

Zhang, X.-Y., Gong, Y.-j., Zhan, Z., Chen, W.-N., Li, Y., and Zhang, J. (2015) Kuhn-Munkres parallel genetic algorithm for the set cover problem and its application to large-scale wireless sensor networks. IEEE Transactions on Evolutionary Computation.

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Deposited on: 04 May 2016

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# Kuhn-Munkres Parallel Genetic Algorithm for the Set Cover Problem and Its Application to Large-Scale Wireless Sensor Networks

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Abstract-Operating mode scheduling is crucial for the lifetime of wireless sensor networks. However, the growing scale of networks has made such a scheduling problem more and more challenging, as existing set cover and evolutionary algorithms become unable to provide satisfactory efficiency due to the curse of dimensionality. In this paper, a Kuhn-Munkres parallel genetic algorithm is developed to solve the set cover problem and is applied to lifetime maximization of large-scale wireless sensor networks. The proposed algorithm schedules the sensors into a number of disjoint complete cover sets and activates them in batch for energy conservation. It uses a divide-and-conquer strategy of dimensionality reduction, and the polynomial Kuhn-Munkres algorithm are hence adopted to splice the feasible solutions obtained in each subarea to enhance the search efficiency substantially. To further improve global efficiency, a redundant-trend sensor schedule strategy is developed. Additionally, we meliorate the evaluation function through penalizing incomplete cover sets, which speeds up convergence. Eight types of experiments are conducted on a distributed platform to test and inform the effectiveness of the proposed algorithm. The results show that it offers promising performance in terms of the convergence rate, solution quality, and success rate.

*Index Terms*—parallel genetic algorithm, set cover problem, large-scale wireless sensor networks, Kuhn-Munkres algorithm.

#### I. INTRODUCTION

WIRELESS sensor networks (WSNs) have been widely used in a number of fields to satisfy various requirements, such as road traffic monitoring [1], environmental observation [2], healthcare sensing [3], and asset monitoring [4]. Typically, hundreds or even thousands of sensors, each with a series of transceivers, a battery and a micro central processing unit, are deployed in a target area. Since it is impossible to recharge or replace the battery in some scenarios, how to extend the lifetime of WSNs becomes a critical task [5].

Manuscript received ...; revised ...; accepted ... This work was partially supported by the National Natural Science Foundation of China (NSFC) Key Project No. 61332002 and by the NSFC Youth Projects No. 61502542 and No. 61300044 (Y.-J. Gong and J. Zhang are equally contributed corresponding authors, email: gongyuejiao@gmail.com; junzhang@ieee.org).

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Existing ways for lifetime enhancement are classified into five categories: operation mode control [6], data processing [7][8], sink relocation [9]-[11], topology control [12][13], and optimal routing [14]-[16]. There are various definitions of the network lifetime. In this paper, the lifetime of a WSN refers to the duration of time that the network is able to carry out its set mission. Normally, the networks can fulfill its mission if it can guarantee the specified coverage requirements by the sensors deployed, i.e., the set cover condition is satisfied [17].

As summarized in [18]-[20], the deployment methods for sensors in WSNs vary with applications, which can be categorized into deterministic deployment and random deployment. Deterministic deployment is applied to a smallor medium-scale network in a friendly sensory environment [21]-[23]. The set cover problem here can be transformed into a minimum set cover problem or its dual problem. There are some certain theoretical developments [24]-[26] and optimization algorithms [18][21] related to this field. Since this problem is NP-hard, evolutionary-computation based solvers are potentially promising because of their powerfulness in dealing with NP-hard problems. However, the optimal number of sensors cannot be known in advance, which increases the difficulty of applying an evolutionary algorithm, such as the genetic algorithm (GA). The variable length chromosome puts a great challenge to the crossover operation of GA. Nevertheless, this issue has been well solved recently by using a bi-objective GA [27].

When the environment is inaccessible or unfriendly, or the number of sensors is too large, sensors are often scattered from an aircraft or by other means of transportation, which in effect results in random deployment. In order to guarantee coverage and connectivity, the sensors are to be densely deployed in target areas. To construct an energy-efficient WSN in this case, sensors are assigned to different cover sets independent of one another [28]. Activating them in batch ensures that only one cover set is active at a time and the others are scheduled to sleep. This scientific problem is known as the *Set K-cover* problem [29] or *Disjoint Set Covers* problem [30], which is a nondeterministic polynomial complete (NP) problem and hence its optimization is NP-hard.

A common objective of solving the *Set K-cover* problem is to maximize the lifetime of the WSN, but they differ slightly in terms of coverage constraints. In [31], sensors are aimed to be scheduled into K disjoint sets while guaranteeing that the coverage ratio of each set is as high as possible by modeling

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the *Set K-cover* problem as an *N*-person card game, and solutions are obtained after a gaming process. A heuristic method, termed the Most Constrained-Minimally Constraining Covering (MCMCC), is proposed in [29]. The essence of MCMCC is to minimize the coverage of sparsely covered areas within one cover set. It requires that each cover set is able to cover the target area completely. We focus on complete cover sets in this paper.

Owing to their success in solving nondeterministic polynomial problems, GAs [32]-[34] and other evolutionary algorithms (EAs) [35]-[37] have been applied to the lifetime problem in WSNs recently. Lai et al. [38] propose a GA for maximum disjoint set covers (GAMDSC), which applies a scattering operator to the EA offspring to keep critical sensors from joining the same cover set. Hu et al. [39] propose a schedule transition hybridized genetic algorithm (STHGA), which adopts a forward encoding scheme for chromosomes and utilizes redundancy information via designing a series of transition operations. Ant-colony optimization for maximizing the number of connected cover (ACO-MNCC) is proposed in [40], which maximizes the lifetime of heterogeneous WSNs. These algorithms are shown competitive in solving small to medium sized WSNs. However, the performance of existing methods decreases substantially when dealing with a large-scale Set K-cover Set problem.

To improve set cover efficiency and large-scale WSN performance over existing algorithms, a Kuhn-Munkres scheduled parallel GA (KMSPGA) is developed in this paper, so as to provide the following features and benefits:

- A divide-and-conquer strategy to achieve dimensionality reduction in a simple but effective way and solve the small-scale *Set K-cover problem* separately in each subarea;
- The merging of local feasible solution is modeled as a Maximum Weight Perfect Matching (MWPM) problem such that Kuhn-Munkres (KM) algorithm [41][42] is applicable;
- A redundant-trend sensor schedule strategy (RTSS) to further improve global search efficiency, which is easier to implement than the existing auxiliary schedule strategies [39], and
- A modified fitness index by introducing a coverage inadequate penalty to guide convergence better.

The rest of this paper is organized as follows. In Section II, we describe the *Set K-cover* problem for WSNs, with assumptions and definitions given. In Section III, we develop KMSPGA in detail. This parallel genetic algorithm is comprehensively tested for performance through simulations in Section IV. Finally, conclusions are drawn in Section V.

## II. PROBLEM FORMULATION

# A. Sensor Model

In this section, we introduce the sensor model adopted in our algorithm, which is crucial for the coverage, connectivity, and energy consumption.



Fig. 1. Boolean Disk model. Considering two points  $z_1$  and  $z_2$  in the target area, we have  $f(D(s, z_1)) = 1$  and  $f(D(s, z_2)) = 0$ .

A sensor coverage model is an abstract concept to measure the sensing capability and to quantify how well a sensor can monitor the occurrence of events within its sensing range. Various models have been proposed, such as the Boolean sector coverage model and the attenuated disk coverage model [43]. Comparisons of existing models are made in [44]-[46].

In this paper, we adopt the Boolean disk model because of its concise definition and wide applicability [19][43], which is illustrated in Fig. 1. This model ignores the dependency of environmental conditions. A point in the space of the target area is considered to be covered only if it is within the range of at least one sensor. A coverage function is formalized as:

$$f(D(s,z)) = \begin{cases} 1, & \text{if } D(s,z) \le R_s \\ 0, & otherwise \end{cases}$$
(1)

$$D(s,z) = \sqrt{(s_x - z_x)^2 + (s_y - z_y)^2}$$
(2)

where  $s = (s_x, s_y)$  is the central coordinate of a sensor and  $(z_x, z_y)$ is the coordinate of point z in a 2-dimensional space,  $R_s$  is the sensing radius and D(s, z) calculates the Euclidean distance between s and z. Once a point is covered by a sensor, the coverage function value is 1, otherwise, it is 0. In [47], a geometric analysis of the relationship between coverage and connectivity is provided, showing that the connectivity can be guaranteed inherently if 1) coverage is satisfied and 2) the communication range of sensors is no less than twice of the sensing range. Afterwards, many related publications for solving the Set K-cover problems are based on this proof [31][39] that they use the sensors with communication range at least twice than the sensing range and consider the coverage constraint only. In this paper, we also make such an assumption and meet the connectivity constraint by satisfying the coverage constraint.

In WSNs, the energy consumption of a specific sensor can be determined by the following aspects: 1) the communication when sending and receiving data; 2) the data type and size acquired for processing; 3) the differences in the inherent characteristic of the battery; 4) the different local working condition, such as temperature. These are, however, diverse and highly related to the application scene. In this paper, we focus on the Set K-cover problem. Instead of fully formulating the above issues influencing the energy consumption, we simplify the model by making the following three assumptions: 1) for every active sensor, the amount of energy consumption is identical per unit time; 2) for every sleep sensor, they consume negligible energy compared with the active ones; 3) every sensor carries the same amount of energy in the initial stage. Based on these assumptions, the lifetime of each complete cover set is identical to each one another. Nevertheless, it is to be noted that the energy consumption model in use have no direct relationship with the operations or parameters of our proposed KMSPGA. In practical application,

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there is still room for KMSPGA to adapt this framework to fitting new forms of energy constraints.

## B. Set K-Cover Problem

In this paper, we focus on maximizing the lifetime of a large-scale WSN using a static optimization strategy. The proposed KMSPGA is an off-line algorithm that pre-calculates K disjoint complete cover sets at a time. By alternatively activating the K cover sets in batch, the lifetime of the WSN can be K times larger than the lifetime of a single cover set. Hence, maximizing the lifetime of the WSN is equivalent to finding the maximum disjoint complete cove sets.

We consider the *Set K-cover* problem a scheduling problem. For a positive integer K, sensors are judiciously scheduled into K disjoint cover sets such that each cover set is able to meet the coverage requirement. In this paper, we focus on a complete coverage. The premise of finding K disjoint complete cover sets is that every element of the target area is covered by at least K sensors. Sufficiency of this premise is proven in the following of this subsection. Fig. 2 illustrates the concept of an element and a K-covered element. In Fig. 2 (a), the grey square area is divided into eight elements. Element  $E_x$  in Fig. 2 (b) is covered by four sensors; hence it is called 4-covered.

Without loss of generality, assume that a target area  $\Gamma$  is a rectangle and that *N* sensors  $S_1, S_2, ..., S_N$  are randomly deployed in  $\Gamma$ . A constraint of the coverage requirements is a complete coverage. Given a positive number *K*, sensors are scheduled into *K* cover sets  $C = \{C_1, C_2, ..., C_K\}$ . For each cover set  $C_i$  ( $i = \{1, 2, ..., K\}$ ), if every element of  $\Gamma$  is covered by at least one sensor in  $C_i$ , then *C* would be considered as a feasible solution of the *Set K-cover* problem. The *Set K-cover* problem can be formalized as:

$$\bigcup_{\forall S_k \in C_i} E(S_k) \supseteq \Gamma \tag{3}$$

$$\bigcup_{i=1}^{K} C_i \subseteq S \tag{4}$$

$$C_i \cap C_j = \emptyset, \ i \neq j, \ i, j \in \{1, 2, \dots, K\}$$
(5)

where  $E(S_k)$  represents the element sensed by sensor  $S_k$ , k is the sensor index, S is the collection of sensors. Assume that the target area is partitioned into M elements  $E_1, E_2, ..., E_M$ . To make a clear explanation of how to calculate the upper bound of K, we first give the following proposition and its proof:

*Proposition* 1: The prerequisite of finding *K* disjoint cover sets is that each element is covered by at least *K* sensors.

*Proof*: Assume that  $E_{\tau}$  is covered by Q sensors  $SC_{\tau} = \{S_{\tau,1}, S_{\tau,2}, ..., S_{\tau,Q}\}$ , where  $SC_{\tau}$  is the collection of sensors covering element  $E_{\tau}$ . There still exist K disjoint complete cover sets with  $Q \le K$ .

Let *C* be a feasible solution of the *Set K-cover* problem. Then, we have |C| = K. Considering the same element  $E_{\tau}$  in the assumption,  $E_{\tau}$  is expected to be covered in each  $C_i \in C$  according to (3). Then, *K* cover sets are considered as *K* covering tasks, and we have *Q* sensors that can perform this task. Then, *K* tasks are assigned to *Q* sensors. Since we have  $|C| = K > |SC_{\tau}| = Q$ , there exist cover sets  $C_p$ ,  $C_q \in C$ , and sensor  $S_{\tau,m} \in SC_{\tau}$  ( $m \in \{1, 2, ..., Q\}$ ) satisfying  $S_{\tau,m} \in C_p \cap C_q$  according to the drawer principle, and therefore it contradicts

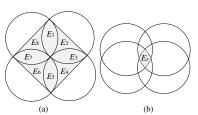


Fig. 2. Notion of an element and a *K*-covered element. (a) The shaped area is divided into eight elements  $E_1, E_2, ..., E_8$ , where four elements are covered by two sensors and the remaining four by one only. (b) Element  $E_x$  is called *K*-covered, if it is covered by *K* sensors, in which case K = 4.

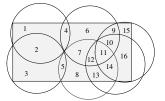


Fig. 3. Elements  $E_1, E_3, E_6, E_8$ , and  $E_{15}$  are called critical elements, as they are covered by the minimum number of sensors.

(5). In conclusion, the assumption is invalid and hence the proposition is proven to be tenable.

In order to better explain how to estimate the upper limit of K, the notion of critical element and critical sensor is introduced as follows. The target area is partitioned into a number of elements by thousands of densely deployed sensors. An element covered by the minimum number of sensors is called a 'critical element' and the corresponding sensors 'critical sensors'. Fig. 3 illustrates critical elements and critical sensors, where six sensors divide the rectangular area into sixteen elements  $E_1$ - $E_{16}$ . Being covered by one sensor only, elements  $E_1$ ,  $E_3$ ,  $E_6$ ,  $E_8$ , and  $E_{15}$  are critical elements.

Let  $E_c$  be a critical element. Assume that the number of sensors covering  $E_c$  is  $\hat{U}$ . According to *Proposition* 1, we can at most find  $\hat{U}$  disjoint complete cover sets only if the  $\hat{U}$  critical sensors are chosen in  $\hat{U}$  disjoint cover sets which guarantee that  $E_c$  is covered by every cover set, and therefore  $\hat{U}$  is regarded as the upper limit of K.

#### C. Critical Parameters

In this subsection, we discuss some critical parameters related to a large-scale WSN. A redundant rate represents the density of sensors deployed in the target area. The redundant rate in the 2D ideal plane model is computed as (6) according to [39], where  $area(\Gamma)$  is the area of  $\Gamma$  and N is the number of sensors.

$$\eta = \frac{N \cdot \pi \cdot R_s^2}{\hat{U} \cdot area(\Gamma)} \tag{6}$$

It is difficult to compute the coverage ratio of  $\Gamma$  accurately when applying the Boolean disk model. For this reason, we divide the target area into *T* smaller square grids  $(g_1, g_2, ..., g_T)$ , *T* being computed as (7), where *d* is the width of the grid. A coverage ratio is defined as (8), where  $N_g(S_i)$  represents the collection of grids covered by sensor  $S_i$  and  $|N_g(S_i)|$  is the number of grids in  $N_g(S_i)$ . Equation (9) indicates that a grid belongs to no more than one collection in case of a repeat count.

$$T = \frac{area(\Gamma)}{d^2} \tag{7}$$

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$$\delta = \frac{1}{T} \sum_{i=1}^{N} |N_g(S_i)| \tag{8}$$

$$N_g(S_i) \cap N_g(S_j) = \emptyset, \quad i \neq j, i, j \in \{1, 2, \dots, N\}$$
(9)

The coverage criteria stipulates that grid  $g_j$  is covered by sensor  $S_i$  only if all its four vertices are within the sensing range of  $S_i$ . Fig. 4 shows how to calculate the coverage ratio. This calculation method is widely used to estimate the coverage ratio [39][40]. However, the grid width can influence the computation of the coverage ratio in terms of computational complexity and accuracy. Fig. 5 gives an example to explain this special case, where the grey grid is apparently covered by the WSN. Unfortunately, it is regarded as uncovered due to the above coverage criteria whether a grid is covered. If the width of this grid is halved, two resultant grids become covered. However, the calculation of the coverage ratio is of an  $O(N \times T)$  computational complexity. A shorter width means a higher computational complexity according to (7).

In our work, we adopt a simple strategy to determine the grid width in a preprocessing step. Given  $\Psi$  kinds of  $d_i$  ( $i = \{1, 2, ..., \Psi\}$ ,  $d_i < d_{i+1}$ ) in the process of estimating the upper limit of K, we choose the smallest  $d_1$  first to obtain an exact value  $\hat{U}_1$ , because  $d_1$  is small enough to guarantee accuracy. Then  $d_i$  ( $i = \{2, 3, ..., \Psi\}$ ) is adopted to work out  $\hat{U}_i$  in sequence. The largest  $d_i$  (ensuring  $\hat{U}_i = \hat{U}_1$ ) is used for calculating the coverage ratio. Algorithm 1 presents a set of pseudocode of this preprocess of choosing a proper grid width. In this paper, we adopt 5 kinds of grid widths: ( $d_1$ ,  $d_2$ ,  $d_3$ ,  $d_4$ ,  $d_5$ ) = (0.625, 0.78125, 1, 1.25, 1.5625).

#### III. PROPOSED PARALLEL GENETIC ALGORITHM

## A. Kuhn-Munkres Parallel Genetic Approach

KMSPGA is designed on a divide-and-conquer strategy, and the polynomial KM algorithm is adopted to splice the feasible solutions obtained in each subarea. The framework of KMSPGA is shown in Fig. 6. In the first step, we uniformly divide the target area into a number of subareas and encode them. Assuming that the number of sensors within subarea  $A_i$ is  $N_i$ , it satisfies:

$$\sum_{i=1}^{L \times W} N_i = N \tag{10}$$

where L and W denote the number of partitions along horizontal and vertical directions respectively, and hence  $L \times$ W denotes the number of subareas obtained. Therefore, the Set *K*-cover problem size of subarea  $A_i$  is whittled down to  $N_i$ . After the partition process, the small-scale Set K-cover problem within each subarea is separately solved by a parallel processing module. When local solutions of each small-scale Set K-cover problem reach a predefined state, they are spliced through a KM combination operation to achieve global optimization efficiently. There are two termination conditions: 1) FEs reaches its upper limit and/or 2) the number of complete cover sets reach Û. KMSPGA deploys a termination-controller in its master process. The controller checks whether the current process has reached the termination condition at the end of every generation. If so, the master process broadcasts a termination signal to all the slave

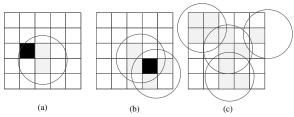


Fig. 4. Calculation of the coverage ratio. (a) The light grey grids are covered while the black one is not covered because one of its vertices is beyond the sensing range of the sensor. (b) The black grid is covered by two sensors  $S_p$  and  $S_q$ , but only belongs to either of  $N_g(S_i)$ ,  $i \in \{p, q\}$ , in the case of a repeat count. (c) The coverage ratio is 11/25 = 0.44.

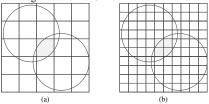


Fig. 5. An example to show how the grid width influences the computation of the coverage ratio. (a) The grey grid is regarded as uncovered due to the coverage criteria. (b) The grey grid becomes covered by either of the two sensors after the grid width is shortened.

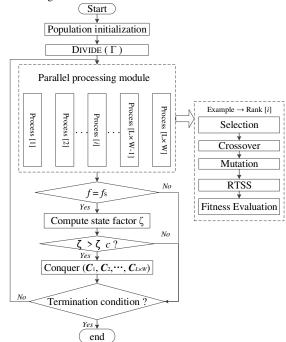


Fig. 6. Flowchart of the Kuhn-Munkres parallel GA framework.

Algorithm 1 Preprocessing of choosing a proper grid width

1: Procedure WIDTHCHOOSE  $\{d_1, \ldots, d_{\psi}\}$ 2:  $d \leftarrow d_1;$ 3: Compute  $\hat{U}_1$ ; 4: for  $i = \Psi \rightarrow 2$  do 5:  $d \leftarrow d_i$ ; Compute  $\hat{U}_i$ 6: if  $\hat{U}_i = \hat{U}_1$  then 7:  $d \leftarrow d_i$ ; 8: Q٠ break :

).	01
10:	end if

11:	end for
12:	return d ;

6:

4:

5:

6:

7:

8:

9:

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processes and, afterwards, all the processes of KMSPGA terminate. Pseudocode of this divide-and-conquer strategy is given in **Algorithm 2**.

Divide() contains two steps. Firstly, the target area  $\Gamma$  is uniformly partitioned into  $L \times W$  subareas  $A = (A_1, A_2, ..., A_{L \times W})$ . Fig. 7 gives four examples where  $\Gamma$  is divided into 2, 8, 32, and  $L \times W$  subareas in Fig. 7 (a), (b), (c), and (d), respectively. Then, the centers of all sensors are traversed to obtain a classification  $SE = \{SE_1, SE_2, ..., SE_{L \times W}\}$ , where  $SE_i$  is the collection of sensors falling in subarea  $A_i$ . Every sensor within  $SE_i$  satisfies that its central coordinate  $(s_x, s_y) \in A_i$  ( $k = \{1, 2, ..., N_i\}$ ). Additionally, if the center of a sensor falls on the boundaries of two or more subareas, it will be randomly scheduled into any one of the subareas. In order to keep the concision of KMSPGA, we adopt a uniform partition here instead of other ways such as clustering techniques. Besides, the uniform partition is convenient for the following combination operation.

In the parallel processing module, each process evolves a sub-population to obtain a feasible solution. The collection of sub-populations is formulized as  $SP = (SP_1, SP_2, ..., SP_{L\times W})$ . Each sub-population size is  $N_p$ . They are evolved independently through a selection, crossover, mutation, and RTSS operation in *processing()*. We estimate the state of each sub-population through periodically sampling the information of the best individual at a sampling frequency  $f_s$ . A state factor  $\zeta$  is computed as follows:

$$\zeta = \frac{1}{\hat{U}} \sum_{i=1}^{L \times W} U_i \tag{11}$$

where  $U_i$  is the number of disjoint complete cover sets obtained by the best individual of  $SP_i$ . Since  $U_i \leq \hat{U}$  (i = 0, 1, ...,  $L \times W$ ), we have  $\zeta \leq L \times W$ . Therefore, the upper limit of the threshold value,  $\zeta_c$ , is set to  $L \times W$ -1. The value of  $\zeta$ determines whether the KM combination operation will be applied. The independent evolution process will be terminated until  $\zeta$  reaches  $\zeta_c$ . Therefore, instead of performing the KM operation every generation, the execution timing of KM is adaptively adjusted based on the state factor. This way, the effectiveness of the operation is improved, and hence the computational cost is substantially reduced. The threshold  $\zeta_c$ determines the frequency of performing the KM operation (merging the local feasible solutions) and then checking for the termination condition. This procedure however does not affect the main evolution process of solutions to much extent. Thus, different settings of  $\zeta_c$  will not change the output solution quality of the proposed algorithm, but only influence the execution time. The standard uniform crossover and uniform mutation are adopted in the evolutionary process. RTSS is conducted right after mutation, and before fitness evaluation, which is introduced in detail in Subsection D. The tournament selection is adopted because of its efficiency, with a tournament size  $T_s$ .

*Conquer*() is a combination operation in order to merge the feasible local solutions  $C = (C_1, C_2, ..., C_{L\times W})^T$ ,  $C_i = (C_{i,1}, C_{i,2}, ..., C_{i,\hat{U}})$ .  $C_i$  is the best solution of the small-scale *Set K*-cover in subarea  $A_i$ . Fig. 8 shows an example of this merging process, where  $\Gamma$  is divided into 2×4 subareas. We need three steps to merge  $(C_1, C_2, ..., C_8)$  into  $C_{1-8}$ . Considering a couple of neighboring sub-populations *p* and *q*, *Conquer*() is expected

#### Algorithm 2 Process of divide-and-conquer.

```
1: Procedure DIVIDE (\Gamma)
```

2: Divide  $\Gamma$  into  $L \times W$  subareas  $(A_1, A_2, \dots, A_{L \times W})$ 

3: **for**  $i = 1 \rightarrow N$  **do** 

4: **for**  $j = 1 \rightarrow L \times W$ 

- 5: **if**  $S_i$  falling in  $A_j$  **then** 
  - Add  $S_i$  into  $SE_i$ ;

7: end if

8: end for

9: end for

10: end procedure

1: **Procedure** PARALLELPROCESSING (SP)

2: DIVIDE  $(\Gamma)$ ;

3: Repeat

for  $process = 1 \rightarrow L \times W$  parallel do

- processing ( $P_{process}$ ); //  $g_c$  is the current generation
- if  $g_c \% f_s = 0$  then // "%" is a modulus operator
- compute the state factor  $\zeta$ .
- if  $\zeta > \zeta_c$  then

CONQUER ( $C_1, C_2, ..., C_{L \times W}$ );

- 10: **end if**
- 11 end if
- 12: end for

13: Until termination conditions are satisfied.

## 14: end procedure

1: **Procedure** CONQUER ( $C_1, C_2, ..., C_{L \times W}$ ).

```
2: Repeat
```

- 3: Compute the weight matrix  $w_E$ .
- 4: Combine neighbor cover sets  $(C_p, C_q)$ ;
- 5: Until all local cover sets are merged into a global one.

## 6: end procedure

	A7	$A_8$	]	A29	A30	$A_{31}$	A32	
A2	A7	A18		A25	A 26	A27	$A_{28}$	
A2	As	$A_6$		$A_{21}$	A22	$A_{23}$	$A_{24}$	
	A	A6		A17	$A_{18}$	A 19	A20	
	A3	$A_4$		A13	A14	A15	$A_{16}$	
A	213	214		$A_9$	A10	$A_{11}$	A12	
211	A	$A_2$		As	$A_6$	$A_7$	$A_8$	
	211	712		$A_1$	$A_2$	$A_3$	$A_4$	
(a) 1×2	(	(b) 2×4			(c) -	4×8		(d) $L \times W$
E-7 Distator	£ 41		1			£ 41	1.	The terrest

Fig. 7. Division of the target area and encoding of the subareas. The target area is divided into 2, 8, 32, and  $L \times W$  subareas in figure (a), (b), (c), and (d).

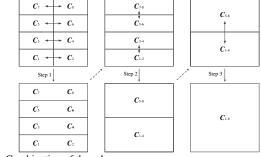


Fig. 8. Combination of the subareas.

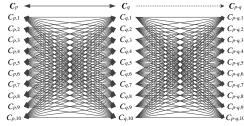


Fig. 9. Bipartite combinations of solutions  $C_p$  and  $C_q$  obtained by neighbor sub-populations, such that the Kuhn-Munkres algorithm can be applied.

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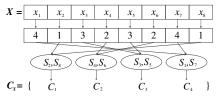


Fig. 10. Representation of chromosome. *C* is an equivalent way of representing a chromosome, where genes of the same value form a cover set. to find a best combination  $C_{p-q}$  for each dimension of  $C_p$  and  $C_q$  ensuring that the coverage ratio summation of  $C_{p-q,k}$  is maximal. The total number of matching combinations is  $\hat{U}$ !. As shown in Fig. 9, given  $\hat{U} = 10$ , we have 10! kinds of matching ways of  $C_p$  and  $C_q$ . This combination problem can be modeled as a MWPM problem in graph theory, which can now be solved using the KM algorithm.

In the literature, KM algorithm has been successfully applied to a number of fields, such as allocation of vehicle-to-infrastructure and vehicle-to- vehicle links [48], group role assignment [49], and user grouping for grouped OFDM-IDMA [50]. Given a bipartite graph G = (V, E) and weight function w(e), MWPM aims at finding a perfect matching of maximum weight. The weight of the matching M is formulized as (12). Cover sets  $(C_{i,1}, C_{i,2}, ..., C_{i,\hat{U}})$  are considered as the vertices of  $G (i \in \{p, q\})$ . Weight  $w_{i,j}$  of edge e between  $C_{p,i}$  and  $C_{q,j}$  is computed as (13), where  $|N_g(C_{p,i} \cup C_{q,j})|$  represents the number of grids covered by the sensors within  $C_{p,i}$  and  $C_{q,j}$ , and  $|N_g(A_p \cup A_q)|$  is the number of grids covered within  $A_p$  and  $A_q$ . Then, the weight matrix  $W_E$  is represented as (14). KM combination is constantly conducted until all local solutions are totally combined into a global solution.

$$w(M) = \sum_{e_k \in M} w(e_k) \tag{12}$$

$$w_{i,j} = \frac{|N_g(C_{p,i} \cup C_{q,j})|}{|N_g(A_p \cup A_q)|}$$
(13)

$$\boldsymbol{W}_{\boldsymbol{E}} = \begin{pmatrix} w_{1,1} & \dots & w_{1,\hat{U}} \\ \vdots & \ddots & \vdots \\ w_{\hat{U},1} & \dots & w_{\hat{U},\hat{U}} \end{pmatrix}$$
(14)

#### B. Chromosome Representation

In this subsection, we describe the chromosome representation in KMSPGA. We compute the value of  $\hat{U}$  in the whole target area. It is worth mentioning that even if we divided the target area into subareas, we still have to ensure that the whole target area can be covered by  $\hat{U}$  complete cover sets. Therefore,  $\hat{U}$  is also the upper limit of the number of complete cover sets for each subarea, then, sub populations  $SP_1, SP_2, \ldots, SP_{L\times W}$  share the same  $\hat{U}$ . Taking Population  $SP_k$  as an example, each chromosome is encoded as  $X = (x_1, x_2, \ldots, x_n)$ , where  $x_i$  represents the batch number of sensor i, and n is the number of sensors falling in  $A_k$  ( $n = N_k$ ). Since there is at most  $\hat{U}$  batches, we have  $x_i \in \{1, 2, ..., \hat{U}\}$ .

For chromosome X, sensors with same batch number are chosen to a same cover set. Therefore, X is transformed into  $C_X$ =  $(C_1, C_2, ..., C_0)$ , which is a candidate solution of *Set K-cover* problem. Fig. 10 gives an example of chromosome representation and shows the relationship between X and  $C_X$ , Algorithm 3 Fitness evaluation  $SF_k$ : a flag variable in case of repeat count. <u> $RF_i$ : a flag variable representing whether  $S_i$  is redundant.</u> 1: **Procedure** CHROMOSOMEEVALUATION  $(x_1, ..., x_N)$ 2: for  $i = 1 \rightarrow N$  do

2:	for $i = 1 \rightarrow N$ do
3:	$SF_i \leftarrow 0;$
4:	$RF_i \leftarrow \text{true}; // S_i \text{ is redundant if } RF_i \text{ is true}$
5:	end for
6:	for $i = 1 \rightarrow T$ do
7:	for $j = 1 \rightarrow N$ do
8:	$k \leftarrow x_j;$
9:	if $g_i \subset N_g(S_j)$ and $SF_k = 0$ then
10:	$CN_k \leftarrow CN_k + 1$ ;
11:	$SF_k \leftarrow 1;$
12:	$RF_i \leftarrow false;$
13:	end if ;
14:	end for
15:	end for
16:	$f \leftarrow 0;$
17:	for $i = 1 \rightarrow \hat{U}$ do
18:	$\delta_i \leftarrow CN_i T^1;$
19:	$f \leftarrow f + \delta_i \cdot P(\delta_i);$
20:	end for
21:	end procedure

where *n* is eight and  $\hat{U}$  is four. On the contrary, *X* can be easily transformed from  $C_X$ . Therefore, *X* and  $C_X$  are equivalent on representing an individual. In the remainder of this section, we adopt the  $C_X$  structure in representing a chromosome because this form is more convenient for introducing and descripting the operations of KMSPGA while *X* is adopted in the practical implementation of KMSPGA.

#### C. Improved Fitness Index

In STHGA [39], the evaluation function is defined as (15), where  $\delta_i$  represents the coverage ratio of cover set  $C_i$ . The computation of  $\delta_i$  is shown in (16), where the value of  $\delta_{i,k}$  is 1 if grid *k* is covered by  $C_i$ . The value of  $\delta_i$  ( $i \in \{1, 2, ..., \hat{U}\text{-}1\}$ ) is 1 in STHGA because of the forward encoding scheme. The evaluation function of GAMDSC [38] is shown in (18), where  $f_B$  represents the number of disjoint complete cover sets and [x]denotes the floor of *x*.

$$f_A = \frac{1}{T} \sum_{i=1}^{U} \delta_i \tag{15}$$

$$\delta_i = \frac{L \cdot W}{T} \sum_{k=1}^{T \cdot (L \cdot W)^{-1}} \delta_{i,k}$$
(16)

$$\delta_{i,j} = \begin{cases} 1, & \text{if grid } j \text{ is covered by } C_i \\ 0, & \text{otherwise} \end{cases}$$
(17)

$$f_B = \sum_{i=1}^{\hat{U}} \lfloor \delta_i \rfloor \tag{18}$$

In order to improve the convergence rate, we adopt a penalty function  $P(\delta_i)$  of (19), where  $\lambda$  is the penalty coefficient. The fitness evaluation function of KMSPGA is given in (20). Hence, the contribution of an incomplete cover set is lower than a complete one because of this penalty. Therefore, individuals with more incomplete cover sets are

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eliminated more easily than those with more complete cover sets. Algorithm 3 shows the pseudocode of the fitness evaluation in KMSPGA.

$$P(\delta_i) = \begin{cases} 1, & \text{if } d_i = 1\\ \lambda, & \text{otherwise} \end{cases}$$
(19)

$$f = \sum_{i=1}^{\hat{U}} \delta_i \cdot P(\delta_i)$$
(20)

#### D. Redundant-Trend Sensors Schedule Operation

In RTSS, the redundant information is indirectly utilized in order to improve search efficiency. The redundant information is collected in the fitness evaluation process. In **Algorithm 3**, steps 2 to 15 give this collection process. As can be noted from the pseudocode, the collection process is embedded in the fitness evaluation in case of increasing computational complexity. Note that after collecting the redundant information, RTSS is not applied directly after the fitness evaluation, but, as introduced in subsection A, it is performed after crossover and mutation operations, the landscape of the chromosome may change. Thus, the redundant information utilized in RTSS is hysteretic.

Considering that  $S_k$  is a member of  $C_j$ , whether  $S_k$  is redundant for  $C_j$  depending on its contributions to  $C_j$ .  $S_k$  is considered to be redundant only if it has no contributions to  $C_j$ , which is judged in the fitness evaluation. However,  $S_k$  may not still be redundant, because crossover and mutation operations may change the members of  $C_j$ . Therefore,  $S_k$  is called redundant-trend sensors in RTSS due to this uncertainty of the redundant state. Fig. 11 shows an example of this uncertainty. The grey sensor is considered to be redundant after the fitness evaluation operation. However, it is uncertain whether it is still redundant after crossover or mutation.

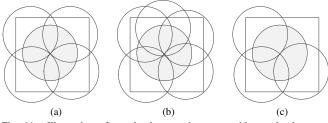


Fig. 11. Illustration of a redundant-trend sensor, with a redundant state uncertain after crossover and mutation operations. (a) Since the area covered by this sensor has already been covered by other sensors in the same cover set, the grey sensor is considered to be a redundant one. Figure (b) and (c) show two situations of redundant state of the grey sensor after the crossover and mutation operations. The grey sensor is redundant in (b) while it is no longer redundant in (c).

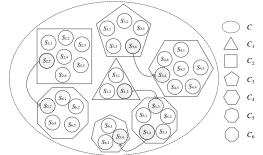


Fig. 12. Redundant-trend sensors transition between disjoint cover sets, where grey sensors are redundant-trend sensors. Different cover sets are represented

The process of RTSS is described as follows. Suppose that the cover set is  $C = \{C_1, C_2, ..., C_{\hat{U}}\}$ . Firstly, we traverse the redundant information of the sensors in  $C_i$ . Assuming  $S_{i,k}$  is the kth member of  $C_i$ , if  $S_{i,k}$  is judged to be redundant for  $C_i$  in Algorithm 3, we then consider it a redundant-trend sensor in RTSS. A cover set  $C_m$  ( $m \in \{1, 2, ..., \hat{U}\}$ ) will receive  $S_{i,k}$ through a tournament selection, where  $N_c$  candidates are randomly selected and the one with the lowest coverage ratio is chosen to receive  $S_{i,k}$  as one of its members. Fig. 12 illustrates this schedule strategy between disjoint cover sets. In Fig. 12, different polygons represent different cover sets, the grey sensors are redundant-trend sensors, and the direction of arrow represents the schedule direction. RTSS has twofold functions. It helps enhance the coverage ratio through the schedule strategy if the redundant-trend sensor is actually a redundant one. However, if the redundant sensor is no longer redundant for the current cover set, the scheduling operation becomes a disturbance for the population. This kind of stochastic disturbance enriches the diversity of the population.

#### IV. EXPERIMENTAL RESULTS

#### A. Experimental Setup

In this section, experiments are conducted to ascertain the performance of KMSPGA. In Subsection B, we compare KMSPGA with the state-of-the-art algorithms. MCMCC, GAMDSC, and STHGA are serial algorithms which perform well in solving the Set K-cover problem. The experiments and comparisons are used to verify the effectiveness of our proposed KMSPGA algorithm for lifetime maximization of large-scale WSNs. In Subsection C, we compare KMSPGA with a traditional pure parallel genetic algorithm (PGA). Further, the performance of the PGA embedding only RTSS (SPGA) or KM combination (KMPGA) are also tested in order to study the effectiveness of the two operations. Experiments in Subsection D and E are designed to evaluate the robustness of KMSPGA with different partitions and the redundant rates. In Subsection F, we conduct parameter investigation and give their suggested values. Finally, experiments in Subsections G-I are conducted with new and different testing scenarios to further verify the performance of KMSPGA.

KMSPGA and other algorithms are tested on a computer cluster of 25 nodes (with a total of 100 processing cores), which is homogenous with the same Intel core i3-3240 CPU running at 3.40 GHz, 4GB memory and Ubuntu 12.04 LTS

	TABLE I	
	THE PARAMETER SETTING OF KMSPGA	
Parameter	Description	Value
Ψ	Number of grid width classification.	5
λ	Penalty coefficient in cost evaluation.	0.2
L	Number of partitions along horizontal direction.	{2,4}
W	Number of partitions along vertical direction.	{2,4,8}
$f_s$	Sampling frequency.	3000
$P_{c}$	Crossover probability.	0.6
$P_m$	Mutation probability.	0.001
$N_c$	Number of candidates in RTSS.	[3 <sup>-1</sup> Û]
$N_p$	Sub-population size.	15
$T_s$	Tournament size	5
	Threshold of the state factor with $2 \times 2$ partitions.	3.0
~	Threshold of the state factor with 2×4 partitions.	7.0
$\zeta_c$	Threshold of the state factor with 4×4 partitions.	15.0
	Threshold of the state factor with $4 \times 8$ partitions.	30.0

	EXPERIMENT ON 12 INSTANCES IN CO												OOKITIW3						
_				~			GAMDS	SC				STHG	A		KMSPGA				
INSTANCE	Ν	$R_s$	η	U	FEs	l	J	t-test <sup>1</sup>	ç	FEs	U		t-test <sup>2</sup>	ç	FEs	U		ç	
					Mean	Mean	Std	p-value	$\mathfrak{Z}_r$	Mean	Mean	Std	p-value	$\mathfrak{I}_r$	Mean	Mean	std	$S_r$	
I1-1	5000	5	4.760	33	15000	2.0	0	0	0	10248	33	0	N/A	1.00	2688	33	0	1.00	
I1-2	5000	8	4.845	83	15000	7.4	0.67	3.31e-61	0	11507	82.9	0.25	1.6e-01	0.93	2922	83	0	1.00	
I2-1	10000	5	4.488	70	30000	1.8	0.50	4.10e-65	0	25466	69.9	0.25	1.6e-01	0.93	8418	70	0	1.00	
I2-2	10000	8	4.493	179	30000	10.2	0.57	1.76e-73	0	27077	178.8	0.38	2e-02	0.83	9324	179	0	1.00	
I3-1	15000	5	4.446	106	45000	1.3	0.50	1.15e-61	0	44887	103.9	1.13	8.90e-09	0.03	28350	105.8	0.55	0.87	
I3-2	13000	8	4.435	272	45000	12.5	0.57	6.76e-79	0	40946	271.8	0.50	2e-02	0.8	15488	272	0	1.00	
I4-1	20000	5	4.303	146	60000	1.1	0.35	9.21e-76	0	59956	143.7	1.33	1.75e-09	0.06	33086	145.9	0.25	0.93	
I4-2	20000	8	4.347	370	60000	13.2	0.76	3.12e-79	0	59802	367.9	1.36	3.7e-00	0.1	35838	370	0	1.00	
I5-1	25000	5	4.462	176	75000	1.6	0.51	9.25e-62	0	72679	175.4	0.63	3.8e-01	0.5	26761	175.7	1.27	0.93	
15-2	23000	8	4.458	451	75000	16.1	0.57	2.46e-85	0	74496	449.7	0.87	8.22e-09	0.2	36955	451	0	1.00	
I6-1	30000	5	4.446	212	90000	2.5	0.50	1.032e-77	0	87592	211.1	0.74	1.35e-07	0.3	22845	212	0	1.00	
I6-2	50000	8	4.570	528	90000	20.8	0.92	3.30e-81	0	86594	527.6	0.62	1e-02	0.67	34785	528	0	1.00	
10-2		0	÷.570	520	70000	20.0		5.500-81	5	00374	521.0	0.02	10-02		34/03		0	1.0	

TABLE II EXPERIMENT ON 12 INSTANCES IN COMPARISON WITH SERIAL ALGORITHMS

1 Two tailed t-test of the null hypothesis that GAMDSC is equal to KMSPGA. 2 Two tailed t-test of the null hypothesis that STHGA is equal to KMSPGA.

64-bit operating system. The parallel programming practice uses the Message Passing Interface (MPI). Table I shows the parameter settings.  $P_c$  and  $P_m$ , the constant crossover and mutation rates, are set as  $P_c = 0.6$  and  $P_m = 0.001$ , respectively. The value of  $\lambda$  determines the degree of punishment for the individuals with incomplete cover sets. In all of the test instances, we adopt  $\lambda = 0.2$  as default value.  $N_c$ , the number of candidates in RTSS, is empirically setting to 3<sup>-1</sup>Û. Sample frequency  $f_s$  is set to 3000 function evaluations (FEs). In Subsections *B*-*H*,  $\Gamma$  is a 50×50 square area, and in Subsection *I*,  $\Gamma$  is a 3-D surface. P<sub>L×W</sub> denotes the partition way of  $\Gamma$ .  $\zeta_c$  is set to 3.0, 7.0, 15.0, and 30.0 for  $P_{2x2}$ ,  $P_{2x4}$ ,  $P_{4x4}$ , and  $P_{4x8}$ , respectively. Thirty trials are performed for each instance and the results are averaged over the trials. A two-tailed t-test of the null hypothesis is conducted in Subsections B, G, H, and I. The null hypothesis will be rejected if *p*-value is smaller than the significance level  $\alpha = 0.05$ .

#### B. Comparison with State-of-the-Art Serial Algorithms

We conduct experiments on 12 instances to verify the performance of KMSPGA in comparison with the serial algorithms: MCMCC, GAMDSC, and STHGA.  $P_{2x4}$  is adopted in this subsection.

The experimental results are listed in Table II. Mean and Std are the mean quality and standard deviation.  $S_r$  is the success rate. The best solutions are marked in bold. KMSPGA outperforms the other algorithms in terms of the convergence rate and solution quality. Furthermore, KMSPGA produces significant increases both in the convergence rate and in the success rate in the higher dimensional space. The success rate of instances I1-1, I1-2, I2-1, I2-2, I3-2, I4-2, I5-2, I6-1, and I6-2 reach 1.00. The mean FEs is far less than the serial algorithms. MCMCC is not available because it took an unacceptable time before termination. The worst runtime of MCMCC is  $O(N^2)$  [39], where N is the number of sensors. This heuristic method performs efficiently when the number of sensors is small or medium. However, the computational complexity of MCMCC becomes so high in terms of large-scale WSNs that it fails to work out a feasible solution in an acceptable time, i.e., I2-1, a single run of MCMCC exceeds eight hours. GAMDSC is another genetic algorithm used in this experiment. Since GAMDSC lacks an efficient search strategy to handle such a large number of sensors, solutions obtained by GAMDSC are fewer than U. STHGA possesses high quality of solutions in the low-dimensional spaces when

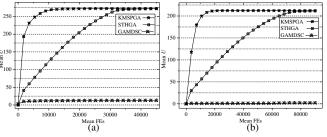


Fig. 13. Convergence curves of the compared algorithms. (a) I3-2. (b) I6-1. the number of sensors is less than 5000. The complicated local search operations help STHGA search the problem space efficiently. However, the performance of this serial GA is badly influenced by curse of dimensionality. The success rate of instance I3-1 obtained by STHGA is 0.03. At the same time, the large number of function evaluations indicates that the searching efficiency reduces due to the curse of dimensionality.

Fig. 13 shows the convergence curves of the compared algorithms on instances I3-2 and I6-1, where the *x*-axis represents the mean FEs and the *y*-axis the mean *U* over 30 trials. The convergence rate of KMSPGA is higher than STHGA and GAMDSC. Furthermore, KMSPGA generates higher-quality solutions. In Fig. 13 (a), KMSPGA obtains the optimal value at about 15,000 FEs, STHGA reaches the near optimal value at 35,000 FEs, and GAMDSC evolves very slowly with a low-quality solution. Similarly, in Fig. 13 (b), KMSPGA converges to the optimal value at about 20,000 FEs, STHGA obtains the near optimal value at about 20,000 FEs, and GAMDSC still evolves very slowly with a low-quality solution.

In this subsection, we investigate the effectiveness and reliability of KMSPGA and the other compared algorithms with the same maximum number of function evaluations so that the comparison is fair. Some of the compared algorithms, e.g., STHGA, are not suitable for parallelism. The reasons are as follow. As can be noted from our description of the proposed parallel genetic framework, the disjoint cover sets within the same chromosome are supposed to be peer to each other. In STHGA,  $C_i$  is a complete cover set while  $C_{U+1}$  is incomplete due to the forward encoding scheme, which makes  $C_i$  ( $i \in \{1, 2, ..., U\}$ ) is not peer to cover set  $C_{U+1}$ . Combination between any  $C_{p,i}$  and  $C_{q,j}$  (i, j < U) makes no sense to these already complete cover sets. The complicated auxiliary search

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EXPERIMENTAL RESULTS ON 12 INSTANCES IN COMPARISON WITH PARALLEL ALGORITHMS OF A 2×4 PARTITIC											TION.						
		PG.	A			KMP	PGA			SPO	GA		KMSPGA				
INSTANCE	FEs	l	IJ	Sr	FEs	ι	IJ	c	FEs	i	IJ	c	FEs	i	IJ	S <sub>r</sub>	
	Mean	Mean	Std	$\mathfrak{I}_r$	Mean	Mean	Std	$S_r$	Mean	Mean	Std	$S_r$	Mean	Mean	Std	$\mathfrak{I}_r$	
I1-1	15000	28.7	1.34	0.00	14719	31.37	1.24	0.17	15000	28.9	1.73	0.00	2688	33	0	1.00	
I1-2	15000	71.0	2.06	0.00	1500	78	1.68	0.00	4114	83	0	1.00	2922	83	0	1.00	
I2-1	30000	63.5	2.19	0.00	29998	68.4	1.04	0.13	14688	69.8	0.50	0.80	8418	70	0	1.00	
I2-2	30000	119.5	5.36	0.00	30000	159.8	3.77	0.00	29640	176.7	1.18	0.03	9324	179	0	1.00	
I3-1	45000	83.2	3.85	0.00	45000	96.7	2.03	0.00	45000	100.3	2.12	0.00	28350	105.8	0.55	0.87	
I3-2	45000	160.7	6.30	0.00	45000	231.3	6.01	0.00	43619	270.4	1.30	0.23	15488	272	0	1.00	
I4-1	60000	106.1	4.90	0.00	60000	122.4	3.01	0.00	56098	144.2	1.19	0.17	33086	145.9	0.25	0.93	
I4-2	60000	156.0	6.24	0.000	60000	290	3.06	0.00	60000	358.2	4.03	0.00	35838	370	0	1.00	
I5-1	75000	137.8	3.96	0.00	75000	156.7	3.06	0.00	40688	175.8	0.41	0.80	26761	175.7	1.27	0.93	
I5-2	75000	198.7	9.10	0.00	75000	370.1	6.10	0.00	74700	447.6	1.99	0.03	36955	451	0	1.00	
I6-1	90000	138.8	6.18	0.00	90000	170.2	5.30	0.00	31485	211.9	0.18	0.97	22845	212	0	1.00	
I6-2	90000	277.4	8.40	0.00	90000	428.3	7.66	0.00	39085	528	0	1.00	34785	528	0	1.00	

TABLE IV

EXPERIMENTAL RESULTS ON 12 INSTANCES IN COMPARISON WITH PARALLEL ALGORITHMS OF A 4×4 PARTITION.

		PGA	ł			KMP	GA			SPC	iΑ		KMSPGA					
INSTANCE	FEs		U c		FEs	U	J	c	FEs	L	7	c	FEs	U	7	c		
	Mean	Mean	Std	$S_r$	Mean	Mean	Std	$S_r$	Mean	Mean	Std	$S_r$	Mean	Mean	Std	$S_r$		
I1-1	14900	28.9	1.76	0.03	14597	31.8	0.86	0.20	14920	29.4	1.57	0.03	3660	33	0	1.00		
I1-2	15000	70.8	2.59	0.00	15000	78.7	1.57	0.00	6414	82.9	0.25	0.93	4314	83	0	1.00		
I2-1	30000	65.3	1.54	0.00	29599	65.8	3.93	0.07	25997	68.5	1.07	0.20	3960	70	0	1.00		
I2-2	30000	135.7	5.02	0.00	30000	170.0	2.03	0.00	30000	175.1	1.89	0.00	11748	179	0	1.00		
I3-1	45000	90.4	2.33	0.00	45000	101.1	1.46	0.00	45000	99.6	1.85	0.00	13230	106	0	1.00		
I3-2	45000	198.1	6.48	0.00	45000	258.7	3.42	0.00	44580	268.7	1.91	0.03	23876	271.9	0.18	0.97		
I4-1	60000	129.2	2.85	0.00	60000	132.6	9.10	0.00	60000	141.6	2.04	0.00	22782	145.9	0.25	0.93		
I4-2	60000	219.4	9.36	0.000	60000	340.0	4.69	0.00	60000	361	2.36	0.00	29412	370	0	1.00		
I5-1	75000	169.2	2.36	0.00	72395	174.8	1.28	0.30	31588	175.8	0.41	0.80	8625	176	0	1.00		
I5-2	75000	305.1	8.55	0.00	75000	410.3	5.78	0.00	71599	449.1	1.53	0.20	29802	451	0	1.00		
I6-1	90000	203.2	2.57	0.00	89399	209.7	1.70	0.07	25586	211.9	0.18	0.97	9975	212	0	1.00		
I6-2	90000	453.3	5.29	0.00	90000	504.5	4.73	0.00	26985	528	0	1.00	24285	528	0	1.00		

operations adopted by STHGA also increases the difficulty of parallelizing the algorithm.

# C. Comparison with Parallel Algorithms

In this subsection, we compare KMSPGA with the pure PGA, KMPGA, and SPGA. The PGA combined only with the KM combination or RTSS form KMPGA or SPGA. The experimental instances in Subsection *B* are adopted here. We conduct the experiments under two different partitions:  $P_{2x4}$  and  $P_{4x4}$ .

The experimental results are given in Table III and Table IV. KMSPGA outperforms KMPGA and SPGA in all of the instances. Although  $S_r$  of KMPGA and SPGA is 0 in the majority of the instances, the mean U obtained by KMPGA and SPGA is much larger than PGA, which reveals that the KM combination and RTSS contribute to the enhancement of the solution quality, which is made available by KMSPGA. Take instance I4-2 of partition 2×4 as an example, where  $\hat{U}$  is 370. The mean U obtained by PGA is 156.0, accounting for only 42.16%. As for KMPGA and SPGA, the mean U obtained are 290 and 358.2, accounting for 78.38% and 96.81%. SPGA obtains larger mean U than KMPGA in the majority of the instances, and therefore contribution of RTSS is larger than that of KM combination when it comes to the degree of the improvement of solution quality.

It is worth mentioning that parallel evolutionary algorithms are suitable for the problems of a high dimensionality or of complex and time-consuming computation features [51], such as large-scale air traffic flow optimization [52], discrete resource allocation in classic economic field [53], and large-scale function optimization [54][55]. They are adopted to either speed up the optimization or enhance the solution

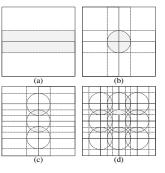


Fig. 14. Visual illustration of correlative areas under representative partitions. Grey areas represent the special correlative areas where sensors are relevant with the largest number of subareas. Sensors falling in grey areas are correlation with 2, 4, 6 and 6 subareas in (a), (b), (c), and (d) separately.

quality through a dimensionality reduction strategy. In this paper, the dimensionality and computational complexity of *Set K*-*cover* problem become so high in large-scale WSN that we adopt a parallel evolutionary algorithm for performance enhancement.

#### D. Discussion on Correlation between Sensors and Subareas

We investigate how the number of partitions affects the solution quality in this subsection. It can be intuitively noted that the number of subareas covered by one sensor increases as the target area is divided more finely. A larger number of subareas covered by one sensor means a higher correlation between the sensors and subareas. The high correlation influence the performance of parallel processing module and the KM combination. Fig. 14 gives a visual illustration of a special kind of correlative area (marked as grey), and sensors

	EXPERIMENTAL RESULTS UNDER DIFFER									ERENT PARTITIONS ON 12 INSTANCES							
	$P_{2\times 2}$	(ω <sub>-5</sub> = 1.3	8, ω <sub>-8</sub> =1.	67)	P <sub>2×4</sub> (	ω <sub>-5</sub> = 1.79,	ω <sub>-8</sub> =2.4	1)	P <sub>4×4</sub> (	ω <sub>-5</sub> =2.30,	, ω <sub>-8</sub> =3.4	-3)	$P_{4\times 8} (\omega_{-5} = 3.41, \omega_{-8} = 5.55)$				
INSTANCE	FEs	U	J	ç	FEs	L	J	ç	FEs	U	J	ç	FEs	U	ſ	ç	
	Mean	Mean	Std	$S_r$	Mean	Mean	Std	$\mathfrak{Z}_r$	Mean	Mean	Std	$S_r$	Mean	Mean	Std	$S_r$	
I1-1	4965	33	0	1.00	2688	33	0	1.00	3660	33	0	1.00	4935	33	0	1.00	
I1-2	2730	83	0	1.00	2922	83	0	1.00	4314	83	0	1.00	2016	83	0	1.00	
I2-1	14604	69.9	0.57	0.83	8418	70	0	1.00	3960	70	0	1.00	4650	70	0	1.00	
I2-2	9010	179	0	1.00	9324	179	0	1.00	11748	179	0	1.00	3744	179	0	1.00	
I3-1	41490	104	1.24	0.37	28350	105.8	0.55	0.87	13230	106	0	1.00	15159	106	0	1.00	
I3-2	3718	272	0	1.00	15488	272	0	1.00	23876	271.9	0.18	0.97	37464	265.1	4.55	0.23	
I4-1	59445	143.8	1.44	0.1	33086	145.9	0.25	0.93	22782	145.9	0.25	0.93	45330	145.7	0.65	0.8	
I4-2	52500	369.7	0.71	0.76	35838	370	0	1.00	29412	370	0	1.00	57060	346.6	6.89	0.07	
I5-1	67695	174.5	1.98	0.43	26761	175.7	1.27	0.93	8625	176	0	1.00	10230	176	0	1.00	
I5-2	66675	449.3	1.82	0.37	36955	451	0	1.00	29802	451	0	1.00	71010	433.2	5.71	0.07	
I6-1	87660	209.6	2.22	0.2	22845	212	0	1.00	9975	212	0	1.00	12084	212	0	1.00	
I6-2	78615	526.1	2.83	0.43	34785	528	0	1.00	24285	528	0	1.00	68115	521.4	5.06	0.30	

TABLE V EXPERIMENTAL RESULTS UNDER DIFFERENT PARTITIONS ON 12 INSTANC

falling in the grey area will be correlate with 2, 4, 6 and 6 subareas in (a), (b), (c), and (d), respectively. A sensor covers at least two subareas when it falls into the correlative areas. The correlative areas influence the independence of the evolution of each sub-population. The average number of subareas covered by one sensor is adopted to reveal the degree of correlation between sensors and subareas. A Monte Carlo method is utilized to estimate this value. One hundred thousand sensors (i.e., N = 100,000) are randomly deployed into the target area, then the number of subareas covered by each sensor is computed. The average number of subareas covered by one sensor is computed as  $\omega$ :

$$\omega = \frac{1}{N} \sum_{i=1}^{N} N_{sa}(S_i) \tag{22}$$

where  $N_{sa}(S_i)$  is the number of subareas covered by sensor  $S_i$ . A larger value of  $\omega$  means a stronger correlation between sensors and subareas. The target area is expected to be divided into more subareas in order to achieve dimensionality reduction as the number of sensors increases. However, it is unreasonable to increase the number of subareas without constraints, because the efficiency of combination strategy decreases as the number of subareas increases.

Table V lists the results with different partitions:  $P_{2\times 2}$ ,  $P_{2\times 4}$ ,  $P_{4\times4}$ , and  $P_{4\times8}$ , where  $\omega_{-r}$  represents the value of  $\omega$  with sensing radius r.  $P_{4\times 4}$  achieves the best performance in the majority of test instances in terms of convergence rate, solution quality, and success rate. The success rate is low considering the performance of  $P_{2\times 2}$  and  $P_{4\times 8}$ . However, the reason is completely different. As for  $P_{2\times 2}$ , the partition quantity is not enough to reduce the dimensionality to an acceptable level. On the contrary, the partition quantity of  $P_{4\times8}$  is so large that the correlation between sensors and subareas becomes too high to apply the divide-and-conquer strategy. As can be noted in instance Ix-1 and Ix-2, the success rate of the former is clearly higher than the later because of the smaller value of  $\omega$  in Ix-1 than that in Ix-2. Consequently, the number of partitions is restricted by the value of  $\omega$ . In order to help the divide-and-conquer strategy work efficiently, the partition quantity is limited to an appropriate range.

#### E. Experiments with Different Redundant Rate

In this subsection, we study the influence of different redundant rates on the solution quality. Two groups of experimental instances are adopted in this experiment, where the sensor radius is 5. Instances prefixed by "J" represent the number of sensors is 20,000, whereas instances prefixed by "H" represent the number of sensors is 25,000. Although the number and the sensing radius of sensors are fixed, the redundant rate can be different because of the random deployment strategy. It is quite difficult to generate an instance with a specified  $\hat{U}$ . Instead, to generate this test set with different redundant rates, we create a relatively large number of candidate instances, calculate their  $\hat{U}$ s, and then select the candidate instances with appropriate  $\hat{U}$  to the test set. The redundant rate ranges from 3.997 to 5.003.

Performance of different partitions, i.e. P<sub>2x4</sub>, P<sub>4x4</sub>, and P<sub>4x8</sub>, is tested. STHGA is also adopted for comparison. Results are summarized in Table VI. KMSPGA ( $P_{2x4}$ ,  $P_{4x4}$ , and  $P_{4x8}$ ) achieves a high success rate in a large range of redundant rates, which indicates that KMSPGA offers very promising performance with robustness.  $P_{4\times4}$  achieves the fastest convergence rate, the largest mean U, and the highest  $S_r$  in the majority of the instances in Table VI. As far as STHGA is concerned, the success rate declines sharply when the redundant rate decreases. The x-axis of Fig. 15 (a) and (b) is the redundant rate, the y-axis of Fig.15 (a) represents mean  $S_r$ , and the y-axis of Fig. 15 (b) represents the ratio of mean U to  $\hat{U}$ . The best solution obtained by KMSPGA among different partitions in each instance is represented by  $P_{BEST}$ . In Fig. 15 (a), KMSPGA ( $P_{2\times4}$ ,  $P_{4\times4}$ , and  $P_{4\times8}$ ) obtains a high  $S_r$  in most of the instances while the  $S_r$  of STHGA declines sharply as  $\eta$ decreases. Fig. 15 (b) also indicates that KMSPGA possesses high solution quality within a large-scale range of  $\eta$ . In fact,  $P_{\text{BEST}}$  maintains a value of 1.00 in both Fig. 15 (a) and (b).

# F. Parameter Investigation

We investigate the penalty coefficient  $\lambda$  and the threshold value of state factor  $\zeta_c$  in this subsection in order to show how these parameters influence the performance of KMSPGA.

The penalty coefficient  $\lambda$  is adopted to improve the convergence rate. We adopt test instances I4-2 and I5-1 with partition P<sub>4×4</sub>. The value of  $\lambda$  is increased from 0 to 1 with step 0.05. Based on the experimental results, we give the relationship between  $\lambda$  and FEs in Fig. 16. As can be noted from the results, the value of FEs when  $\lambda$  is set within [0.05, 0.95], the required FEs is significantly less than that when  $\lambda$  is set to 1 (no penalty). The experimental results indicate that the penalty coefficient  $\lambda$  has an effect of improving the convergence rate of KMSPGA. Meanwhile, KMSPGA is generally insensitive to the value of  $\lambda$ , and it works identically well when  $\lambda$  is set to [0.2, 0.7].

TABLE VI EXPERIMENTAL RESULTS WITH DIFFERENT REDUNDANT RATES ON 20 INSTANCES

			EXPERIMENTAL RESULTS WITH DIFFERENT REDUN							JIDANI KATES UN 20 INSTANCES									
IN	ISTAN	CE		STH	GA			$P_{2\times 4}$				P4>	(4		P <sub>4×8</sub>				
Na.	Û		FEs	U	J	c	FEs	U	J	$S_r$	FEs	L	J	$S_r$	FEs	U	r -	c	
Ina.	U	η	Mean	Mean	Std	$S_r$	Mean	Mean	Std	$S_r$	Mean	Mean	Std	$S_r$	Mean	Mean	Std	$S_r$	
J01	129	4.871	51407	128.8	0.38	0.83	2824	129	0	1.00	2508	129	0	1.00	3394	129	0	1.00	
J02	133	4.724	52332	132.9	0.25	0.93	5802	133	0	1.00	3684	133	0	1.00	3904	133	0	1.00	
J03	135	4.654	53898	134.8	0.38	0.83	6334	135	0	1.00	4176	135	0	1.00	6222	135	0	1.00	
J04	136	4.620	56295	135.8	0.50	0.80	9714	136	0	1.00	5856	136	0	1.00	6664	136	0	1.00	
J05	139	4.520	54113	138.8	0.38	0.83	9210	139	0	1.00	7410	139	0	1.00	7485	139	0	1.00	
J06	145	4.333	59789	143.1	1.05	0.06	18705	145	0	1.00	13350	145	0	1.00	15204	145	0	1.00	
J07	149	4.217	59281	147.8	1.07	0.30	29544	149	0	1.00	16740	149	0	1.00	10512	149	0	1.00	
J08	155	4.080	60000	146.1	1.58	0.00	60000	150.5	2.13	0.00	25875	155	0	1.00	53992	154.0	1.48	0.53	
J09	157	4.002	60000	152	1.69	0.00	57819	155.5	1.57	0.37	35235	157	0	1.00	59925	151.7	2.63	0.03	
J10	158	3.977	60000	154.0	1.52	0.00	37884	158	0	1.00	33045	157.9	0.40	0.93	57420	156.2	1.21	0.17	
H01	157	5.003	56796	157	0	1.00	2674	157	0	1.00	4680	157	0	1.00	4035	157	0	1.00	
H02	159	4.940	62349	158.9	0.35	0.86	3524	159	0	1.00	4485	159	0	1.00	5250	159	0	1.00	
H03	162	4.848	63270	161.9	0.31	0.90	3954	162	0	1.00	4485	162	0	1.00	4635	162	0	1.00	
H04	164	4.789	63223	163.9	0.31	0.90	5054	164	0	1.00	5085	164	0	1.00	5805	164	0	1.00	
H05	167	4.703	67287	166.9	0.35	0.87	7305	167	0	1.00	4635	167	0	1.00	6030	167	0	1.00	
H06	173	4.540	72310	172.5	0.63	0.57	15330	173	0	1.00	7632	173	0	1.00	16059	173	0	1.00	
H07	175	4.488	73608	174.1	0.92	0.40	19980	175	0	1.00	11430	175	0	1.00	19960	175	0	1.00	
H08	178	4.412	74741	176.4	1.04	0.10	19685	178	0	1.00	10932	178	0	1.00	15810	178	0	1.00	
H09	181	4.340	74003	179.9	0.96	0.33	20730	181	0	1.00	10575	181	0	1.00	15510	181	0	1.00	
H10	185	4.245	74570	183.4	1.10	0.13	41559	185	0	1.00	18375	185	0	1.00	24435	185	0	1.00	

The threshold value of the state factor,  $\zeta_c$ , determines the frequency of executing the KM operation. We adopt test

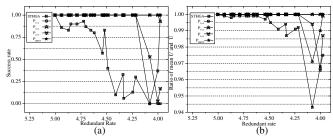


Fig. 15. Influence of redundant rate on the performance of KMSPGA. (a) Influence on the mean success rates. (b) Influence on the mean ratios.

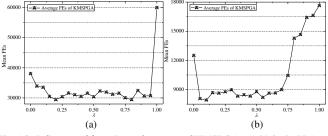
instance I2-2 to investigate the effect of this parameter. For partition  $P_{2\times4}$  and  $P_{4\times4}$ ,  $\zeta_c$  is increased from 0 to 7 with step 0.5 and from 0 to 15 with step 1, respectively. The experimental results show that, KMSPGA achieves a 100% success rate with different  $\zeta_c$ . Therefore,  $\zeta_c$  has no influence on the solution quality. Fig. 17 shows the influence of the value of  $\zeta_c$  on the FEs and a practical execution time of KMSPGA. It can be observed that  $\zeta_c$  has negligible influence on the convergence rate. Based on the experimental results, we suggest that the value of  $\zeta_c$  is set to  $|0.94 \times L \times W|$ .

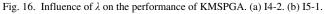
## G. Proof-of-Principle Experiments

To conduct proof-of-principle experiments, we first apply a deterministic deployment strategy to generate *K* complete cover sets. The superposition of these *K* complete cover sets results in a *Set K-cover* instance, to which an optimal solution is known. Then KMSPGA and other compared algorithms are applied to solving this instance.

Here, it is to be noted that deterministically deploying the least number of circles to cover any polygon is an NP-hard problem as discussed in [21], [56], and [57]. Nevertheless, if we relax the requirement of that the layout should be "theoretically best with the least circles", there exists an efficient deterministic node placement strategy to cope with this issue. Fig. 18 shows a "tessellation" placement strategy. First, the target area is completely tiled by a number of

compact polygons. Then, sensors are placed at the vertices of polygons. If each polygon is covered by at least one sensor, the





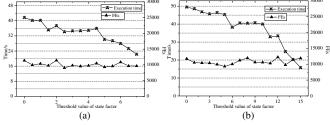


Fig. 17. Influence of threshold value  $\zeta_c$  on the performance of KMSPGA. (a)  $P_{2x4}$ . (b)  $P_{4x4}$ .

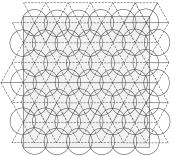


Fig. 18. Illustration of "tessellation" placement strategy, and the polygons adopted here is triangle.

target area is completely covered. Fig. 18 is adopted as 1-cover set (unit set) since it is the optimal tessellation requiring the minimum number of sensors. We then repeatedly add K unit sets into a cover set to generate a K-cover as the instance of the proof-of-principle experiment. Six new instances are

	EXPERIMENTAL RESULTS OF THE PROOF-OF-PRINCIPLE EXPERIMENT ON 6 INSTANCES																
				G	AMDSC					STHGA	1		KMSPGA				
INSTANCE	R	Û	FEs	3	U	c	t-test <sup>1</sup>	FF	Es	U	c	t-test <sup>2</sup>	F	Es	U	6	
			Mean	Std	Mean	$S_r$	p-value	Mean	Std	Mean	$S_r$	p-value	Mean	Std	Mean	$S_r$	
DD1-1	5	100	13500	0	1	0	4.14e-80	598	0.46	100	1.00	1.09e-22	463.5	25.94	100	1.00	
DD1-2	8	100	6600	0	4.1	0	3.31e-70	598	0.58	100	1.00	1.09e-30	328.5	27.39	100	1.00	
DD2-1	5	300	40500	0	1	0	1.30e-86	1798	0.72	300	1.00	3.13e-40	779.5	47.13	300	1.00	
DD2-2	8	300	19800	0	6.2	0	3.07e-78	1798	0.50	300	1.00	2.56e-44	494.0	44.56	300	1.00	
DD3-1	5	600	81000	0	1	0	5.18e-82	3598	0.52	600	1.00	1.87e-37	1285.5	136.35	600	1.00	
DD3-2	8	000	39600	0	8.0	0	8.54e-82	3598	0.46	600	1.00	3.08e-49	661.0	67.76	600	1.00	

TABLE VII EXPERIMENTAL RESULTS OF THE PROOF-OF-PRINCIPLE EXPERIMENT ON 6 INSTANCE

TABLE	VIII

			Exp	PERIMENT	TAL RESULTS	WITH SE	NSORS OF I	DIFFERENT	RADIUS	es on 5 Inst	ANCES				
	Û			C			STHGA	KMSPGA							
INSTANCE		FEs	U		t-test <sup>1</sup>	FEs	U		t-test <sup>2</sup>	c	FEs	L	J	c	
		Mean	Mean	Std	p-value	$S_r$	Mean	Mean	Std	p-value	$\mathbf{S}_{\mathrm{r}}$	Mean	Mean	Std	Sr
DR1	159	45000	5.1	0.48	1.98e-74	0	44815	157.7	0.71	3.6e-11	0.1	11714	159	0	1.00
DR2	220	45000	8.1	0.71	1.65e-73	0	44148	219.0	0.96	2.28e-06	0.33	12954	220	0	1.00
DR3	276	45000	15.8	0.92	8.4e-73	0	41145	275.6	0.61	2.8e-03	0.7	14794	276	0	1.00
DR4	163	45000	5.4	0.50	4.0e-74	0	40232	162.8	0.40	1.2e-02	0.8	6914	163	0	1.00
DR5	232	45000	7.7	0.58	4.31e-75	0	44762	230.7	0.99	8.3e-07	0.16	25714	231.9	0.25	0.93

1 Two tailed t-test of the null hypothesis that GAMDSC is equal to KMSPGA. 2 Two tailed t-test of the null hypothesis that STHGA is equal to KMSPGA.

EXPERIMENTAL

TABLE IX													
D		D	<b>D</b>										

EXPERIMENTAL RESULTS IN 3-D ENVIRONMENT ON 6 INSTANCES																	
INSTANCE	Ν	$R_s$	Û	GAMDSC							KMSPGA						
				FEs	L	J	t-test <sup>1</sup>	ç	FEs	l	U	t-test <sup>2</sup>	$\mathbf{S}_{\mathrm{r}}$	FEs	U		$S_r$
				Mean	Mean	Std	p-value	$S_r$	Mean	Mean	Std	p-value		Mean	Mean	Std	
3D1-1	5000	5	25	15000	1	0	0	0	12673	24.9	0.31	8e-02	0.9	4785	25	0	1.00
3D1-2		8	76	15000	3.5	0.51	2.84e-64	0	13449	75.9	0.25	16e-01	0.93	4618	76	0	1.00
3D2-1	10000	5	54	30000	1	0	4.91e-69	0	30000	50.5	0.97	8.3e-18	0	16309	53.9	0.25	0.93
3D2-2		8	163	30000	3.2	0.43	4.12e-67	0	30000	157.2	1.31	2.09e-18	0	20756	162.5	0.62	0.67
3D3-1	3D3-1 3D3-2 15000	5	84	45000	1	0	2.34e-63	0	45000	75.1	1.18	1.70e-25	0	41736	83.6	0.62	0.67
3D3-2		8	250	45000	3.5	0.51	1.35e-66	0	45000	234.7	2.17	5.03e-24	0	34290.5	249.1	1.35	0.63

1 Two tailed t-test of the null hypothesis that GAMDSC is equal to KMSPGA. 2 Two tailed t-test of the null hypothesis that STHGA is equal to KMSPGA.

generated in this way with K setting to 100, 300, and 600. The numbers of sensors covering the target area are 45 and 22 when  $R_s$  is 5 and 8, respectively.

The experimental results are given in Table VII. Both STHGA and KMSPGA achieve 100% success rate in the tests. However, KMSPGA achieves a higher convergence rate than STHGA. The experimental results also indicate that these ideal and regular test instances are easy to solve by the two algorithms. The reasons are presented as follow. Aiming at using the least number of circles to realize complete coverage, the deterministic deployment strategy tends to minimize the overlapping area of neighboring circles. Thus, for each sensor, the coverage ratio of each cover set will be quite different considering whether or not the sensor is assigned to the right cover set. This feature makes the fitness evolution possess good differentiation for different individuals (candidate solutions) and hence provides a promising guidance for the search. Furthermore, this feature also benefits the proposed redundancy-based schedule strategy, owing to the uniqueness of sensor within each unit set.

# H. Experiments with Sensors of Different Radiuses

Although in the above experiments, sensors of identical radius are assumed for simplicity. However, as the proposed KMSPGA algorithm does not contain any radius-related parameters or operators, it is a generic algorithm suitable for both application scenarios with homogenous or heterogeneous sensors deployed. In this subsection, we conduct experiments using sensors of different radiuses to investigate the performance of KMSPGA. Five new instances are generated and tested, in which the radiuses of sensors follow Gaussian distribution with different mean values and standard deviations. The radius of sensor *j* in instance *i* is set to  $R_i+r_j$ , where  $r_j$  is a random number following a standard normal distribution. For  $i \in \{0, 1, 2\}$ ,  $R_i$  is set to 6, 7, and 8, respectively. For i = 3 and i = 4,  $R_i$  is randomly chosen within an integer interval [5, 7] and [6, 8], respectively. The number of each instance is 15000.

From the experimental results given in Table VIII, it is still observed that KMSPGA achieves significantly higher solution quality and success rate among the compared algorithms in most of the instances. Besides, KMSPGA achieves smallest standard deviations, which indicates that KMSPGA possesses high stability. The results confirm that KMSPGA works well with sensors of different radiuses.

# I. Experiment in a 3-D Environment

In the literature of WSN lifetime maximization, sensors are assumed to be uniformly distributed on an ideal square plane. However, in practice, this is not always the case. Instead, sensors are often deployed on a 3-D surface so that the sensor distribution is no longer uniform, but it is highly dependent to the shape/landscape of the surface [58]. This uneven distribution makes it even harder to optimally schedule the sensors into the right complete cover sets. In this subsection, we implement KMSPGA to test its effectiveness and reliability on this application scene.

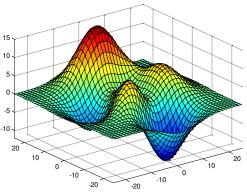


Fig. 19. Simulated free flowing contour.

The multi-peak function used to simulate the 3-D surface is given as

$$\Gamma_{3-D}(x,y) = (3 \cdot (1 - \frac{x}{10})^2 \cdot e^{-(\frac{x}{10})^2 - (\frac{y}{10} + 1)^2} - 10 \cdot (\frac{x}{50} - (\frac{x}{10})^3 - (\frac{y}{5})^5) \cdot e^{-(\frac{x}{10})^2 - (\frac{y}{10})^2} - \frac{1}{3} \cdot e^{-(\frac{x}{10} + 1)^2 - (\frac{y}{10})^2}) \cdot 2$$
(22)

which consists of three peaks and three valleys. Its 3-D view is shown in Fig. 19. The number of sensors is set to 500, 10000, and 15000. GAMDSC and STHGA are also modified to solve the *Set K-Cover* problem in this 3-D environment. Experimental results given in Table IX indicate that these instances are significantly hard to solve. The performance of all the algorithms decreases on this test set. GAMDSC and STHGA are unable to achieve an acceptable solution within the limited FEs. In comparison, the proposed KMSPGA still offers a high success rate and solution quality, which further confirms the effectiveness and reliability of KMSPGA.

#### V. CONCLUSION

Due to the curse of dimensionality, existing set cover algorithms are unable to provide satisfactory efficiency for large-scale WSNs scheduling. In this paper, we have developed a Kuhn-Munkres parallel genetic algorithm, KMSPGA, for the set cover problem and applied it to lifetime maximization of large-scale WSNs. The KMSPGA framework is based on a divide-and-conquer strategy of dimensionality reduction. Firstly, the target area is divided into several subareas, and then individuals are evolved independently in each subarea until the state factor reaches a predefined value. The polynomial Kuhn-Munkres algorithm is then utilized to splice the solutions obtained in each subarea so as to generate global optimal solution of the entire problem. KMSPGA also includes a novel schedule operation for further improvement in performance.

Eight types of experiments have been conducted to verify the design and effectiveness of KMSPGA. The experimental results indicate that KMSPGA achieves a higher convergence rate, solution quality, success rate, and scalability. Further, by investigating the influence of different partitions, redundant rate, penalty coefficient on the performance of KMSPGA, KMSPGA is also seen to offer high robustness. Finally, experimental results on new testing scenarios indicate that KMSPGA achieves wide applicability. Future work includes the development of an improved matching algorithm for the combination operation, of distributed, cloud and multi-objective versions of KMSPGA and their applications to various kinds of real-world problems.

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