kargupta [26] have developed a technique for uniformly sampling data distributed over a large-scale peer-to-peer network. Bhaduri et al. [13], Sharfman et al. [57], and Wolff et al. [69] have developed techniques for threshold monitoring over a large, distributed set of data streams.

Finally, some work has gone into more complex data mining tasks: association rule mining [67], facility location [44], decision tree induction [14], classification through meta-learning [46] (all four based on local majority voting), genetic algorithms [23], k-means clustering [27] [68], Web user community formation [25], hidden variable distribution estimation in a wireless sensor network [49], and outlier detection in distributed data streams [50] [62].

In the last two papers, Otey et al. [50] and Su et al. [62] address outlier detection over multiple data streams and wireless sensor networks in particular. Otey et al. agree with our analysis of two of the main challenges in computing outliers in a sensor network: The impracticality of collecting all the data to one site and the dynamic nature of the data. But their work differs from ours in that it ignores the problems associated with partial message loss and possible sensor failure during the computation. Therefore, the algorithms they describe use complete passes over all the (distributed) data. Such passes require that the sensors remain active during every execution of an algorithm and are sensitive to the dropping of even a few messages. Furthermore, the algorithms require some form of synchronization between all sensors (in order to move from one pass to the next) whereas ours only require the minimal level of synchronization – between the transmitter of a message and its nearby receivers. Achieving global near-synchronization in a wireless sensor network is not a trivial task. Finally, Otey et al. focus on their own definition of an outlier, *links*-based. While this semantic may be

suitable for some applications, we choose to focus on implementing more standard definitions of an outlier.

Su et al. take a very different approach. They employ a more standard definition of an outlier, kernel based density estimation, which is an approximation for distance based density. Their work differs from ours and that of Otey et al. mainly in its assumption of a hierarchical sensor network architecture in which leaders collect data from their subject nodes. While a hierarchical architecture greatly simplifies many algorithms, it is often unsuitable for real deployments of wireless sensor networks. The leaders, almost by definition, require many more resources than their subjects and therefore quickly deplete their power or suffer from communication congestion. When that happens, new leaders have to be selected; much system time is thus invested in maintaining the computational architecture. Therefore, the work of Su et al. might be better than ours for hybrid scenarios in which some sensors (e.g., those with fixed power connections) are naturally the leaders. Our work, in comparison, might be more suitable for the classic scenario in which identical sensors are randomly deployed and must self-organize to form a network.

## 4 Preliminaries

### 4.1 Outlier detection defined

Let  $\mathbb{D}$  be a data space. Following a common approach, our outlier detection method relies on a ranking function,  $R$ . This function maps  $x \in \mathbb{D}$  and finite  $D \subseteq \mathbb{D}$  to a non-negative real number  $R(x, D)$ , indicating the degree to which  $x$  can be regarded as an outlier with respect to a dataset  $D$ . Some common examples of  $R$  include the density metrics mentioned in the previous section. We assume that a fixed total linear order,  $\prec$ , on  $\mathbb{D}$  is used as a tie-breaking mechanism to ensure that  $R(\cdot, Q)$  creates a total linear ordering on  $\mathbb{D}$  for any finite  $Q \subseteq \mathbb{D}$ .

$R$  is assumed to satisfy the following two axioms. Given  $x \in \mathbb{D}$ , for all finite  $Q_1 \subseteq Q_2 \subseteq \mathbb{D}$ : **anti-monotonicity**,  $R(x, Q_1) \geq R(x, Q_2)$ ; **smoothness**, if  $R(x, Q_1) > R(x, Q_2)$ , then there exists  $z \in Q_2 \setminus Q_1$ , such that  $R(x, Q_1) > R(x, Q_1 \cup \{z\})$ . The anti-monotonicity axiom is similar to the *a-priori rule* in frequent itemset mining [2]. The smoothness axiom expresses the intuition that  $R$  changes gradually. As more points are added to  $Q_1$ , the rating function changes gradually to  $R(x, Q_2)$ . Of the examples in the previous paragraph, all but LOF satisfy these assumptions, assuming, as we do, the use of a tie-breaking mechanism as described in the previous paragraph.

Given  $n$ , a user-defined parameter, and a finite dataset  $D \subseteq \mathbb{D}$ , the outliers of  $D$  are denoted  $O_n(D)$  and are defined to be the top  $n$  points in  $D$  with respect to  $R(\cdot, D)$  (if  $|D| < n$ , then  $O_n(D)$  is defined to be  $D$ ).

### 4.2 Distributed system set-up

The distributed system architecture we assume consists of a collection of sensors,  $p_i$ , each producing a finite dataset  $D_i \subseteq \mathbb{D}$ .  $D_i$  only contains points that originated at sensor  $p_i$ , for example, a data point associated with a measurement made by  $p_i$ . Sensors communicate by exchanging messages with their immediate neighbors as defined by an undirected graph. We assume that messages are reliable, i.e., the sender will be

informed if a message is not received, and each sensor  $p_i$  can accurately maintain the list of its immediate neighbors,  $\Gamma_i$ , in the graph. Our algorithms work as long as there exists a path, possibly unknown, from each sensor to every other sensor. Note that message reliability is difficult to fully maintain in a WSN – some message dropping is expected. While our algorithm assumes no message dropping, the WSN simulator we used in our experiments realistically includes message dropping. We used an end-to-end acknowledgement mechanism to largely mitigate the effect of message dropping. A small number of messages are still lost, but the algorithm achieves 99% accuracy.

### 4.3 Notation Summary

Table 1 is provided as an aid to understanding the notation we use throughout the paper. A summarized definition is provided with each symbol listed, as well as, the subsection where a complete definition can be found. Note that some of the symbols listed in the table, *e.g.*,  $D_{i,j}^i$ , are only fully defined later in this paper.

$\mathbb{D}$ $R(.,.)$ $O_n(.)$	Space from which all data points are drawn. Outlier ranking function for a data point (first argument) with respect to a set of points (second). Top $n$ outliers in a set of data points with respect to $R(.,.)$ .	Section 4.1.
$p_i$ $\Gamma_i$ $D_i$	The $i^{th}$ sensor. The immediate neighbors of $p_i$ in the communication graph. The data points produced by $p_i$ .	Section 4.2.
$D$ $D_{i,j}^i$ $D_{j,i}^i$ $P^i$	All data points in the network. The data points $p_i$ is holding that $p_i$ sent to $p_j$ . The data points $p_i$ is holding that $p_j$ sent to $p_i$ . All the data points $p_i$ is holding.	Section 5 before 5.1.
$N(.,.)$	The nearest neighbor of a data point (first argument) among a set of points (second).	Section 5.1.
$[. .]$	The support set of a data point or set of points (second argument) with respect to another set of points (first).	Section 5.2.
$\bigcup_j D_j[\text{hops to } i \leq d]$ $x.hop$ $x.rest$ $Q^{\leq h}$ $[Q]^{min}$	The union of all $D_j$ where the hop distance between $p_j$ and $p_i \leq d$ . An addition field on data point $x$ to account for hop distance. The remaining fields on data point $x$ . The subset of data points in $Q$ with hop field $\leq h$ . The result of replacing all points in $Q$ that differ only in their hop field by the point with the smallest hop field.	Section 6 before 6.1.

**Table 1** Notation table: symbol, summarized definition, and section where complete definition appears.

## 5 Global distributed outlier detection algorithm

In this section, we describe a distributed algorithm by which sensors, each assumed to know  $R$  and  $n$ , compute  $O_n(D)$  where  $D = \bigcup_i D_i$  (*global* outlier detection). In a wireless sensor network, it can be desirable for sensors to find outliers only with respect to the data contained in nearby sensors, rather than in the entire network (*semi-global* outlier detection). In section 6, we describe how to modify the global outlier detection algorithm to act in a semi-global manner.

At any point in time,  $p_i$  keeps track of the data points it has sent to or received from its neighbor  $p_j$  at some past time. Let  $D_{i,j}^i$  denote the set of points sent from  $p_i$  to  $p_j$ , and,  $D_{j,i}^i$  denote the set of points sent from  $p_j$  to  $p_i$ . Importantly,  $(D_{i,j}^i \cup D_{j,i}^i)$  denotes the data points that  $p_i$  can be sure are commonly held with  $p_j$  (there may be more). Let  $P_i$  denote  $D_i \bigcup_{j \in \Gamma_i} D_{j,i}^i$ , the set of points  $p_i$  is holding at the current time.<sup>2</sup>  $p_i$  uses  $P_i$  to compute an estimate of the overall correct answer,  $O_n(D)$ . The *estimate* of  $p_i$  is  $O_n(P_i)$ , the set of outliers based on all the information available to  $p_i$  at the current time.

The algorithm does not assume any special sensors. Each sensor,  $p_i$ , asynchronously waits for an *event* to occur: (i) the algorithm is initialized, (ii)  $D_i$  changes, (iii) a message is received from a neighbor, or (iv) a link goes up/down, causing the immediate neighborhood of  $p_i$  to change (however, algorithm correctness requires that we assume the network remain connected). Note that events for  $p_i$  are entirely local and can be detected without the aid of any other sensors beyond the immediate neighborhood. Once  $p_i$  detects an event, it will decide which of the points it is currently holding ( $P_i$ ), if sent, could cause its neighbor,  $p_j$ , to change its estimate.  $p_i$  then sends these points and adds them to  $D_{i,j}^i$  ( $p_i$  carries out this process separately for all of its neighbors).

Gradually, the set of points held by each sensor becomes large enough to obtain overlap so that each sensor's estimate is the correct answer,  $O_n(D)$ . This will be guaranteed to occur once each sensor, individually, decides that none of the points it is currently holding need be sent to its neighbors. At this point, the algorithm is terminated. To see how all of this works, consider an example.

### 5.1 Example

Let  $R$  be the distance to the nearest neighbor and, given  $x \in \mathbb{D}$  and finite  $P \subseteq \mathbb{D}$ , let  $N(x, P)$  denote the nearest neighbor of  $x$  among points in  $P$ . Given finite  $Q \subseteq \mathbb{D}$ , let  $N(Q, P)$  denote  $\bigcup_{x \in Q} N(x, P)$ . Let  $n = 1$  and consider a network of two sensors,  $p_i$  and  $p_j$ , each initially holding the following one-dimensional datasets. The correct answer the algorithm will compute is  $O_n(D) = \{0.5\}$ .

- $D_i = \{0.5, 3, 6, 10, 11, \dots, a\}$ .
- $D_j = \{4, 5, 7, 8, 9, a + 1, a + 2, \dots, a + b\}$ .
- $D = D_i \cup D_j = \{0.5, 3, 4, 5, 6, 7, 8, 9, 10, 11, \dots, a, a + 1, \dots, a + b\}$ .

Initially,  $P_i = D_i$ ,  $P_j = D_j$ , and  $D_{i,j}^i = D_{j,i}^i = D_{i,j}^j = D_{j,i}^j = \emptyset$ . For simplicity, we will describe the algorithm in synchronous fashion starting with  $p_i$ . But the ideas extend nicely to asynchronous operation.

<sup>2</sup> Note the distinction between  $D_i$  and  $P_i$ .  $D_i$  is the set of points that *originated* at sensor  $p_i$ , while  $P_i$  is the set of all points that  $p_i$  is holding including  $D_i$  and those originating at other sensors but propagated to sensor  $p_i$  through message passing.

1.  $p_i$  computes its estimate as  $O_n(P_i) = \{6\}$  and then must compute the set of its data points that might cause  $p_j$  to change its estimate if sent. We call these the *sufficient points* from  $P_i$  for sensor  $p_j$ . Formally, we define a set  $Z_j \subseteq P_i$  to be sufficient for  $p_j$  if

$$[O_n(P_i) \cup N(O_n(P_i), P_i)] \cup [N(O_n(D_{i,j}^i \cup D_{j,i}^i \cup Z_j), P_i)] \subseteq Z_j. \quad (1)$$

The rationale for the first part is simple.  $O_n(P_i)$  is necessary for  $p_j$ , because if  $p_i$  is right in its estimate, then  $p_j$  ought to know about it. Moreover,  $p_i$  must also send  $N(O_n(P_i), P_i)$ , because this allows  $p_j$  to determine if any of its points can cause the ranking of the points in  $O_n(P_i)$  to change.

The rationale for the second part is somewhat more complex. In brief, if  $p_i$  were to send  $Z \subseteq P_i$  to  $p_j$ , then  $p_i$  would also need to send  $N(O_n(D_{i,j}^i \cup D_{j,i}^i \cup Z), P_i)$ . To avoid re-sending,  $p_i$  requires that  $N(O_n(D_{i,j}^i \cup D_{j,i}^i \cup Z), P_i)$  be contained in  $Z$ . To understand the reasoning for all of this, consider that  $(D_{i,j}^i \cup D_{j,i}^i \cup Z)$  is the total set of points  $p_i$  knows  $p_j$  would have if  $Z$  had been sent to  $p_j$ . Thus,  $O_n(D_{i,j}^i \cup D_{j,i}^i \cup Z)$  is the best approximation  $p_i$  would have to the estimate of  $p_j$  if  $Z$  were sent to  $p_j$ . Hence,  $p_i$  computes its nearest neighbors to these, because, if  $p_i$  is right in its approximation, it must ensure that  $p_j$  have these neighbors since they could cause  $p_j$  to change its estimate.

Getting back to our example, observe that  $O_n(P_i) \cup N(O_n(P_i), P_i) = \{3, 6\}$ . Moreover,  $N(O_n(D_{i,j}^i \cup D_{j,i}^i \cup \{3, 6\}), P_i) = N(O_n(\{3, 6\}), P_i) = \{3\} \subseteq \{3, 6\}$ . Hence,  $Z_j = \{3, 6\}$ , as this satisfies (1) above. So,  $p_i$  sends  $\{3, 6\} \setminus (D_{i,j}^i \cup D_{j,i}^i) = \{3, 6\}$  and updates  $D_{i,j}^i$  to  $\{3, 6\}$ .

2.  $p_j$  will receive these points and update  $D_{i,j}^j$  to  $\{3, 6\}$  (currently  $D_{j,i}^j = \emptyset$ ). It is thus implicit that  $P_j$  now denotes  $D_j \cup D_{i,j}^j = \{3, 4, 5, 6, 7, 8, 9, a+1, a+2, \dots, a+b\}$ .  $p_j$  computes  $O_n(P_j) \cup N(O_n(P_j), P_j) = \{3, 4\}$  (assuming appropriate tie-breaking). Moreover, it can be seen that this satisfies (1). So,  $p_j$  sends  $\{3, 4\} \setminus (D_{i,j}^j \cup D_{j,i}^j) = \{3, 4\} \setminus \{3, 6\} = \{4\}$  and updates  $D_{j,i}^j$  to  $\{4\}$ .
3.  $p_i$  receives these points and updates  $D_{j,i}^i$  to  $\{4\}$  (currently  $D_{i,j}^i = \{3, 6\}$ ). It is thus implicit that  $P_i$  now denotes  $D_i \cup D_{j,i}^i = \{0.5, 3, 4, 6, 10, 11, \dots, a\}$ .  $p_i$  computes  $O_n(P_i) \cup N(O_n(P_i), P_i) = \{0.5, 3\}$ . Moreover, it can be seen that this satisfies (1). So,  $p_i$  sends  $\{0.5, 3\} \setminus (D_{i,j}^i \cup D_{j,i}^i) = \{0.5, 3\} \setminus \{3, 4, 6\} = \{0.5\}$  and updates  $D_{i,j}^i$  to  $\{0.5, 3, 6\}$ .
4.  $p_j$  will receive these points and update  $D_{i,j}^j$  to  $\{0.5, 3, 6\}$  (currently  $D_{j,i}^j = \{4\}$ ). It is now implicit that  $P_j$  now denotes  $D_j \cup D_{i,j}^j = \{0.5, 3, 4, 5, 6, 7, 8, 9, a+1, a+2, \dots, a+b\}$ .  $p_j$  computes  $O_n(P_j) \cup N(O_n(P_j), P_j) = \{0.5, 3\}$ . Moreover, it can be seen that this satisfies (1). So,  $p_j$  sends  $\{0.5, 3\} \setminus (D_{i,j}^j \cup D_{j,i}^j) = \{0.5, 3\} \setminus \{0.5, 3, 4, 6\} = \emptyset$ . In other words, nothing is sent.

At this point the algorithm has terminated, both sensors are waiting for an event to occur, and there are no messages in flight.  $P_i$  and  $P_j$  denote  $\{0.5, 3, 4, 6, 10, 11, \dots, a\}$  and  $\{0.5, 3, 4, 5, 6, 7, 8, 9, a+1, a+2, \dots, a+b\}$ , respectively. Therefore  $O_n(P_i) = \{0.5\} = O_n(P_j)$ , which in turn, equals the correct answer  $O_n(D)$ . Observe that the total amount of communication (data points sent) was 4. The naive approach, which centralized all the data on either  $p_i$  or  $p_j$ , requires  $\min\{a-6, b+5\}$  communication. For large  $\min\{a, b\}$ , the distributed algorithm requires much less communication.

## 5.2 The algorithm

To translate the previous example into a formal algorithm for general  $R$  (satisfying the anti-monotonicity and smoothness axioms), we must provide some definitions generalizing the role of  $N(.,.)$ . Given  $x \in \mathbb{D}$  and finite  $P \subseteq \mathbb{D}$ ,  $Q_1 \subseteq P$  is called a *support set of  $x \in \mathbb{D}$  over  $P$*  if  $R(x, P) = R(x, Q_1)$ . Intuitively, we can see that all other points from  $P$  can be discarded without affecting the rank of  $x$ . Using cardinality and the tie-breaking mechanism discussed earlier ( $\prec$  a total linear order on  $\mathbb{D}$ ), we can define a unique, smallest support set of  $x$  with respect to  $P$ , denoted  $[P|x]$ .<sup>3</sup> Given  $Q \subseteq P$ , let  $[P|Q]$  denote  $\bigcup_{x \in Q} [P|x]$ . In the previous example,  $R$  was the distance to the nearest neighbor, so,  $[P|x]$  equaled  $N(x, P)$  using  $\prec$  to break ties.

With these more general definitions, we define a set  $Z_j \subseteq P_i$  to be sufficient for  $p_j$  if

$$(O_n(P_i) \cup [P_i|O_n(P_i)]) \cup \left( [P_i|O_n(D_{i,j}^i \cup D_{j,i}^i \cup Z_j)] \right) \subseteq Z_j. \quad (2)$$

Due to the broadcast nature of wireless sensor network communication,  $p_i$  cannot send points to one immediate neighbor without the other neighbors receiving them as well. In light of this, the algorithm accumulates all points (tagged with recipient IDs) to be sent to all immediate neighbors in a single packet,  $M$ . When an immediate neighbor,  $p_j$ , receives  $M$ , the neighbor extracts those points that are tagged with ID  $j$ . If no points are tagged as such,  $p_j$  does not regard receipt of  $M$  as an event.

$p_i$  detects an event if one of the following occurs: (i) the algorithm is initialized, (ii)  $D_i$  changes, (iii)  $M$  is received and contains points tagged with  $i$  (i.e., points are received from a neighbor), or (iv) a link goes up/down, causing  $p_i$ 's immediate neighborhood to change (however, algorithm correctness requires that we assume the network remain connected). In response,  $p_i$  carries out the following algorithm whose pseudo-code is given in the ‘‘Global Outlier Detection’’ Figure. First,  $P_i$  is updated to account for all  $p_j$  from which points were received in  $M$ . Only points not already in  $P_i$  are added to  $D_{j,i}^i$ . The first two steps in the main for-loop (‘‘For each  $j \in \Gamma_i$ , do’’) compute a  $Z_j$  satisfying Equation (2), although the result is not guaranteed to be the smallest set to do so. The ‘‘If...then’’ in the main for-loop tests whether there are any points found sufficient for  $p_j$  that  $p_i$  cannot already be sure  $p_j$  has, i.e., points in  $Z_j$  but not in  $D_{j,i}^i \cup D_{i,j}^i$ . If any such points are found, they are added to  $M$  along with their recipient ID  $j$ .

## 5.3 Streaming data and peer addition/deletion

We now extend the algorithm to one supporting a sliding window over a stream of data sampled at each sensor. To support this, we assume each point is time-stamped when it is sampled. We further assume that sensor clocks are synchronized to the extent required by the application for whom outlier detection is a preceding step. Thus, sensor  $p_i$  can delete all points in  $P_i$  (regardless of where they were originally sampled) once their time-stamp indicates they have expired.

The algorithm still needs to be modified to accommodate the addition of sensors during operation. All that is required is to treat the arrival of a new sensor as an event

<sup>3</sup> Formally, given  $Q_1, Q_2 \subseteq P$  support sets of  $x$  with respect to  $P$ ,  $Q_1$  is smaller than  $Q_2$  if  $|Q_1| < |Q_2|$  or ( $|Q_1| = |Q_2|$  and  $Q_1$  is lexicographically smaller than  $Q_2$  with respect to  $\prec$ ).

**Fig. 2** Global Outlier Detection**Algorithm 1**


---

```

– Set  $M = \emptyset$ , and, update  $P_i$  accounting for all neighbors  $p_j$  from which points were
received. For each point  $x$  received from  $p_j$ , do the following. If  $x$  is not already  $P_i$ , then
add  $x$  to  $D_{i,j}^i$ .
– For each  $j \in \Gamma_i$ , do
– – Set  $Z_j = O_n(P_i) \cup [P_i | O_n(P_i)]$ .
– – Repeat until no change:  $Z_j = Z_j \cup [P_i | O_n(D_{i,j}^i \cup D_{j,i}^i \cup Z_j)]$ .
– – If  $Z_j \setminus (D_{i,j}^i \cup D_{j,i}^i)$  is non-empty, then
– – – Append  $\langle j, Z_j \setminus (D_{i,j}^i \cup D_{j,i}^i) \rangle$  to  $M$ .
– – – Add points in  $Z_j \setminus (D_{i,j}^i \cup D_{j,i}^i)$  to  $D_{i,j}^i$ .
– – End If.
– End For.
– If  $M$  is non-empty, broadcast it to all sensors in  $\Gamma_i$ .
end

```

for the new sensor and for all its immediate neighbors. The algorithm can also be modified to accommodate the removal of sensors (e.g., when their battery is depleted), assuming that the network remains connected. In the sliding window model, a simple strategy is to merely allow points that originated with the removed sensor to age out of the window. The consequence of that would be that the result would not be the same as would be computed from the data of all *active* sensors. This usually poses no practical problem. A more general and complex solution is to propagate messages into the network, causing sensors to explicitly delete those points that originated with the removed sensor. We leave the details of this approach to future work.

#### 5.4 Algorithm correctness

The correctness of the algorithm can be proved in the following sense: if the data and network links remain static (and the network is connected), then communication will eventually stop, at which point all the sensors' outlier estimates will equal  $O_n(D)$ . This does *not* mean that the algorithm cannot handle dynamic changes or network links; it merely means that the algorithm will respond to such changes or links by converging to the correct answer. Since convergence takes time, and since during that time further changes are possible convergence to a globally correct value is only assured on the condition that the data and network remain unchanged long enough.

It is easy to see that, barring changes in the data or network, the algorithm will always terminate. So, the proof proceeds in two steps. First, upon termination, all sensors have the same outlier estimates and support (Theorem 1). Second, the consistent estimates shared by all sensors are indeed the correct ones (Theorem 2). Proofs of Theorems 1 and 2 are provided in Appendix A.

**Theorem 1** *Assuming a connected network, if for all sensors  $p_i$ ,  $D_i$  and  $\Gamma_i$  do not change, then upon termination of the algorithm, all sensors' outlier estimates and supports agree. I.e., for all  $p_i, p_j$ ,  $O_n(P_i) = O_n(P_j)$  and  $[P_i | O_n(P_i)] = [P_j | O_n(P_j)]$ .*

**Theorem 2** *Assuming a connected network, if for all sensors  $p_i$ ,  $D_i$  and  $\Gamma_i$  do not change, then upon termination of the algorithm, all sensors' outlier estimate will be correct. I.e., for all  $p_i$ ,  $O_n(P_i) = O_n(D)$ .*

**Comments: 1)** Theorem 1 holds without the smoothness axiom; hence, for any anti-monotonic  $R$ , upon convergence, all sensors will agree on their outlier estimate and support. However, without the smoothness axiom, Theorem 2 does not hold, i.e., the consistent outlier estimates might not be the correct ones. Counter-examples exist showing how an anti-monotonic but not smooth  $R$  causes the algorithm to terminate without all sensors agreeing upon the correct set of outliers.

**2)** For an arbitrary  $R$ , it is not clear how to efficiently compute  $[P|x]$  and we do not address the issue. However, efficient computation is straightforward for the  $R$  we consider in our experiments: average distance to the  $k^{th}$  nearest neighbor.

**3)** The correctness proof requires reliable communication, e.g., no message dropping. The WSN simulator we used in our experiments includes message dropping. We used an end-to-end acknowledgement mechanism to largely mitigate the effect of message dropping, and the algorithm achieved 99% accuracy.

## 6 Semi-Global Distributed Outlier Detection Algorithm

For many applications, computation only takes into account samples which were taken close to one another. In such applications outliers should also be computed respective to the data sampled by nearby sensors, rather than by the entire network. In this section, we describe how to modify the global outlier detection algorithm to act in a semi-global manner. Under this approach, each sensor computes outliers only from within those points sampled in its spatial proximity.

To account for spatial locality, we use *hop distance*: the number of hops between two sensors along their shortest path in the underlying communication network. Given integer  $d$  and sensor  $p_i$ , let  $\bigcup_j D_j[\text{hops to } i \leq d]$  denote the union of all  $D_j$  such that  $p_j$  and  $p_i$  have hop distance no greater than  $d$ . The semi-global outlier detection problem requires each  $p_i$  to compute  $O_n(\bigcup_j D_j[\text{hops to } i \leq d])$ . Setting  $d$  to infinity yields the global outlier detection problem discussed earlier.

To account for hop distance, each data point  $x$  has an additional field  $x.hop$  (at birth  $x.hop$  is set to zero). Let  $x.rest$  denote all the remaining fields – these are the ones used by the rating function  $R$ . Given a set of points  $Q$ , for  $0 \leq h \leq d$ , let  $Q^{\leq h}$  be the set of points  $x \in Q$  with  $x.hop \leq h$ . Let  $[Q]^{min}$  be the result of replacing all points that differ only in their hop field by the point with the smallest hop field. For example, consider  $Q = \{w, v, x, y, z\}$  where  $w.rest = v.rest$ ,  $x.rest = y.rest = z.rest$ , and  $v.rest \neq x.rest$ . If  $w.hop < v.hop$  and  $x.hop < y.hop, z.hop$ , then  $[Q]^{min} = \{w, x\}$ .

### 6.1 Semi-global outlier detection algorithm

The basic idea of the semi-global outlier detection algorithm is that each sensor  $p_i$  will run the global outlier detection algorithm over only those points arising on sensors within  $d$  hops. At first glance, the following simple modification of the global algorithm seems adequate. Before  $p_i$  sends a copy of a point,  $x$ , to its neighbors, it first increments  $x.hop$  and sends only if  $x.hop \leq d$ . Unfortunately, such a simple modification will not work. It does not take into account the fact that  $x$  must not have any effect on the outlier determination process of sensors  $p_j$  whose distance from  $p_i$  is more than  $d - x.hop$ . Examples can be demonstrated wherein this omission causes an incorrect overall result.

**Fig. 3** Semi-Global Outlier Detection**Algorithm 2**

```

– Set  $M = \emptyset$ , and, update  $P_i$  accounting for all neighbors  $p_j$  from which points were
received. For each point  $x$  received from  $p_j$ , do the following. If there does not exist  $y$  in
 $P_i$  with  $x.rest = y.rest$ , then add  $x$  to  $P_i$  (and update  $D_{i,j}^i$ ), otherwise if  $x.hop < y.hop$ ,
then replace  $y$  with  $x$  in  $P_i$  (updating as needed  $D_i$  and  $D_{f,i}^i$  for each  $f \in \Gamma_i$ ).
– For each  $j \in \Gamma_i$ , do
– – For  $h = 0$  to  $d - 1$ 
– – – Set  $Z_j^h = O_n(P_i^{\leq h}) \cup [P_i^{\leq h} | O_n(P_i^{\leq h})]$ .
– – – Repeat until no change:  $Z_j^h = Z_j^h \cup [P_i^{\leq h} | O_n(D_{i,j}^{i,\leq h} \cup D_{j,i}^{i,\leq h} \cup Z_j^h)]$ .
– – – Increment the hop field for each point in  $Z_j^h$ .
– – End For.
– – Set  $Z_j = [\bigcup_{h=0}^{d-1} Z_j^h]^{min}$ .
– – Remove points  $x$  from  $Z_j$  such that there exists  $y \in (D_{i,j}^i \cup D_{j,i}^i)$  with  $x.rest = y.rest$ 
and  $y.hop \leq x.hop$ .
– – If  $Z_j$  is non-empty, then
– – – Append  $\langle j, Z_j \rangle$  to  $M$ .
– – – Update  $D_{i,j}^i$  by adding the points in  $Z_j$ .
– – End If.
– End For.
– If  $M$  is non-empty, broadcast it to all sensors in  $\Gamma_i$ .
end

```

To avoid this problem,  $p_i$  must partition  $P_i$  into  $d$  parts:  $P_i^{\leq h}$  for  $0 \leq h \leq d - 1$ . For each of these parts, the global outlier detection algorithm is applied independently. Upon detecting an event (defined as before),  $p_i$  executes the algorithm whose pseudo-code is given in Figure 2.

First,  $P_i$  is updated to account for all  $p_j$  from which points were received in  $M$ . Because of the hop fields, the update step is somewhat more complicated than that of the global outlier detection algorithm. A point  $x$  from  $p_j$  is added to  $D_{j,i}^i$  if there does not exist  $y \in P_i$  with  $x.rest = y.rest$  ( $x$  does not already appear in  $P_i$ ). Or, if there does exist  $y \in P_i$  with  $x.rest = y.rest$ , but  $x.hop < y.hop$ , then  $x$  replaces  $y$  in  $P_i$  (updating as needed  $D_i$  and  $D_{f,i}^i$  for each  $f \in \Gamma_i$ ). Note that there cannot be more than one  $y$  with the same rest fields as  $x$  since all but the point with the smallest hop would have been removed earlier.

Next, for each neighbor  $p_j$  and each  $0 \leq h \leq d - 1$ , a set  $Z_j^h$  is computed which satisfies (2) with  $Z_j$ ,  $P_i$ ,  $D_{i,j}^i$ , and  $D_{j,i}^i$  replaced by  $Z_j^h$ ,  $P_i^{\leq h}$ ,  $D_{i,j}^{i,\leq h}$ , and  $D_{j,i}^{i,\leq h}$ , respectively. This computation is done by the first steps inside the nested for loops. Then, the hop field for each point in  $Z_j^h$  is incremented in preparation for sending to  $p_j$ .

Once all  $0 \leq h \leq d - 1$  have been processed for  $p_j$  (the inner for loop completes),  $Z_j^1 \dots Z_j^{d-1}$  are unioned and redundancies are eliminated. For any pair of points  $x, y$  in  $\bigcup_{h=0}^{d-1} Z_j^h$ , if  $x.rest = y.rest$  and  $x.hop < y.hop$ , then  $y$  is dropped (this action is signified by the ‘min’ superscript in the step immediately after the inner for loop). Then, all points  $x$  are removed from  $Z_j$  if there exists a point  $y$  in  $(D_{i,j}^i \cup D_{j,i}^i)$  with the same rest fields but  $y.hop \leq x.hop$ . If the resulting  $Z_j$  is non-empty, then these points are added to  $M$  (along with ID  $j$ ) for broadcast to neighbors, and  $D_{i,j}^i$  is updated by adding the points in  $Z_j$ .

## 7 Performance Evaluation

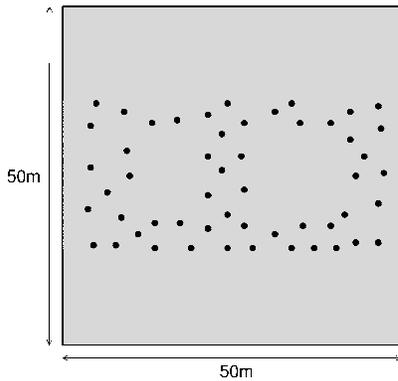
### 7.1 Experimental setup

We used the SENSE wireless sensor network simulator [20] to evaluate the performance of the global and semi-global outlier detection algorithms. Specifically, we analyzed the following metrics: (1) the accuracy of the algorithms in detecting outliers; (2) the average energy consumption per node per sampling period, as well as its breakdown to transmission related, and reception related consumption; and (3) the minimum and maximum energy consumption in the network. We observed both the global and semi-global outlier detection algorithms to be highly accurate; nodes converged upon the correct results approximately 99% of the time. We attribute any detection error to dropped packets (despite the "end-to-end" acknowledgement mechanism being used). Since average detection accuracy was consistent across all simulation parameters, we did not include any accuracy-related plots.

Various scenarios were used to analyze the performance of our algorithms. First, we compared our algorithms' energy consumption to that of a purely centralized outlier detection algorithm. Here, all nodes periodically sent their sliding window contents to a central node, which detected outliers using the unioned data sets and returned the outliers back to the nodes. For simplicity, we configured the centralized algorithm to calculate only global outliers because for this algorithm, energy consumption is independent of whether global or semi-global outliers are detected. Also, all algorithms (including the centralized solution) were evaluated using the following two outlier ranking functions ( $R$ , in our terminology): *distance to nearest neighbor* (NN) and *average distance to  $k$  nearest neighbors* (KNN).

We chose a centralized algorithm for our comparison because, to the best of our knowledge, there exist no comparable distributed solutions for WSN outlier detection. We find such a comparison to still be valid as many WSN deployments continue to employ centralized configurations, citing ease of administration as well as maintenance of a single (and "standard") point of interface with the growing number of applications and systems in the sense-and-respond computing domain. Such reasons, however, do not preclude the utility of a distributed algorithm such as ours, since it remains very useful as a general data processing solution for a wide range of applications, centralized are not.

For our data sets, we used real-world recorded sensor data streams from [35], in which distributed data samples are both spatially and temporally correlated. The data set is composed of a series of data samples describing environmental phenomena such as heat, light, and temperature from 53 sensors which periodically transmitted individual data samples to a central base station. The data set has some missing data points, which to the best of our knowledge is largely due to packet loss. Following accepted practices, we replaced missing data points with the average values of the data points within the sliding windows that preceded the missing points. This helped retain the temporal trends of the data streams. The data points contained the following features: (1) ID of the sensor that produced the data point; (2) epoch (sequential number denoting the data point's position in the sensor's entire stream); (3) data value (we specifically used temperature); and (4) x,y location coordinates. We used the data points' temperature value and location coordinates as inputs into the outlier rating functions. The location coordinates can represent the place where the measurements were taken or an estimate of a target's position or some other spatial information. It is important to note that



**Fig. 4** 53 node network topology (from [35]) used in all simulations reported.

these coordinates are attributes of the data on which our algorithm works in the example. They might become erroneous or anomalous just like any other data attribute, due to inaccurate initialization, power degradation, or a transmission error. The algorithm itself, however, would work the same regardless of whether such coordinates are given.

We originally simulated two networks, one containing all of the 53 nodes, with each node placed according to the coordinates in the data (thus matching the real topology from [35]), and the other containing 32 randomly selected nodes. By comparing the two networks we intended to study the scale-up properties of the algorithms. However, since the simulations revealed no real performance difference between the two networks, we eventually chose to only include results from the simulation of the full, 53 node network. The topology of the network is shown in Figure 4.

The terrain dimensions in our network simulation are  $50\text{m} \times 50\text{m}$ . Most hardware specifications claim that a sensor node's transmission range typically reaches up to approximately 250m, when properly elevated. However, when placed on the ground the reported ranges are much smaller [73] and, for reliable indoor communication using Crossbow motes, the effective range drops to a few meters [66]. Therefore, we configured all nodes to have a uniform transmission range of approximately 6.77m. We also used the hardware energy model based on the Crossbow mote specifications [24], with a transmit/receive/idle power setting of 0.0159W/0.021W/3e-6W (assuming a 3V power source). We simulated the wireless transport medium using the free-space signal propagation model.

Two protocols were used for routing. For the distributed algorithms, we used simple broadcast (as opposed to unicast) transmission with promiscuous listening that allowed all nodes to send data points to all their adjacent neighbors using one transmission. For the centralized algorithm, we used the well-accepted AODV [52] wireless routing protocol for multi-hop communication. We note that a simple end-to-end acknowledgment mechanism was also used to reinforce reliable communication. While alternative protocols do exist for more data-centric and energy-efficient communication, our main goal was straightforward comparison of the algorithms without having to balance the various advantages of each.

All simulations were run for 1000 seconds of simulated time and were repeated up to 12 times using different random number generator seed values to obtain averaged results (most results are based on 12 simulation runs, some were based on 6). Error bars are included with each point in each plot showing 95%, Student's T, confidence intervals (some bars are very small and appear not to be present). As shown in the following plots (and the tables in Appendix B), we collected results for different values of the following algorithm parameters: (1) the length of the node's sliding window,  $w$ ; and (2) the number of outliers to be reported,  $n$ . Additionally, for the distributed localized outlier detection algorithm, the hop diameter was varied from one to three. Plot data are labeled as follows: (1) *Centralized* for results obtained with the centralized algorithm; (2) *Global-NN* and *Global-KNN* for results obtained using distributed global outlier detection with NN and KNN outlier detection ranking functions ( $R$ ), respectively; and (3) *Semi-global, epsilon = x* for all results obtained using distributed localized outlier detection with hop diameter  $x$ . For brevity, we will often refer to the different algorithms by these labels.

## 7.2 Experimental results

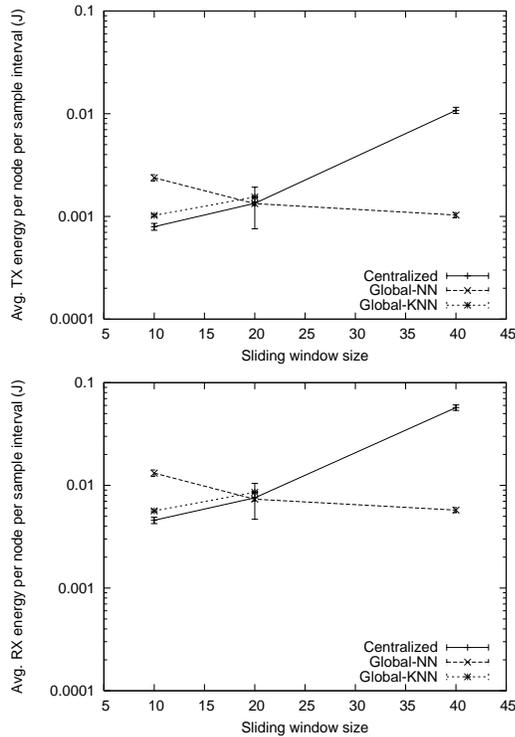
The numbers (averages, confidence intervals) from all of the following plots are displayed in tables in Appendix B.

### 7.2.1 Effect of sliding window size

Figure 5 compares network energy consumption per sampling interval with increasing  $w$  but  $n$  and  $k$  fixed at 4. Here, we show separate plots for transmission (TX) and reception (RX) energy for the reader who is interested in the disparity between the energy consumption due to different radio operations. We note that data points are missing for Global-KNN at  $w=40$ , due to a computational limitation of our simulator. However, preliminary results based on similar simulations and statistics are shown in [17]. Since the trends between both sets of results are nearly identical for the non-missing data points, it is reasonable to extrapolate the values of the missing points here.

Both figures show that as  $w$  increases, Global-NN is the only algorithm that reduces its energy consumption. Figure 5 shows that Global-NN eventually becomes the most energy-efficient solution given the domain of  $w$ . We attribute Global-NN's reduction in energy consumption to an increase in incoming data redundancy as the size of the sliding window increases. Since Global-NN only uses one supporting point to determine an outlier, the probability of finding new outliers or supports with this scheme in each new time interval as the sliding window increases is low.

Global-KNN and Centralized consume, as can be seen in Figure 5, more energy as  $w$  increases. However, given comparable results in [17], we can extrapolate that Global-KNN's energy consumption is a concave increasing function of  $w$ , whereas Centralized's is a convex increasing function, which makes the latter comparatively less energy-efficient in that it approaches a point of network failure at a higher rate. By contrast, the energy trend for Global-KNN for this data set indicates that when global outliers are defined by multiple supporting points (in this case 4), the increasing time interval from which the points are chosen has a less drastic effect on the number of messages required for the algorithm to converge. Overall, in cases where a user prefers to use a

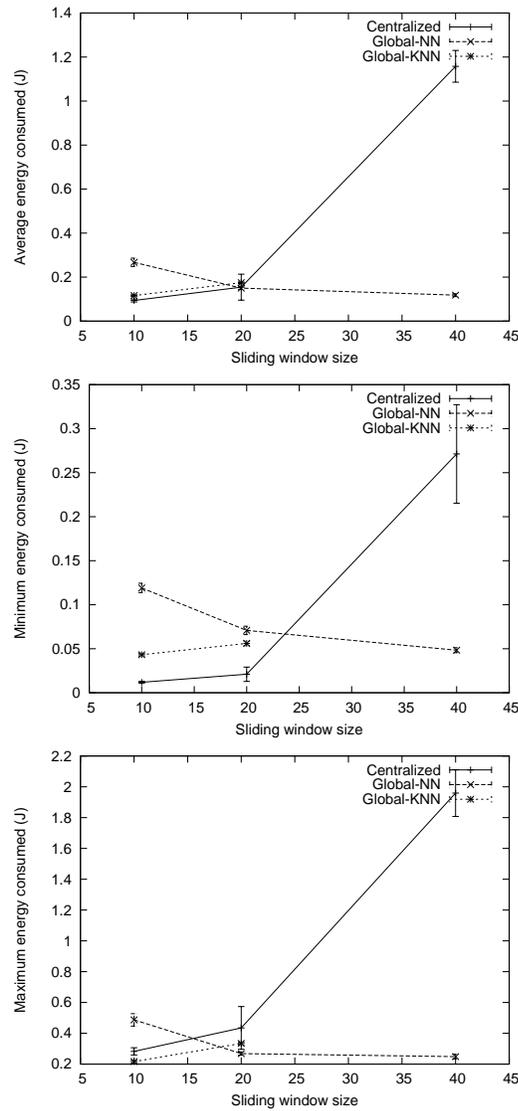


**Fig. 5** Average transmission and reception energy consumption per node per sample interval vs.  $w$  ( $n=4$ ,  $k=4$ ) for global outlier detection.

higher number of supporting points and a larger sliding window to define outliers, the energy consumption of Global-KNN will be higher, but it is still more energy efficient than Centralized.

Balanced energy consumption is important for the longevity of a sensor network as an operating system. Figure 6 compares the minimum, average, and maximum energy consumption for a sensor node as  $w$  increases. This view further accentuates the advantage of using the Global-NN outlier detection solution over Centralized for large window sizes. In Centralized, some nodes consume by far more energy than others. This would lead, in a real application to the quick depletion of their energy. Since those nodes usually consume that much energy because they are critical to the algorithm, this would endanger the livelihood of the system. Another observation is that the range of energy consumption for different nodes running the same detection algorithm is larger for the centralized solution than for the distributed solution. Figure 7 illustrates this clearly by showing the values from Figure 6, normalized with respect to the average energy consumption. For  $w=10$ , the most energy consuming node consumed nearly three times more energy than the average node in a centralized algorithm and less than twice the energy of the average node in both distributed algorithms.

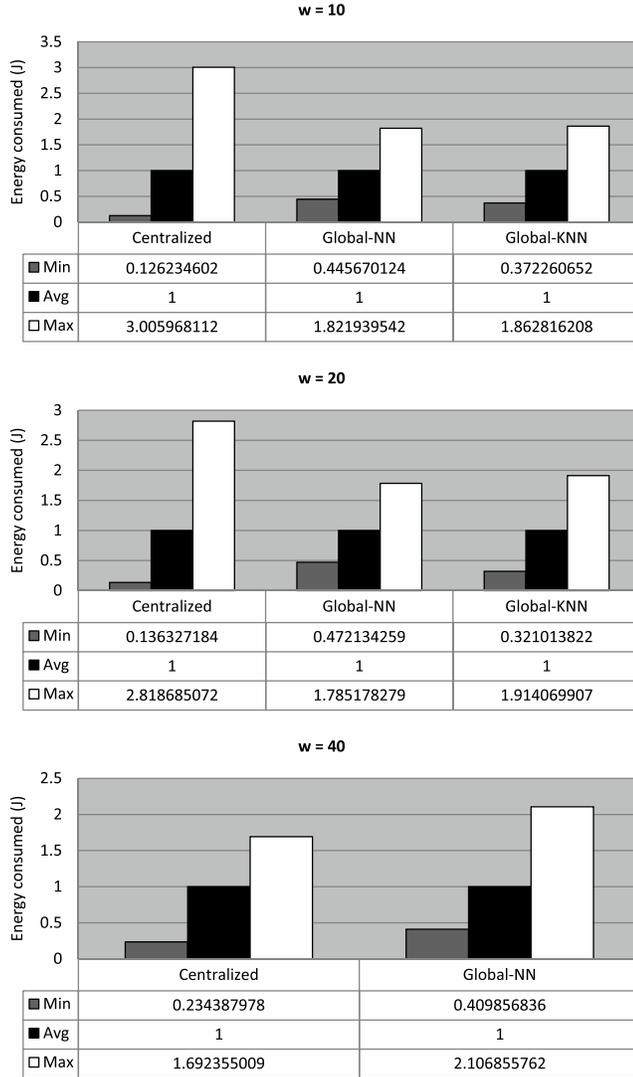
For the partial information for  $w=40$ , the normalized range of energy consumption is actually lower for the centralized algorithm than for the distributed one. However, referring back to Figure 6, we see that the average energy consumption for a node in



**Fig. 6** Average, minimum, and maximum energy consumption of a node for global outlier detection.

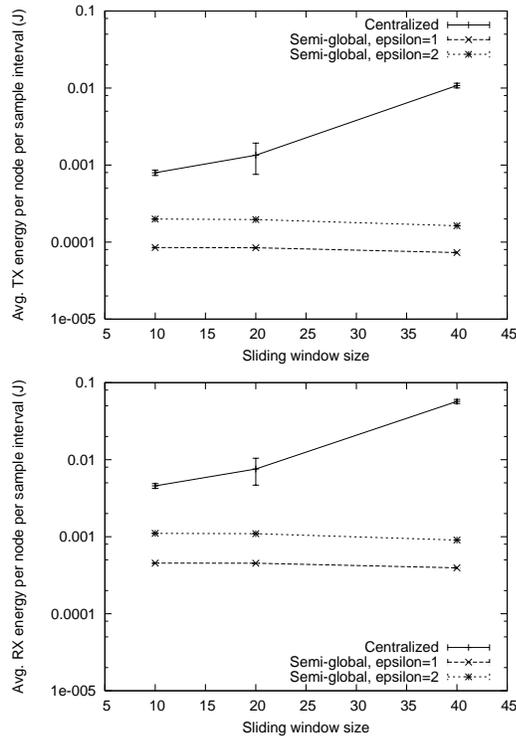
the centralized case is much higher than for the distributed case. Hence, in this case, the normalized maximum value does not convey the full picture of energy quality of the compared algorithms.

The plots in Figure 8 compare energy consumption rates for the centralized and distributed localized outlier detection algorithms. Since the results of using NN and KNN outlier detection methods are nearly identical, only results for the former are shown. Again, the centralized algorithm consumes much more energy per sample interval than the distributed algorithms. For the distributed localized algorithms, the



**Fig. 7** Normalized average, minimum, and maximum amount of energy consumed by a node for global outlier detection.

rate of energy consumption increases along with the values of epsilon (the hop range). This is expected because, as epsilon increases, so does the message passing overhead as data points travel farther from their place of origin. The behavior of the distributed algorithm in the localized case for nearest neighbor outlier detection is similar to that of global case for the same detection method. Energy consumption generally decreases as  $w$  increases. As before, we attribute this behavior to the increased data redundancy as the size of the sliding window increases. In general, the extent of the spatial area

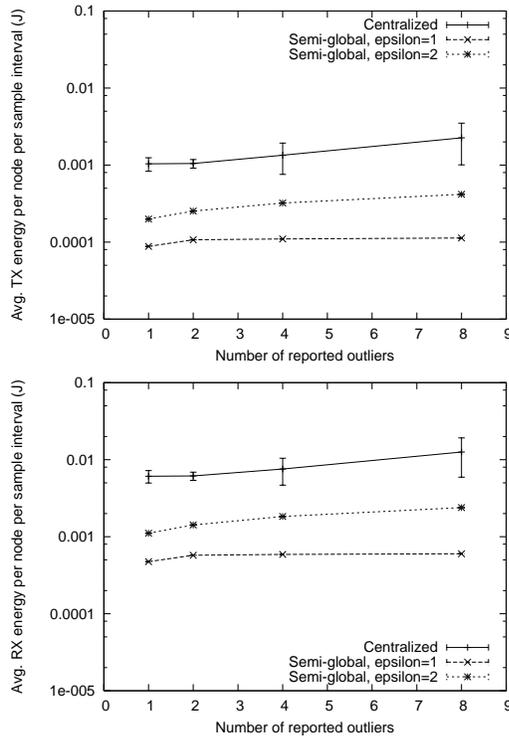


**Fig. 8** Average transmission and reception energy consumed per node per sample interval vs.  $w$  ( $n=4$ ) for localized outlier detection using nearest neighbor outlier detection.

over which outliers are defined affects the energy consumption trends of the algorithm, but not notably.

### 7.2.2 Effect of the number of reported outliers

We now investigate how the number of outliers produced affects energy consumption. Figure 9 illustrates the performance of the localized outlier detection algorithms under increasing values for  $n$  for KNN outlier detection. Similar plots for NN detection are omitted due to space restrictions and similarity of results; NN detection is negligibly less energy efficient most likely due to a lower rate of convergence. The energy consumption trends for these algorithms are straightforward and expected. Energy consumption increases along with both  $n$  and  $\epsilon$ , which both cause more message passing overhead with increasing value. We also noticed that the rate of increase was related to  $\epsilon$ . This is expected since the effects of larger  $\epsilon$  and  $n$  values on energy consumption compound one another.



**Fig. 9** Average transmission and reception energy consumption per node per sample interval vs.  $n$  ( $w=20$ ,  $k=4$ ) for localized outlier detection using  $k$  nearest neighbor outlier detection.

## 8 Conclusions

We addressed the problem of unsupervised outlier detection in WSNs. Our solution offers flexibility in the heuristic used to define outliers and computes the results in-network to reduce bandwidth and energy consumption. Its use of single-hop communication permits very simple node failure detection and message reliability assurance mechanisms. Dynamic data updates are also seamlessly accommodated by our method.

We evaluated the outlier detection algorithm's behavior on real-world sensor data using a simulated wireless sensor network. These initial results show promise for our algorithm in that it outperforms a strictly centralized approach under some very important circumstances. When the unabridged data from the entire sensor network are sent to a single location, the node collecting this data as well as its nearest neighbors become a bottleneck of the entire system. Indeed, the density of traffic in this region is proportional to the area of coverage of the entire network while the average node has traffic density proportional to the area covered by its communication range. In the example simulated in the paper, the traffic in the area of the collecting node was about 50 times more dense than in the other parts of the network. The immediate consequence can be shorter system lifetime, as the battery operated nodes closest to the collection point will die of energy exhaustion while the nodes farthest from the collection point will still have a lot of energy left. This is simply the result of uneven

spatial distribution of paths through which data are collected. If all nodes, except the central node, are battery operated, the least used nodes will have only 2% of energy used when the most used nodes will start dying. The second consequence of the imbalance is traffic congestion, which leads to interferences and message-collision. Fewer messages are accepted the first time they are transmitted and the algorithm will either degrade the quality of the outcome or would be forced to expand greater effort, costlier protocols, and more energy, to make sure messages are eventually delivered. In short, using the centralized algorithm with its inordinately imbalanced traffic density will put even the best routing protocols under severe stress. Our distributed and localized outlier detection algorithms avoid these difficulties.

Our approach is well suited for applications in which the confidence of an outlier rating may be calculated by either an adjustment of sliding window size or by the number of neighbors used in a distance-based outlier detection technique. These applications are critical for resource-constrained sensor networks for two reasons. First, communication is a costly activity: only the most accurate data should be transmitted to a client application. Second, emerging safety-critical applications that utilize wireless sensor networks will require the most accurate data, including outliers. This work represents our contribution toward enabling efficient data cleaning solutions for these types of applications.

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## A Appendix: Correctness Proofs for the Global Outlier Detection Algorithm

In this section, we provide detailed proofs of Theorems 1 and 2. Before doing so, a few technical lemmas are needed. The first two isolate a couple of useful properties following from the axioms of  $R$ .

**Lemma 1** For any  $P \subseteq Q \subseteq D$  where  $|P| \geq n$ , if  $O_n(P) \neq O_n(Q)$ , then there exists  $x \in O_n(P)$  such that  $R(x, P) > R(x, Q)$ .

*Proof* Assume  $O_n(P) \neq O_n(Q)$ . Since  $|O_n(P)| = |O_n(Q)| = n$ , then there exists  $x \in (O_n(P) \setminus O_n(Q))$  and  $y \in (O_n(Q) \setminus O_n(P))$ . Recall that we assume a tie-breaking mechanism is used to ensure  $R(\cdot, P)$  and  $R(\cdot, Q)$  are one-to-one. Thus, by definition of  $O_n(\cdot)$  it follows that  $R(x, P) > R(y, P)$  and  $R(y, Q) > R(x, Q)$ . The anti-monotonicity axiom implies  $R(y, P) \geq R(y, Q)$ , yielding the desired result.  $\square$

**Lemma 2** For any  $P \subseteq D$ ,  $x \in O_n(P)$ , and  $z \in P$ , we have  $R(x, P) = R(x, [P|O_n(P)]) = R(x, [P|O_n(P)] \cup \{z\})$ .

*Proof* Since, by definition,  $[P|x] \subseteq [P|O_n(P)] \subseteq ([P|O_n(P)] \cup \{z\}) \subseteq P$ , then by the anti-monotonicity axiom it follows that

$$\begin{aligned} R(x, P) &= R(x, [P|x]) \\ &\geq R(x, [P|O_n(P)]) \\ &\geq R(x, [P|O_n(P)] \cup \{z\}) \\ &\geq R(x, P). \end{aligned}$$

$\square$

The last technical lemma shows that once a sensor  $p_i$  completes its local computation, then  $(D_{i,j}^i \cup D_{j,i}^i)$  contains a particular crucial set of points (among others) needed for consistency among sensors' outlier estimates.

**Lemma 3** For any  $p_i$ , once the main for-loop in the algorithm completes,  $[P_i|O_n(D_{i,j}^i \cup D_{j,i}^i)] \subseteq (D_{i,j}^i \cup D_{j,i}^i)$ .

*Proof* Let  $D_{i,j}^i$  (*before*) denote the set of points held by  $p_i$  and sent from  $p_i$  to  $p_j$  immediately before the execution of the ‘‘Repeat until no change: ...’’ step in the main for loop for  $j \in \Gamma_i$ . For  $\ell \geq 1$ , let  $Z_j(\ell)$  denote  $Z_j$  immediately before the  $\ell^{\text{th}}$  iteration in the execution of the ‘‘Repeat until no change ...’’ step in the main for loop for  $j \in \Gamma_i$ , e.g.,  $Z_j(1) = O_n(P_i) \cup [P_i|O_n(P_i)]$ .

By definition,  $Z_j(\ell) \subseteq Z_j(\ell+1) \subseteq P_i$  and  $P_i$  is finite. Thus, let  $\ell^*$  denote the smallest integer such that  $Z_j(\ell^* - 1) = Z_j(\ell^*)$ . Hence, the ‘‘Repeat until no change ...’’ step terminates at the end of iteration  $\ell^*$  and  $Z_j = Z_j(\ell^*)$  in the remainder of the main for loop. Therefore,

$$Z_j(\ell^*) = Z_j(\ell^*) \cup [P_i|O_n(D_{i,j}^i(\text{before}) \cup D_{j,i}^i \cup Z_j(\ell^*))]$$

and

$$D_{i,j}^i \cup D_{j,i}^i = D_{i,j}^i(\text{before}) \cup D_{j,i}^i \cup Z_j(\ell^*).$$

It follows that  $[P_i|O_n(D_{i,j}^i \cup D_{j,i}^i)] \subseteq Z_j(\ell^*) \subseteq (D_{i,j}^i \cup D_{j,i}^i)$ .  $\square$

Now we prove that upon termination of the algorithm, the sensors' estimates are consistent.

*Theorem 1* Assuming a connected network, if for all sensors  $p_i$ ,  $D_i$  and  $\Gamma_i$  do not change, then upon termination of the algorithm all sensors' outlier estimates and supports agree. I.e., for all  $p_i, p_j$ , (i)  $O_n(P_i) = O_n(P_j)$  and (ii)  $[P_i|O_n(P_i)] = [P_j|O_n(P_j)]$ .

*Proof* Since the network is connected, we may assume, without loss of generality, that  $p_i$  and  $p_j$  are neighbors. To prove part (i), we will show that  $O_n(P_i) = O_n(D_{i,j}^i \cup D_{j,i}^i) = O_n(D_{i,j}^j \cup D_{j,i}^j) = O_n(P_j)$ . The middle equality follows from the fact that  $(D_{i,j}^i \cup D_{j,i}^i) = (D_{i,j}^j \cup D_{j,i}^j)$ . By symmetry, it suffices to show the first equality.

Suppose  $O_n(P_i) \neq O_n(D_{i,j}^i \cup D_{j,i}^i)$ . The following contradiction is reached. There exists  $x \in O_n(D_{i,j}^i \cup D_{j,i}^i)$  such that

$$\begin{aligned}
R(x, D_{i,j}^i \cup D_{j,i}^i) &> R(x, P_i) \\
&= R(x, [P_i|x]) \\
&\geq R(x, [P_i|O_n(D_{i,j}^i \cup D_{j,i}^i)]) \\
&\geq R(x, D_{i,j}^i \cup D_{j,i}^i).
\end{aligned}$$

The first inequality follows from Lemma 1 (with  $P = (D_{i,j}^i \cup D_{j,i}^i)$  and  $Q = P_i$ ). The equality follows from the definition of support  $[\cdot|x]$ . The last two inequalities follow from the anti-monotonicity of  $R$  and Lemma 3.

To prove part (ii)  $[P_i|O_n(P_i)] = [P_j|O_n(P_j)]$ , it suffices to show that for any  $x \in O_n(D_{i,j}^i \cup D_{j,i}^i)$ , it is the case that  $[P_i|x] = [P_j|x]$ . This is because  $O_n(P_i) = O_n(D_{i,j}^i \cup D_{j,i}^i) = O_n(P_j)$ . We will prove that  $[P_i|x] = [P_i \cap P_j|x] = [P_j|x]$ . By symmetry it is enough to show the first equality.

From Lemma 3 it follows that

$$\begin{aligned}
[P_i|x] &\subseteq [P_i|O_n(D_{i,j}^i \cup D_{j,i}^i)] \\
&\subseteq (D_{i,j}^i \cup D_{j,i}^i) \\
&\subseteq (P_i \cap P_j).
\end{aligned}$$

Thus, anti-monotonicity implies  $R(x, P_i) \geq R(x, [P_i|x]) \geq R(x, P_i \cap P_j) \geq R(x, P_i)$ , and so,

$$R(x, P_i) = R(x, [P_i|x]) = R(x, P_i \cap P_j).$$

Therefore,  $[P_i|x]$ ,  $[P_i \cap P_j|x]$  are support sets of  $x$  with respect to  $P_i \cap P_j$  and  $P_i$ . Since  $[P_i|x]$  ( $[P_i \cap P_j|x]$ ) is the *unique* smallest support set for  $x$  with respect to  $P_i$  ( $P_i \cap P_j$ ), then it follows that  $[P_i|x] = [P_i \cap P_j|x]$ .  $\square$

Finally, we prove that upon termination the sensors' estimates are equal to the correct answer.

*Theorem2* Assuming a connected network, if for all sensors  $p_i$ ,  $D_i$  and  $\Gamma_i$  do not change, then upon termination of the algorithm, all sensors' outlier estimate will be correct. I.e., for all  $p_i$ ,  $O_n(P_i) = O_n(D)$ .

*Proof* Suppose there exists a sensor  $p_i$  such that  $O_n(P_i) \neq O_n(D)$ . By Lemma 1 (with  $P = P_i$  and  $Q = D$ ), there exists  $x \in O_n(P_i)$  such that  $R(x, P_i) > R(x, D)$ . Moreover, the first equality in Lemma 2 (with  $P = P_i$ ) implies that  $R(x, [P_i|O_n(P_i)]) = R(P_i, x)$ .

Since  $R(x, [P_i|O_n(P_i)]) > R(x, D)$ , then the smoothness axiom (with  $Q_1 = [P_i|O_n(P_i)]$  and  $Q_2 = D$ ), implies there exists  $z \in (D \setminus [P_i|O_n(P_i)])$  such that

$$R(x, [P_i|O_n(P_i)]) > R(x, [P_i|O_n(P_i)] \cup \{z\}).$$

This point  $z$  must be contained in  $P_j$  for some sensor  $p_j$ . Hence, the following inequality is reached.

$$\begin{aligned}
R(x, [P_i|O_n(P_i)]) &> R(x, [P_i|O_n(P_i)] \cup \{z\}) \\
&= R(x, [P_i|O_n(P_j)] \cup \{z\}) \\
&= R(x, [P_j|O_n(P_j)]) \\
&= R(x, [P_i|O_n(P_i)]).
\end{aligned}$$

The inequality above leads to a contradiction: The first equality follows from Theorem 1 part (i). The middle equality follows from the second equality of Lemma 2 (with  $P = P_j$  and noting that  $O_n(P_j) = O_n(P_i)$  by Theorem 1 part (i)). The last equality follows from Theorem 1 part (ii).  $\square$

Transmission energy consumption per node per sample interval (J)						
w (n=4, k=4)	Centralized		Global-NN		Global-KNN	
	Avg.	Confidence interval	Avg.	Confidence interval	Avg.	Confidence interval
10	7.95e-04	±6.14e-05	0.002381	±1.66e-04	0.001024	±3.48e-05
20	0.001343	±5.85e-04	0.001331	±5.25e-05	0.001545	±3.12e-05
40	0.010793	±7.54e-04	0.00103	±5.82e-05	na	na
Receive energy consumption per node per sample interval (J)						
w (n=4, k=4)	Centralized		Global-NN		Global-KNN	
	Avg.	Confidence interval	Avg.	Confidence interval	Avg.	Confidence interval
10	0.004571	±3.29e-04	0.013175	±9.47e-04	0.005645	±1.94e-04
20	0.007566	±0.002894	0.007334	±2.89e-04	0.008564	±1.78e-04
40	0.057144	±0.003496	0.005746	±3.27e-04	na	na
Average energy consumption per node (J)						
w (n=4, k=4)	Centralized		Global-NN		Global-KNN	
	Avg.	Confidence interval	Avg.	Confidence interval	Avg.	Confidence interval
10	0.094	±0.006621	0.267191	±0.018928	0.116109	±0.00388
20	0.154209	±0.059139	0.149974	±0.005806	0.174561	±0.003549
40	1.157647	±0.072147	0.11793	±0.006546	na	na
Minimum energy consumption by a node (J)						
w (n=4, k=4)	Centralized		Global-NN		Global-KNN	
	Avg.	Confidence interval	Avg.	Confidence interval	Avg.	Confidence interval
10	0.011866	±9.71e-04	0.119079	±0.005318	0.043223	±0.00153
20	0.021023	±0.008168	0.070808	±0.004767	0.056036	±0.002701
40	0.271339	±0.055955	0.048334	±0.002853	na	na
Maximum energy consumption by a node (J)						
w (n=4, k=4)	Centralized		Global-NN		Global-KNN	
	Avg.	Confidence interval	Avg.	Confidence interval	Avg.	Confidence interval
10	0.28256	±0.023512	0.486806	±0.040617	0.216289	±0.008319
20	0.434666	±0.138827	0.26773	±0.010854	0.334121	±0.00938
40	1.959149	±0.151924	0.248462	±0.015394	na	na

Transmission energy consumption per node per sample interval (J)						
w (n=4, k=4)	Centralized		Semi-global, epsilon=1		Semi-global, epsilon=2	
	Avg.	Confidence interval	Avg.	Confidence interval	Avg.	Confidence interval
10	7.95e-04	±6.14e-05	8.49e-05	±9.22e-07	1.99e-04	±8.97e-07
20	0.001343	±5.85e-04	8.44e-05	±9.22e-07	1.96e-04	±1.07e-06
40	0.010793	±7.54e-04	7.32e-05	±9.22e-07	1.63e-04	±1.13e-06
Receive energy consumption per node per sample interval (J)						
w (n=4, k=4)	Centralized		Semi-global, epsilon=1		Semi-global, epsilon=2	
	Avg.	Confidence interval	Avg.	Confidence interval	Avg.	Confidence interval
10	0.004571	±3.29e-04	4.54e-04	±5.05e-06	0.001107	±4.73e-06
20	0.007566	±0.002894	4.52e-04	±5.03e-06	0.001096	±6.32e-06
40	0.057144	±0.003496	3.94e-04	±5.05e-06	9.04e-04	±6.15e-06

## B Appendix: Experimental Results Tables

The numbers (averages, confidence intervals) that were used to generate the experimental results plots shown earlier are contained in the following tables.

Transmission energy consumption per node per sample interval (J)						
w (n=4, k=4)	Centralized		Semi-global, epsilon=1		Semi-global, epsilon=2	
	Avg.	Confidence interval	Avg.	Confidence interval	Avg.	Confidence interval
10	7.95e-04	6.14e-05	1.01e-04	9.22e-07	3.06e-04	8.97e-07
20	0.001343	5.85e-04	1.10e-04	9.22e-07	3.22e-04	1.80e-06
40	0.010793	7.54e-04	1.24e-04	9.22e-07	3.12e-04	3.82e-06
n (w=20, k=4)	Centralized		Semi-global, epsilon=1		Semi-global, epsilon=2	
	Avg.	Confidence interval	Avg.	Confidence interval	Avg.	Confidence interval
1	0.001037	2.07e-04	8.79e-05	9.22e-07	1.99e-04	1.36e-06
2	0.001048	1.34e-04	1.07e-04	9.22e-07	2.53e-04	1.81e-06
4	0.001343	5.85e-04	1.10e-04	9.22e-07	3.22e-04	1.80e-06
8	0.002256	0.00125	1.13e-04	9.22e-07	4.17e-04	2.85e-06
Receive energy consumption per node per sample interval (J)						
w (n=4, k=4)	Centralized		Semi-global, epsilon=1		Semi-global, epsilon=2	
	Avg.	Confidence interval	Avg.	Confidence interval	Avg.	Confidence interval
10	0.004571	3.29e-04	5.40e-04	5.02e-06	0.001724	4.66e-06
20	0.007566	0.002894	5.89e-04	5.06e-06	0.001832	1.13e-05
40	0.057144	0.003496	6.67e-04	4.79e-06	0.001761	2.29e-05
n (w=20, k=4)	Centralized		Semi-global, epsilon=1		Semi-global, epsilon=2	
	Avg.	Confidence interval	Avg.	Confidence interval	Avg.	Confidence interval
1	0.006097	0.00112	4.74e-04	5.04e-06	0.00111	8.21e-06
2	0.006145	7.50e-04	5.75e-04	4.94e-06	0.001422	1.12e-05
4	0.007566	0.002894	5.89e-04	5.06e-06	0.001832	1.13e-05
8	0.012555	0.006633	6.01e-04	4.96e-06	0.002388	1.74e-05