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Locational Optimization based Sensor Placement for Monitoring Gaussian Processes Modeled Spatial Phenomena

Linh Van Nguyen, Sarath Kodagoda, Ravindra Ranasighe, and Gamini Dissanayake

Center for Autonomous Systems

Faculty of Engineering and IT

University of Technology, Sydney, Australia

Email: van.l.nguyen-1@student.uts.edu.au; sarath.kodagoda, ravindra.ranasighe, gamini.dissanayake@uts.edu.au

Abstract—This paper addresses the sensor placement problem associated with monitoring spatial phenomena, where mobile sensors are located on the optimal sampling paths yielding a lower prediction error. It is proposed that the spatial phenomenon to be monitored is modeled using a Gaussian Process and a variance based density function is employed to develop an expected-value function. A locational optimization based effective algorithm is employed to solve the resulting minimization of the expected-value function. We designed a mutual information based strategy to select the most informative subset of measurements effectively with low computational time. Our experimental results on real-world datasets have verified the superiority of the proposed approach.

I. INTRODUCTION

Recent advances in micro-electro-mechanical systems and wireless communications empower wireless sensor networks (WSN) to play a key role in a wide range of modern smart technology applications [1]. In particular, mobile robotic wireless sensor networks have attracted a significant attention for monitoring the spatial phenomena due to its distributed spatial properties and the mobility aspects. For example, applications of this technology includes detecting and monitoring forest fires [2], monitoring air quality [3], and monitoring ocean ecology conditions [4].

With on-board sensors, wireless communication and the mobile ability, the mobile wireless robotic sensor networks are capable of providing services required not only for monitoring and exploring the environment but also for exchanging the information. In [5], Cressie proposed a Gaussian Process (GP) model, also known as Kriging interpolation technique, a non-parametric generalization of linear regression, for the spatial phenomena assuming it obeys a Gaussian distribution. It is possible to exhaustively learn a GP model from the sensor readings and then utilizes this model to predict the values at unobserved locations. Combining these we address the sensor placement problem using the mobile wireless robotic sensor networks. Here, sensing observations and locations of sensors are used to generate a GP model for the phenomena.

Recently, there has been considerable activity in studying the wireless sensor placement problem [6] - [11] and designing wireless sensor networks for optimal monitoring. Nevertheless, most of these works are primarily focused on selecting the best subset from among a set of candidate sensor locations which

are selected from the discrete space. For example, in [10], the authors proposed a greedy heuristic algorithm based on entropy based information-theoretic model. The premise behind the entropy approach is to minimize the uncertainty of conditional entropy of unobserved locations, given observations. However, owing to a static sensor network, as shown in [7], the entropy method tends to pick locations along the border of the interested space causing sensed information waste. To address the drawbacks of the entropy approach, Guestrin et al., [7] proposed a new method based on the mutual information. In this method a subset from a set of potential sensor locations is selected such that the mutual information between the selected subset and the rest of the space is maximal. In our previous work [11], a computationally efficient simulated annealing based approximately heuristic algorithm was proposed to solve the sensor selection problem. Nevertheless, solutions of the sensor placement problem, using both mutual information and our previous proposed approach, are not optimal as the subset of sensor locations chosen is fixed.

Motivating on the sensor coverage problem, the authors in [12] proposed an algorithm that allows sensors to concentrate around locations of high event density while satisfying the minimal area coverage constraints. In Cortes et al [13], the authors optimized the mobile sensor network locations with respect to a known event probability density. Their proposed technique guarantees to minimize a cost function locally.

In this paper, we propose to utilize a GP model to develop a variance based density function. The proposed locational optimization based algorithm employed in this work locates the optimal sampling paths for mobile sensors. These trajectories allow to significantly improve the quality of prediction and to reduce the model uncertainty simultaneously. Moreover, a strategy based on mutual information is designed to select the best subset of measurements from all observations. This selection substantially decreases the computation time in developing the density function. In certain circumstances characteristics of the density function may lead the sensors to trap in certain locations. In this case all the mobile sensors are trapped in stationary locations without any further improvement to the sensing quality. An effective technique is developed to drive the mobile sensors to circumcenters instead of moving them to the centroids of the sensor locations' voronoi partitions.

Organization: Section II presents locational optimization

problem. Section III reviews the basics of Gaussian Process regression. Section IV describes the sensor placement based on the locational optimization. A density function and an algorithm for sensor movements are also analyzed and synthesized here. Experimental results are provided in Section V. Section VI shows the conclusions.

Notation: Let \mathbb{R}_+ be a set of non-negative real numbers. The Euclidean distance function is defined by $\|\cdot\|$. We let P be a convex polytope in \mathbb{R}^d and $\text{Ve}(P)$ be a vertex set of P . The set of all robotic sensor locations is denoted $S = \{s_1, \dots, s_n\}$. In other words, the position of i^{th} sensor is denoted s_i . And we define p as an arbitrary position in P . For a set P , we let $\mathbb{P}(P)$ denote the collection of subsets of P . Other notations will be explained in due course.

II. LOCATIONAL OPTIMIZATION PROBLEM

In this section, we describe the basic concepts and results on locational optimization that are used in this work. Readers are referred to [13] and [14] for further details.

A. Voronoi Partition

Consider a network of n mobile robotic wireless sensors that are deployed in a known, convex polytope P . At the fixed locations, the sensors could be divided by the optimal partitions of P , which are a set of Voronoi partitions $\mathcal{V}(S) = \{V_1, \dots, V_n\} \subset \mathbb{P}(P)$ defined by, for each $i \in \{1, \dots, n\}$

$$V_i = \{p \in P \mid \|p - s_i\| \leq \|p - s_j\|, \forall j \neq i\}. \quad (1)$$

In equivalent words, V_i is the set of the positions of P that are closer to s_i than to any of the other points in S .

B. Centroidal Voronoi Tessellation

Take a density function (also called bounded measurable function) $\phi : \mathbb{R}^d \rightarrow \mathbb{R}_+$ into a consideration in a bounded environment of interest $P \in \mathbb{R}^d$. For this work, we only use the value of ϕ limited to P of interest. We can consider ϕ as a function of measuring the information of probability that the events happen over P . The greater the value of $\phi(p)$ is, the more significant the point p is. In this paper, we use a non-increasing and piece-wise continuously differentiable function $f_s : \mathbb{R}_+ \rightarrow \mathbb{R}_+$ as a sensing performance function. This function characterizes the utility at position p taken from i^{th} mobile robotic wireless sensor at the location s_i . One can easily see that the value of this function downgrades with the distance $\|p - s_i\|$ between p and s_i . Therefore, we let this function is defined by a quadratic function $f_s = \|p - s_i\|^2$.

Given a density function ϕ , and a sensing performance function f_s , we formulate the expected-value function over the region P as

$$\mathcal{H}(S) = \sum_{i=1}^n \int_{V_i} f_s \phi(p) dp$$

or

$$\mathcal{H}(S) = \sum_{i=1}^n \int_{V_i} \|p - s_i\|^2 \phi(p) dp. \quad (2)$$

From the notation of the Voronoi partition, we can easily be seen that

$$\begin{aligned} \text{minimize } & \|p - s_i\| = \|p - s_j\| \text{ for all } p \in V_j, \\ \text{subject to } & i \in \{1, \dots, n\}. \end{aligned} \quad (3)$$

Therefore, our task here is to minimize the locational optimization function (2)

$$\text{minimize } \sum_{i=1}^n \int_{V_i} \|p - s_i\|^2 \phi(p) dp. \quad (4)$$

In other words, this optimization problem is to identify an optimal robotic sensor network configuration that makes (2) minimum.

Carrying out the partial derivative of $\mathcal{H}(S)$ with respect to the i^{th} robotic sensor, we get

$$\frac{\partial \mathcal{H}(S)}{\partial s_i} = \sum_{i=1}^n \int_{V_i} \frac{\partial}{\partial s_i} [\|p - s_i\|^2 \phi(p)] dp. \quad (5)$$

Let us recall in Euclidean space. The generalized area and the first mass moment of a set $P \subset \mathbb{R}^d$ with respect to ϕ are specified by

$$A(P) = \int_P \phi(p) dp \quad (6)$$

and

$$FM(P) = \int_P p \phi(p) dp. \quad (7)$$

Therefore, the centroid (center of mass) of a set P is formulated by

$$C(P) = \frac{FM(P)}{A(P)}. \quad (8)$$

Moreover, the polar moment of inertia of P about $s \in P$ is defined by

$$J(P, s) = \int_P \|p - s\|^2 \phi(p) dp. \quad (9)$$

From [15], by using Parallel Axis Theorem, we can easily state that

$$J(P, s) = J(P, C(P)) + A(P) \|s - C(P)\|^2. \quad (10)$$

From (2) and (9), using the Parallel Axis Theorem again, we simplify (2) to

$$\mathcal{H}(S) = \sum_{i=1}^n J(V_i, C(V_i)) + \sum_{i=1}^n A(V_i) \|s_i - C(V_i)\|^2, \quad (11)$$

and (5) to

$$\frac{\partial \mathcal{H}(S)}{\partial s_i} = 2A(V_i)(s_i - C(V_i)). \quad (12)$$

Equation (12) demonstrates that the solution of the locational optimization problem (4) is the set of centroidal Voronoi configurations in P [16]. In other words, the optimal mobile robotic wireless sensor location is the centroid of its Voronoi cell.

III. GAUSSIAN PROCESS BACKGROUND

Consider a two dimensional robotic sensor network that has a \mathcal{Q} , $|\mathcal{Q}| = m$, set of possible locations, which provide point measurements of some physical quantities. If we deploy all mobile sensors in \mathcal{Q} , $S \subset \mathcal{Q}$, sensor readings can still represent the distribution of the physical quantity in the whole space. This can be effectively predicted using Gaussian Process (GP) [17]. Lets denote $\mathcal{Q} = [q_1, q_2, \dots, q_m]$ as a set of locations, and $\mathcal{Z}_{\mathcal{Q}} = [z_1, z_2, \dots, z_m]$ as corresponding random variables at these locations. A joint probability distribution is given by

$$P(\mathcal{Z}_{\mathcal{Q}} = \mathbf{z}_{\mathcal{Q}}) = \frac{1}{(2\pi)^{n/2} \Sigma_{\mathcal{Q}\mathcal{Q}}} e^{-\frac{1}{2}(\mathbf{z}_{\mathcal{Q}} - \mathbf{m}_{\mathcal{Q}})^T \Sigma_{\mathcal{Q}\mathcal{Q}}^{-1} (\mathbf{z}_{\mathcal{Q}} - \mathbf{m}_{\mathcal{Q}})}, \quad (13)$$

where $\mathbf{m}_{\mathcal{Q}}$ is the mean vector and $\Sigma_{\mathcal{Q}\mathcal{Q}}$ is the covariance matrix of random variables, $\mathcal{Z}_{\mathcal{Q}}$.

Rasmussen et al. [17] demonstrated that GP has a marginalization property, which implies for any subset, S , of \mathcal{Q} , the joint distribution on random variables of its locations is Gaussian. Moreover, GP is comprehensively specified by its mean function $\mathcal{M}(q) = \mathbb{E}[z(q)]$, and a symmetric positive definite covariance function $\mathcal{C}(q_1, q_2)$, often called kernel function. One of frequently used kernel function is squared exponential, i.e.,

$$\mathcal{C}(q_1, q_2) = \sigma_f^2 e^{-\frac{\|q_1 - q_2\|^2}{2l^2}}, \quad (14)$$

where $q_1, q_2 \in \mathcal{Q}$, σ_f^2 is the maximum allowable covariance, l is the bandwidth of the kernel, and $\|q_1 - q_2\|^2$ denotes the Euclidean distance between q_1 and q_2 .

We define $S \subset \mathcal{Q}$ as a subset that includes all the selected sensors locations. For any subset S , let \mathcal{Z}_S denote the collections of observations at locations in S . In addition, take into account $\mathcal{U} = \mathcal{Q} \setminus S$ as the set of all elements in \mathcal{Q} but not in S , and $\mathcal{Z}_{\mathcal{U}}$ is a vector of random variables over these unobserved locations. We will assume that sensor measurement has an additive independent identically distributed zero-mean Gaussian noise with variance σ_n^2 . It can be clearly shown that \mathcal{Z}_S and $\mathcal{Z}_{\mathcal{U}}$ are jointly Gaussian distributed as

$$\begin{bmatrix} \mathcal{Z}_S \\ \mathcal{Z}_{\mathcal{U}} \end{bmatrix} \sim \mathcal{N} \left(\begin{bmatrix} \mathbf{m}_S \\ \mathbf{m}_{\mathcal{U}} \end{bmatrix}, \begin{bmatrix} \Sigma_{SS} + \sigma_n^2 I & \Sigma_{SU} \\ \Sigma_{US} & \Sigma_{UU} \end{bmatrix} \right), \quad (15)$$

where \mathbf{m}_S and $\mathbf{m}_{\mathcal{U}}$ (Σ_{SS} and Σ_{UU}) are mean vectors (covariance matrices) of \mathcal{Z}_S and $\mathcal{Z}_{\mathcal{U}}$, respectively. Σ_{US} ($= \Sigma_{SU}^T$), are cross-covariance matrices between \mathcal{Z}_S and $\mathcal{Z}_{\mathcal{U}}$; and I is the $|S| \times |S|$ identity matrix.

In probabilistic terms, we derive the conditional distribution at predicted positions of \mathcal{U} , given \mathcal{Z}_S as follows:

$$\mathbf{m}_{\mathcal{U}|S} = \mathbf{m}_{\mathcal{U}} + \Sigma_{US}(\Sigma_{SS} + \sigma_n^2 I)^{-1}(\mathcal{Z}_S - \mathbf{m}_S), \quad (16)$$

$$\Sigma_{\mathcal{U}|S} = \Sigma_{UU} - \Sigma_{US}(\Sigma_{SS} + \sigma_n^2 I)^{-1}\Sigma_{SU}, \quad (17)$$

where $\mathbf{m}_{\mathcal{U}|S}$ and $\Sigma_{\mathcal{U}|S}$ are mean vector and covariance matrix of $\mathcal{Z}_{\mathcal{U}}$, given \mathcal{Z}_S . As a consequence, using observations at locations in set S , we can predict quantities at unobserved locations, \mathcal{U} . This process is described as the Gaussian regression approach [18].

The quality of prediction is generally measured by calculating the errors at unobserved locations. Therefore, the goal

is to drive robotic sensors to optimal locations $S \subset \mathcal{Q}$ so that it will minimize a certain measure of prediction error at unobserved locations, given observations \mathcal{Z}_S . A typically used function inspired by the notions of G-optimality from optimal design [5], [19] is maximum predictive variance. Specifically, it is given by

$$\begin{aligned} \mathcal{M}(\mathcal{U}) = & \text{maximize } \sigma_f^2 - \Sigma_{qS}(\Sigma_{SS} + \sigma_n^2 I)^{-1}\Sigma_{Sq} \quad (18) \\ & \text{subject to } q \in \mathcal{U} \end{aligned}$$

where Σ_{qS} ($= \Sigma_{Sp}^T$) is the cross-covariance column vector between \mathcal{Z}_S and z_q .

IV. LOCALATIONAL OPTIMIZATION BASED SENSOR PLACEMENT

A. Density Function

Recall that in section II, we employed the density function to describe the measurement of information or probability of the events over P . In this work, we developed a density function based on the variances of the random variables $\mathcal{Z}_{\mathcal{Q}}$.

Consider a location $p \in \mathcal{U}$, $\mathcal{U} \subset \mathbb{R}^2$. From (17), the predicted variance at location p , given observations generated by S , is specified by

$$\Sigma_{p|S} = \Sigma_{pp} - \Sigma_{pS}(\Sigma_{SS} + \sigma_n^2 I)^{-1}\Sigma_{Sp}, \quad (19)$$

where $\Sigma_{pp} = \sigma_f^2$, and Σ_{pS} ($= \Sigma_{Sp}^T$) is the cross-covariance column vector between \mathcal{Z}_S and z_p .

In this paper, we embedded P into a two dimensional space, and we define

$$\phi(p) = \Sigma_{p|S}$$

$$\phi(p) = \Sigma_{pp} - \Sigma_{pS}(\Sigma_{SS} + \sigma_n^2 I)^{-1}\Sigma_{Sp}. \quad (20)$$

The actual measurements substantially influence on the density function as the covariance matrix include the hyper-parameters that are learned from the actual sensor readings.. Therefore when mobile robotic wireless sensors get to new positions, they need to sample the environment to generate a new density function.

B. Computations

There are two key issues associated with the implementation of the proposed algorithm in real mobile robotic wireless sensor networks: (a) limited computing resources to continuously compute the density function (b) moving mobile sensors out of stationary locations due to undesirable characteristics of the density function. As it requires to compute the density function in real-time after each iteration and it depends on the number of observations, calculation of this density function becomes intractable with increasing number of sensors' readings.

1) *Mutual Information*: In order to address the first issue, we employed a method based on the mutual information, a concept from information theory [20], to select the most informative subset of measurements from among all of the potential sensor measurements. After that the selected subset is utilized to develop the density function. For instance, to find k best measurements from whole observations \mathcal{O} , it starts from an empty set of locations $\mathcal{A} = \emptyset$ and greedily adds placements in sequence until $|\mathcal{A}| = k$. The algorithm chooses the next point that produces the maximum increase of mutual information. More specifically, the mutual information between the subset \mathcal{A} and the rest of observations $\mathcal{O} \setminus \mathcal{A}$ can be formulated by

$$F(\mathcal{A}) = I(\mathcal{A}; \mathcal{O} \setminus \mathcal{A}). \quad (21)$$

Once $y \in \mathcal{O} \setminus \mathcal{A}$ is chosen and added to \mathcal{A} , the variation of mutual information can be calculated by

$$\begin{aligned} F(\mathcal{A} \cup y) - F(\mathcal{A}) &= \\ &= H(\mathcal{A} \cup y) - H(\mathcal{A} \cup y | \bar{\mathcal{A}}) - [H(\mathcal{A}) - H(\mathcal{A} | \bar{\mathcal{A}} \cup y)] \\ &= H(y | \mathcal{A}) - H(y | \bar{\mathcal{A}}), \end{aligned} \quad (22)$$

where $\mathcal{A} \cup y$ denotes the $\mathcal{A} \cup y$ and $\bar{\mathcal{A}}$ represents $\mathcal{O} \setminus (\mathcal{A} \cup y)$; $H(y | \mathcal{A})$ is the conditional entropy of y after observing \mathcal{A} . By maximizing $F(\mathcal{A} \cup y) - F(\mathcal{A})$, we will find the next point y to add into \mathcal{A} . Suppose that the variables in \mathcal{O} are distributed normally, then 22 can be calculated in detail by

$$\begin{aligned} F(\mathcal{A} \cup y) - F(\mathcal{A}) &= \\ &= \frac{1}{2} [\log(\sigma_{y|\mathcal{A}}^2) - \log(\sigma_{y|\bar{\mathcal{A}}}^2)], \end{aligned} \quad (23)$$

where $\sigma_{y|\mathcal{A}}^2$ ($\sigma_{y|\bar{\mathcal{A}}}^2$) is the variance of y , given \mathcal{A} ($\bar{\mathcal{A}}$).

2) *Circumcenter*: As discussed in section II-B, the centroid of a sensor's current location will be the sensor's new location in the next movement. Importantly, computation of the centroids substantially depends on the shape of the density function. For instance, if the shape of the current density function is quite similar to that of the previous one or the shape of the density function is nearly flat, the distance between current and new locations is trivial. This leads the mobile robotic wireless sensors to stationary positions, which inadvertently indicates that the mobile sensors have already converge to their optimal locations. In order to address this issue, we utilized an approach derived from a circumcenter notion (refer [21] for more detail). In the event of all mobile robotic wireless sensors are immobile during any iteration without reaching the criterion of sensing quality, mobile sensors are driven from current positions to circumcenters rather than to the centroids of their voronoi cells.

The circumcenter of a bounded set $D \subset \mathbb{R}^2$ is the center of the circle of minimum radius that contains D . Hence, the circumcenter is unique. The calculation of the circumcenter of a polygon $D \subset \mathbb{R}^2$ is to minimize the radius r of the circle centered at p subject to the constraints that the distance between p and each of the polygon vertices is smaller than or equal to r . Mathematically, the problem can be formulated as

$$\begin{aligned} &\text{minimize } r \\ &\text{subject to } \|p - d\|^2 \leq r^2, \text{ for all } d \in \text{Ve}(D) \end{aligned} \quad (24)$$

C. Sensor Movement

Let us define the dynamics of each sensor movement by employing the first order form:

$$\dot{s}_i = u_i. \quad (25)$$

Solving the locational optimization in section II with the density function built in section IV-A, we can find the centroidal Voronoi configuration for a robotic sensor network. As a result, we propose a control law for each mobile sensor as follows:

$$u_i = a(s_i - C(V_i)), \quad (26)$$

where a is a proportional control gain.

Applying this control input, sensors will move to the new locations that generate constantly the Voronoi partitions.

Algorithm 1 Pseudocode for Sensor Movement in Sensor Placement

```

1: Initialize: Proportional control gain  $a$ 
2: Robotic sensors start at random positions  $S$ 
3:  $S^{tot} = S$ 
4: Collect measurements
5: loop
6:   Learn hyperparameters
7:   Develop density function
8:   Generate Voronoi partitions  $\mathcal{V}(S) = \{V_1, \dots, V_n\}$ 
9:   Calculate the centroid  $C(V_i)$ 
10:  → set of all new centroids  $S'$ 
11:  Calculate  $d(i) = \|s'_i - s_i\|$ 
12:  if Max(d) < predefined value then
13:    Remove  $S'$ 
14:    Calculate circumcenter  $CC(V_i)$ 
15:    → set of all new circumcenters  $S'$ 
16:  end if
17:  if  $\|s'_i - s_i^{tot}\| < \text{predefined threshold}$  then
18:    Remove  $s'_i$ 
19:  end if
20:   $S^{tot} = S^{tot} + \text{rest of } S'$ 
21:   $S = S'$ 
22:  Drive sensors to new locations  $S'$ 
23:  Get measurements for rest of  $S'$ 
24:  Select  $k$  out of all  $S^{tot}$  using mutual information
25: end loop

```

The approach to implement the sensor movements in practice is described by algorithm 1.

In this paper, we consider the environments that are constant or slowly changing in a short period of time. In this interval, mobile sensors will collect data, compute the new locations based on the locational optimization and move to these new positions, which have high variances. These movements make the maximum predictive variance (18) to gradually decline. Iteratively, sensors will run on the optimal trajectories that result in minimizing the maximum predictive variance.

V. EXPERIMENTS AND RESULTS

We conducted experiments using a real world data set of an indoor temperature environment from

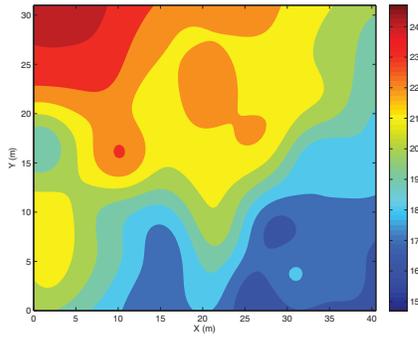


Fig. 1: Spatial temperature field

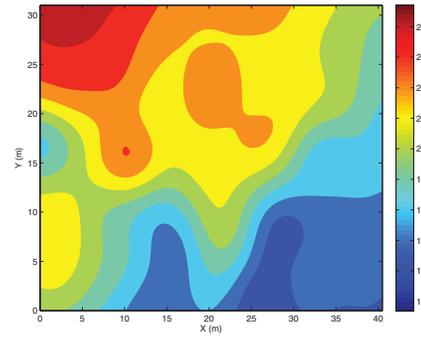


Fig. 3: Predicted temperature field at iteration 5

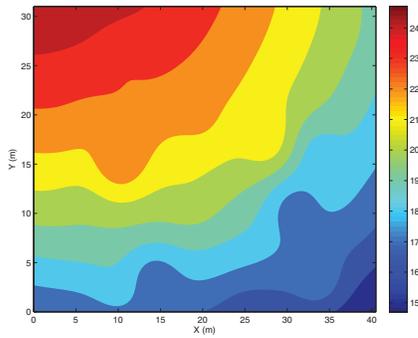


Fig. 2: Predicted temperature field at iteration 1

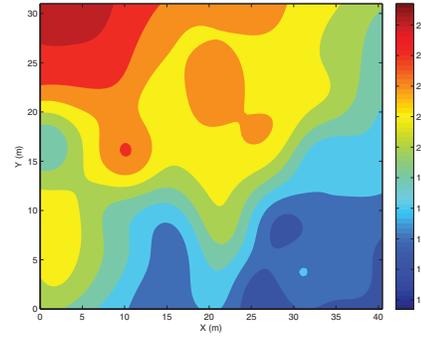


Fig. 4: Predicted temperature field at iteration 10

the Intel Berkeley Research Lab [7] (available at <http://db.scal.mit.edu/labdata/labdata.html>). Fig. 1 illustrates the spatial temperature field generated using this dataset with 52 sensors.

In this paper, in order to learn a Gaussian Process model, we first choose the isotropic square exponential covariance function and the second order trend for the mean function. One can recall that the covariance function is specified by the set of hyper-parameters (σ_f, l, σ_n) . Theoretically, these hyper-parameters could be learned by maximizing the log of the marginal likelihood (see [17] for more details). However, in this experiment, we used the `geOR` package [22] to learn these parameters.

In this experiment, we employ 20 mobile robotic wireless sensors to collect measurements iteratively, starting from random locations as shown in Fig. 5 (in white dots). In this implementation the time period of each iteration is dependent on the configuration of the sensor network, the computational time and the motion time. The predicted temperature fields are demonstrated from Fig. 2 to Fig. 4 at iterations 1, 5, 10. It can be clearly seen that the predictive means are getting closer to the true temperature field Fig. 1 as the number of observations increases. White circles in Figures Fig. 5 to Fig. 7 show the trajectories for the optimal paths of the mobile robotic wireless sensors at iterations 1, 5, and 10. Additionally, Figures Fig. 5 to Fig. 7 also show the prediction error variance corresponding to 1st, 5th and 10th iterations. Note that the white dots in these

figures correspond to the starting locations.

In order to evaluate the effectiveness of the proposed approach, we compare the maximum predictive variance performance metric for the proposed method and random motion method. In random motion method, mobile robotic wireless sensors move along random trajectories. Note that after each iteration, sensors move to new locations and sample the environment. Following that, hyperparameters are learned and variances at all unobserved locations are continuously updated. In Fig. 8, we can see that the maximum predictive variance is degrading gradually as the number of iterations increases as sensors have better knowledge of the environment. In particular, it can be clearly seen that the proposed algorithm exhibits improved performance continuously as each successive motion of sensor nodes is driven by the information gain.

VI. CONCLUSION

In this work we have considered the problem of prediction in spatial phenomena using observations obtained by a robotic mobile sensor network. We modeled the spatial field using Gaussian Processes and proposed a variance based density function for a locational optimization problem. As the number of observations increases, the proposed computationally efficient algorithm allows the mobile sensors to find the most informative locations in collecting measurements minimizing the prediction error variance. The demonstrations of results on real-world datasets have shown that our proposed approach

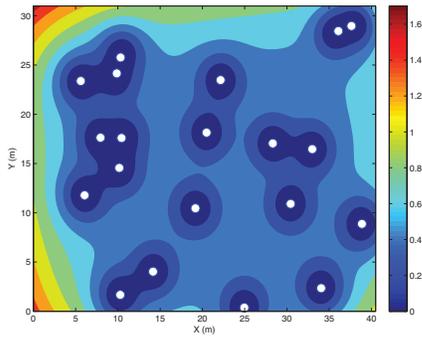


Fig. 5: Prediction error variances at iteration 1

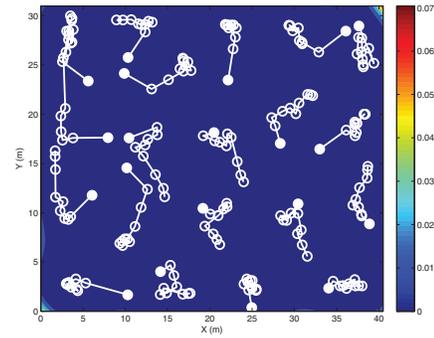


Fig. 7: Prediction error variances at iteration 10

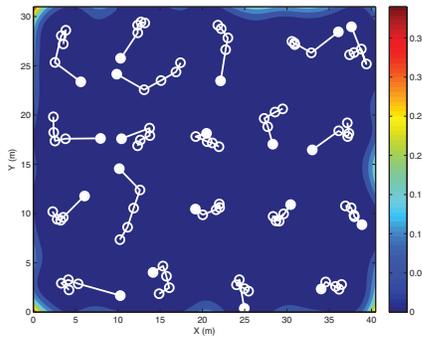


Fig. 6: Prediction error variances at iteration 5

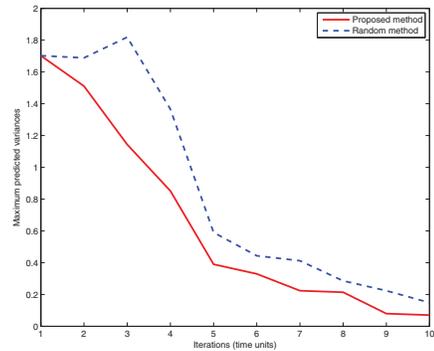


Fig. 8: Maximum prediction error variance

outperforms the random motion method.

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