Dynamic Modeling and Simulation of Ocean Plume for Robotic Control Design.

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Abstract— We present a model of the dynamics of an ocean pollution plume developed specifically for evaluating the performance and stability of robotic command and control algorithms. The key feature of our model is the ability to represent temporal and spatial sparsity as a model parameter. This approach provides a quantifiable, metric-based method for evaluating the affect of chemical concentration intermittency on the performance and stability of autonomous robot control algorithms.

I. Introduction

The study of plume models originated with early research to understand the dynamics of smoke rising from a chimney [2]. Over the 20th century, research has expanded to better understand the temporal and spatial distribution of chemical concentration outflows in a variety of fluid environments. The environmental representation we propose is based upon coastal ocean models which has been thoroughly studied by the oceanographic community. In this work we adapt these standard models to provide predictions of chemical concentrations that would be observed by autonomous marine robots operating in proximity of a pollution source. The purpose of this effort is to develop a virtual environment for evaluating heterogeneous teams of marine robots tasked with characterizing point-source pollution sources such as hydrocarbon spills and leaks; however, the model we propose is sufficiently general to capture the salient dynamics of a variety of pollution scenarios.

A critical aspect of the proposed environmental model is its ability to represent temporal and spatial variability as a parameter of the model. With sufficient averaging in time and space the distribution of chemical agents introduced in the marine environment can be considered to be smoothly distributed based on advection, diffusion, and dissolution. In such an idealized environment control algorithms can make use of observations of smooth gradients to achieve their particular goal, e.g., localizing the pollution source or following the plume front. However, previous studies on chemical plume tracking [14], [19] have demonstrated that the point measurements in such an environment are intermittent, suggesting significant temporal and spatial variability. These chemical *packets* or *filaments* lead to steep localized

gradient observations from mobile platforms, particularly in turbulent environments. Control strategies for multi-robot teams to characterize such environments must be robust with respect to this variability. Therefore the modeling approach we present includes an overt representation of this variability to enable quantifying the robustness of candidate control strategies.

Our research group previously proposed a controller to use with multiple heterogeneous robots to track and monitor an oil spill [15]. The controller is a modified gradient descent which enables the robot to follow the plume front, mapping the extent of the plume as it evolves over time. By testing this control approach with the model presented below we are able to analyze the robustness of such a multi-robot control algorithm to the sparsity of field observations. Furthermore we illustrate quantifying control performance with respect to the variability associated with plumes that exhibit the packet or filament type distributions as expected in field operations.

This paper will present a model that can capture both the time-averaged, idealized structure as well as the instantaneous, realistic structure of a dynamic plume as measured by on-board robot sensors. We describe two defining metrics of an instantaneous plume—the concentration sparsity and concentration peak/mean ratio—and show how to simulate the change from ideal plume to realistic plume by adjusting a single parameter of the model implementation. This method of emulating a pollution plume enables evaluation of the robustness of autonomous robotic sampling strategies with respect to observed variability; the proof-of-concept implementation of the Li et al. approach [15] is used to demonstrate this capability. For this illustrative test case we confirm the control approach performs well for when variability is low, the idealized case, and that the performance degrades as the variability increases. Furthermore we analyze how the performance degrades as a function of observation variability.

Section II-A of this paper details the plume characteristics most important for our use and briefly reviews the basic methods of modeling these characteristics. Section II-B describes other plume models in prominent use and compares their usefulness with respect to our objectives. Section III presents our proposed model as well as the specific implementation. Section III-B illustrates how the implemented model can capture both time-averaged and instantaneous characteristics of plumes and can be used to test robotic controllers. Section IV presents a case-study were the environmental model is used to evaluate the controller proposed by Li et al. [15].

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II. Background and Related Work

A. Fundamentals of Plume Modeling

Models have been developed for plumes in many different environments: atmosphere, water surface, and deep water for example. Over fifty different models specific to oil spills have been developed. Of these, a handful are still widely used [21]. The focus of this paper is models of ocean pollution that capture the temporal and spatial concentration evolution at the air-sea boundary. The dominant components of such models are advection (transfer of pollution by ocean current) and turbulent diffusion (transfer of pollution due to concentration gradients and shear flow). Secondary effects also play a role: evaporation, molecular diffusion, dispersion, emulsification, and interactions with ice and shore.

Evaporation can be an important component when volatile pollutants first reach the water surface and becomes less of an issue later when most of the materials have evaporated. Emulsification is the breakdown of large globules of pollutant into smaller uniform globules. Molecular diffusion is diffusion as a result of molecular properties particular to the pollutant. Interaction with the shore and ice is relevant only when the polluting event is near the coast or in the Arctic. Many of these models do not significantly effect the two metrics—sparsity and peak/mean ratio— or are only relevant in specific situations, such as when the considering cases near floating ice.

Plumes are often generalized as having smooth gradients where the concentration follows a normal curve. This generalization is a time-averaged representation of the concentration and location, usually of three minutes or longer [14]. The Gaussian model, based on partial differential equations, is explicitly designed to model the time-averaged plume [7]. The defining equation for the Gaussian plume model is

$$\frac{\partial c(x,t)}{\partial t} + v^T(x,t) \cdot \nabla c(x,t) = D\nabla^2 c(x,t) \qquad (1)$$

where v is the velocity vector of the environment flow, $\nabla c(x,t)$ is the concentration gradient, $\nabla^2 c(x,t)$ is the concentration divergence at point x and time t and D is a diffusion coefficient of the material.

The Gaussian model is the oldest of the plume modeling techniques and is used extensively to model atmospheric plumes. Because the Gaussian is a time-averaged model, it idealizes some of the more interesting characteristics a plume can have, specifically, the features of an instantaneous plume.

We use two main metrics to quantify the instantaneous characteristics of a plume: the sparsity and the peak/mean ratio.

Definition 1 (Sparsity): The sparsity is defined as the percentage of concentration readings in a stationary point over a certain threshold as a plume migrates over that point.

Sparsity is calculated as the cardinality of the set of concentration readings at a point under a defined threshold, over the cardinality of the set of all samplings at the same point. Sparsity can be modeled as

$$T_x = \{t \in S_x | \ t < \theta\} \qquad (2)$$

$$\text{sparsity} := \frac{|T_x|}{|S_x|} \qquad (3)$$

where S_x is the set of all recent samples and T_x is the set of all samples above a threshold θ [14].

Definition 2 (Peak/mean ratio): The peak/mean ratio is defined as the ratio between the peak concentration value and the time-averaged mean of concentration at a stationary point as a plume migrates over that point,

$$\text{peak/mean ratio} := \max(S_x)/\frac{1}{|S_x|}\sum S_x. \qquad (4)$$

where $\max(S_x)$ is the maximum concentration reading in the set of recent samples, S_x, and $\frac{1}{|S_x|}\sum S_x$ is the mean concentration reading across the same set of samples [4].

Jones showed that for some kinds of plumes, the sparsity at any given location can be upwards of 90% [14]. That is, for 90% of the time, a stationary sensor will read zero (or below a sensing threshold) as the plume moves through the point. A study by Crimaldi et al. found that peak concentration can be three orders of magnitude larger than the mean [4]. These metrics are shown qualitatively in Fig. 1. This figure was generated using the simulation described below and will be used to compare the validity of our model with the metrics defined by Jones and Crimaldi. Notice the peak over 50 samples is normalized to 1 and the mean is almost exactly $1/3$. Qualitatively, the majority of the readings are below the mean.
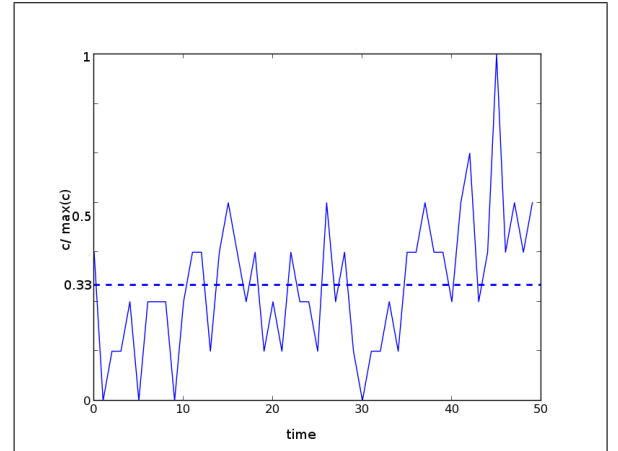

Fig. 1: The Simulation sparsity snapshot

Since the instantaneous concentrations are of great concern for robots navigating a plume, the Gaussian model alone is not sufficient to test control algorithms that rely on instantaneous, fine scale readings of a plume.

The Lagrangian method is a particle based model that depicts the plume as an aggregate of many individual particles, each moving according to the advection laws and a random walk derived from the Gaussian equations. In the limit, as the number of particles approaches infinity and the time step approaches zero, the Lagrangian model approaches the Gaussian model, but when it is far from the limit, the Lagrangian model has the structure very similar to those

metrics described by Jones and Crimaldi. As a result, this method much more capable of representing the instantaneous dynamics of a plume.

The generic defining equation for the Lagrangian random walk method is

$$x_i(t + \Delta t) = x_i(t) + \text{adv}\,(t, x_i(t), \Delta t) + \text{diff}\,(t, x_i(t), \Delta t) \tag{5}$$

where $x_i(t)$ is the location of the ith particle at time t, Δt is the time step, $\text{adv}\,(t, x(t), \Delta t)$ is the advection component and $\text{diff}\,(t, x(t), \Delta t)$ is the diffusion component of the particle movement. Many different ways to model the advection and diffusion equations exist. Usually advection is determined by the velocity vector of the fluid field at location $x(t)$ and time t and the diffusion is a function of the diffusion coefficient for the material and a random variable. Many other higher order and situational effects can be modeled using the Lagrangian method by superposing their contribution to the particle movement or periodically removing individual particles from the map.

B. Related work

The Lagrangian model is now the dominate model used for studies the dynamics of marine pollution. Such fate models focus on the distribution of the pollutant after a certain amount of time. These models use location specific environmental data for currents and wind. Many derivations of the Lagrangian model have been proposed and implemented [21].

General NOAA Operational Modeling Environment (GNOME) is a widely accepted tool produced and verified by NOAA [1]. This software tool uses historical weather data to model the fate of the oil spill. GNOME can be very useful for simulating the transport and fate of oil spills in specific places around the world. It comes equipped with historical current and wind data and can simulate oil spills of many different kinds of oil. The main concern with the GNOME software model is that the step size is on a scale much to large to accurately reflect the instantaneous dynamics of a plume necessary for evaluating robotic remediation strategies. GNOME has a minimum resolution of step-size of 15 minutes. A European model MEDSLICK provides detailed simulation of oil spills very similar to GNOME [5] [6]. It suffers from the same problem as GNOME of using a time and space scale much too large for testing of robotic response.

Researchers have tackled the problem of understanding and containing oil-spills from a number of different perspectives. In the robotics community, researchers study source localization, edge tracking, and autonomous clean-up, with single and multiple robots. Those who study plume source localization generally base their approach on biomimetic techniques, trying to understand and emulate how animals find plume sources [10] [16] [22]. Other strategies focus on the problem by looking at optimal ways to layout robotic sensor networks for spill tracking and monitoring of the plume edge [11] [15] [25] [20]. Some strategies specifically

focus on robotic clean-up and how to generate complete mapping of the dynamic plume [13] [24], and spill finding [17]. Additionally, many of the techniques have been developed for individual and multiple robots [11] [15] [17] [24].

Theoretical development of the above techniques usually use either the Gaussian or Lagrangian model for simulation. Marjovi and Marques use a Gaussian plume model for development of their algorithms for swarm odor plume finding [17]. Sahyoun et al. also use a Gaussian plume model to develop algorithms for plume tracking with mobile sensors [20].

Farrell et al. have attempted to capture the instantaneous structure of plume behavior using a variation of the Lagrangian method [9] [10]. This approach simulation is suited for capturing the instantaneous dynamics of an odor plume and has led to control strategies that emulate how moths find the source of an odor plume [9]. A similar model is used by Li et al. to test source localization algorithms based on moth behavior [10] [16]. Zarzhitsky et al. also used the model developed by Farrell to simulate swarm chemical plume tracing [9] [24] . This model is not applicable to oil spills because it does not model large-scale turbulence and thus presents a largely laminar plume. Though it emulates the metrics described by Jones and Crimaldi, it is incapable of also emulating the idealized plume described by the Gaussian equations.

III. PLUME MODEL AND SIMULATION

A. Model Implementation

Many models for the advection and diffusion component of the Lagrangian method exist. In this paper we use an adaptation of the equations derived by Chao and verified experimentally by Nagheeby et al. [3] [18]. This model accounts for many effects on the plume besides advection and diffusion, such as evaporation, dissolution, shoreline and ice interaction, and emulsification. In the interest of brevity, the examples we discuss below are constructed for a scenario where advection and diffusion are dominate.

We implemented the Lagrangian model (5) in the paper with the x and y components of the advection and diffusion components are as follows. Eqs. (6)-(9) show the advection and diffusion the equations used in this simulation,

$$\text{adv}_x\,(x, t) = k_t u_x + k_w w_x \tag{6}$$
$$\text{adv}_y\,(x, t) = k_t u_y + k_w w_y \tag{7}$$
$$\text{diff}_x\,(x, t) = R_1 \,(12 D \Delta t)^{1/2} \cos(2\pi R_3) \tag{8}$$
$$\text{diff}_y\,(x, t) = R_3 \,(12 D \Delta t)^{1/2} \sin(2\pi R_4) \tag{9}$$

where u and w are the current and wind vectors respectively, k_t and k_w are constants used to capture the impact of the current and wind. Chao recommends these be set to 1 and 0.03 respectively. R_i is a unique uniform random variable on the interval $[0, 1]$. D is the diffusion coefficient, typically set between $0.01 m^2/s$ and $0.5 m^2/s$.

We implemented the above model in Python as a standalone process. This allowed us to use the real-time control

software processes by having the environmental module pass simulated sensor measurements to the controller via inter-process communication. The message passing between processes is implemented using the Lightweight Communications and Marshalling (LCM) library [12]. This work flow enables rapid transition between controller development (in offline and real-time simulation) and field deployment on sensor-equipped mobile robotic platforms (real-time control).

Fig. 2 shows an example of the simulated plume evolution. This density histogram shows the results of the plume emitted from a source located at $(12, 26)$ at $2, 5, 8,$ and 11 hours after the pollution begins.
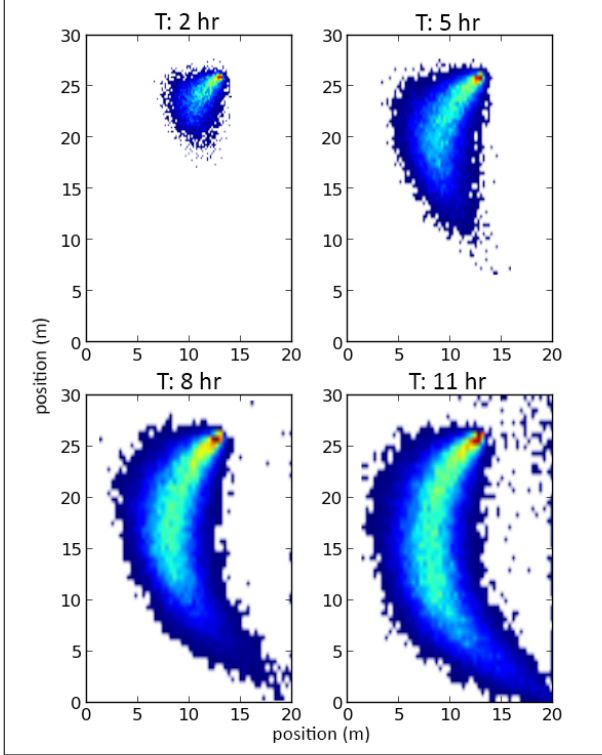

Fig. 2: Two-dimensional histogram of plume concentration at 2, 5, 8, and 11 hours after release

Calculating the concentration is done by counting all of the particles within a certain radius from the point in question,

$$c(x, t) = \text{card}(\{m | m \in B(x, r)\}) \qquad (10)$$

where the concentration at location x and time t is the cardinality of the subset of particles inside a ball centered at x with radius r, $B(x, r)$. For our implementation, we chose the ball using the L_∞ norm.

B. Parameterizing Intermittency

We consider the time-averaged plume and the instantaneous plume as two extremes of what our plume model should be capable of simulating. On one hand, the smooth gradients are completely idealized; on the other hand, the instantaneous plume is extremely sporadic and irregular. The simplicity of the Lagrangian method modeling the plume as composite of many individual particles allows the model to

reach near the Gaussian time averaged plume and near the instantaneous plume, and everywhere in between. Also, the Lagrangian model uses a moving frame of reference, and thus only calculates the concentration of a point based on the particles that are nearest that point, no interpolation and smoothing between gridded blocks occurs in the simulation. This allows for the simulation to produce extremely varying fluctuations in concentration readings over both space and time [4] [14] [23] .
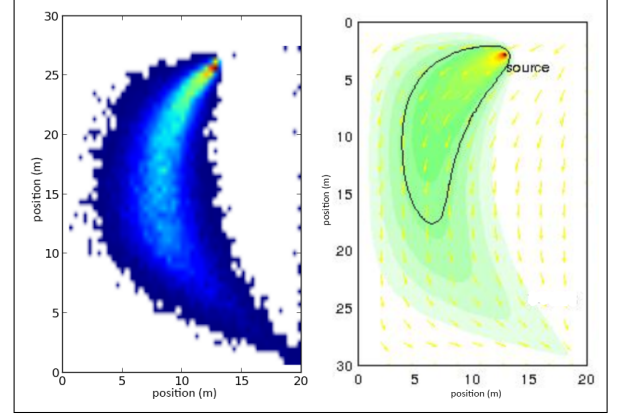

Fig. 3: The Lagrangian model on the left compared with the Gaussian model on the right

In the limit, as the time step approaches zero and the number of particles approaches infinity, the model approaches the Gaussian model. Fig. 3 shows a concentration map of the Lagrangian particle plume generated in this study side-by-side with the Gaussian model originally used to develop the controller under study [15]. This shows qualitatively how similar the two models can be from a macro perspective.

The sampling radius of concentration is used to adjust the intermittency, as measured by the sparsity and peak/mean ratio. For the physical instantiation of this example, we can identify numerical values for the model parameter that yield useful environmental measurement predictions. If the sampling radius is set to 0.5 m the plume. as observed by robotic sensors, will have low variability and seem similar to the idealized Gaussian plume. If the sampling radius is reduced to 0.05 m the predicted plume measurements generated by the model exhibit the sparsity and peak/mean ratio we might expect in field measurements.

Fig. 4 shows the concentration profile over time at a location downstream from the source of the plume. When the sampling radius 0.5 m, the concentration is nearly constant. When the sampling radius is very low, 0.05 m, the concentration profile is very similar to how Jones and Crimaldi described the structure of an instantaneous plume [4] [14]. This graph shows that far from the limit the simulation represents metrics similar to the experimental sampling represented in Fig. 1 suggesting that by varying the sampling radius the model can represent the variability one would expect from field robot experiments.

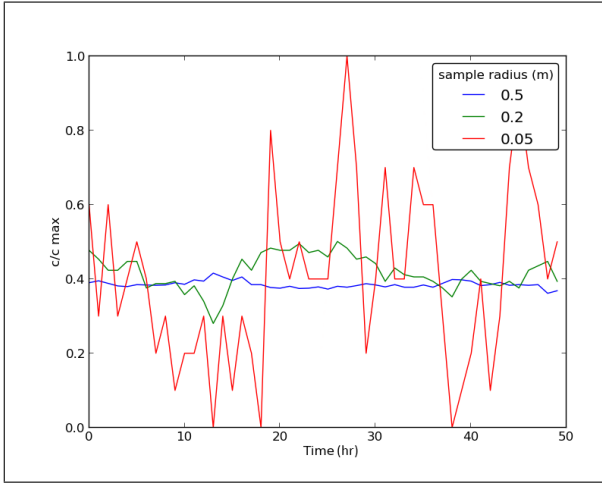Fig. 5 shows the spacial distribution of the plume con-

Fig. 4: Temporal concentration profile for a single location. The location is along the minor axis of the plume 6 m from the source. The y-axis show the concentration normalized by the maximum concentration (c/c_{max})
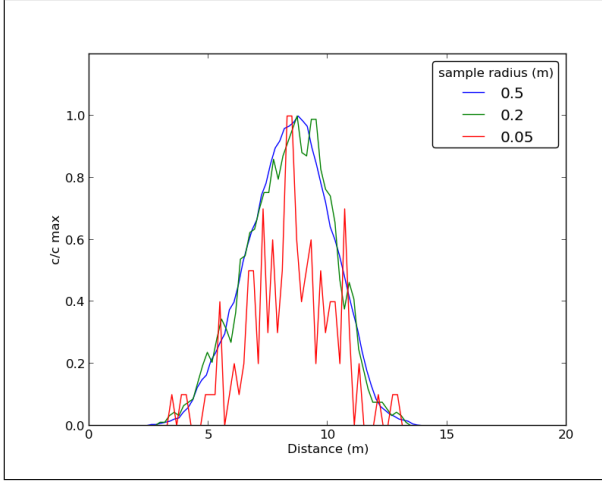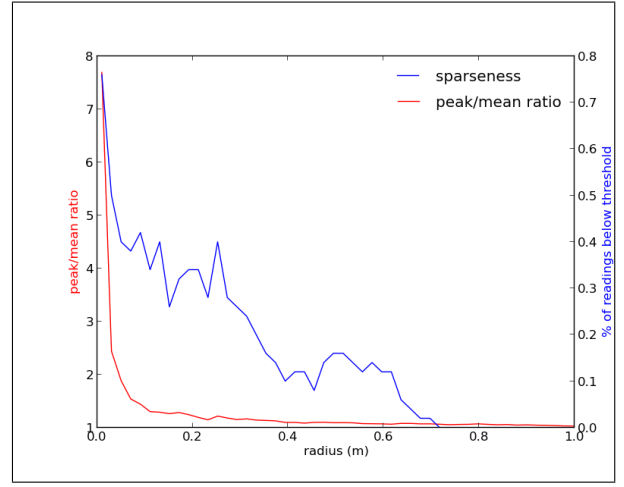

Fig. 6: Sparsity and Peak/mean radius as a function of the sampling radius. The peak/mean ratio values are shown on the right y-axis and the sparsity metric is shown on the left y-axis.

Fig. 5: Spatial variability of predicted concentration measurements as a function of the sampling radius model parameter. the profile is generated for a line parallel to the minor axis of the plume 6 m from the source. The y-axis show the concentration normalized by the maximum concentration (c/c_{max})

centration along the minor axis 6 m down stream from the source. When the radius is at 0.5 m, the concentration profile nearly follows a normal curve, like what we would expect in the limit. When the sampling radius is set to 0.05 m the concentration profile is very irregular. Though it generally follows a bell curve, it has a very large variation. The middle line at 0.2 m shows a point between the two extremes.

Fig. 6 shows the sparsity and peak/mean ratio as the sampling radius moves from 0 to 1. This shows how tuning the radius effects the sparsity and the peak/mean ratio. Additionally, by adjusting the threshold value, we can bring these two curves closer together.

IV. CASE-STUDY

To illustrate the utility of proposed environmental model we present the evaluation of the control algorithm developed in [15]. The goal of the controller is to track the dynamic plume front based on concentration measurements made by mobile robots. This control approach considers the local gradient, but also includes a transport model (advection and diffusion) in the estimation and control framework. Previous work, using a simulated Gaussian plume environmental model (low variability), has shown the controller successfully tracks plume front movement at the water surface. For this case-study we considered the single robot case, but the algorithm can be extended to multiple mobile robots using nearest-neighbor communication. For the details of the controller implementation in the robotic simulation platform, please refer to our companion paper [8].

Two test cases are discussed below: the first with low variability (approximating the Gaussian plume) and the second with high variability (approximating the packet or filament concentration gradients observed in other ocean experiments). For both cases, the control objective is to follow a particular concentration contour (specified as 30 % of the maximum expected concentration). For both cases the underlying model parameters specifying advection and diffusion behavior are identical; the only difference between the two test cases is the value used for the model parameter that controls variability.

The results of the first case, low variability, are illustrated in Fig. 7. For this case the concentration sampling radius of the environmental model is set to 0.5 m to approximate the Gaussian plume model through introduction of spatial smoothing. This value was chosen for the sampling radius because it produces a nearly smooth concentration curve, but still retains some variance. The control strategy successfully guides the mobile robot in a path to follow the plume

front by tracking the gradient contour. These results are consistent with the findings reported in [15] suggesting that the proposed model approximates the Gaussian plume model when the concentration sampling radius is sufficiently high.
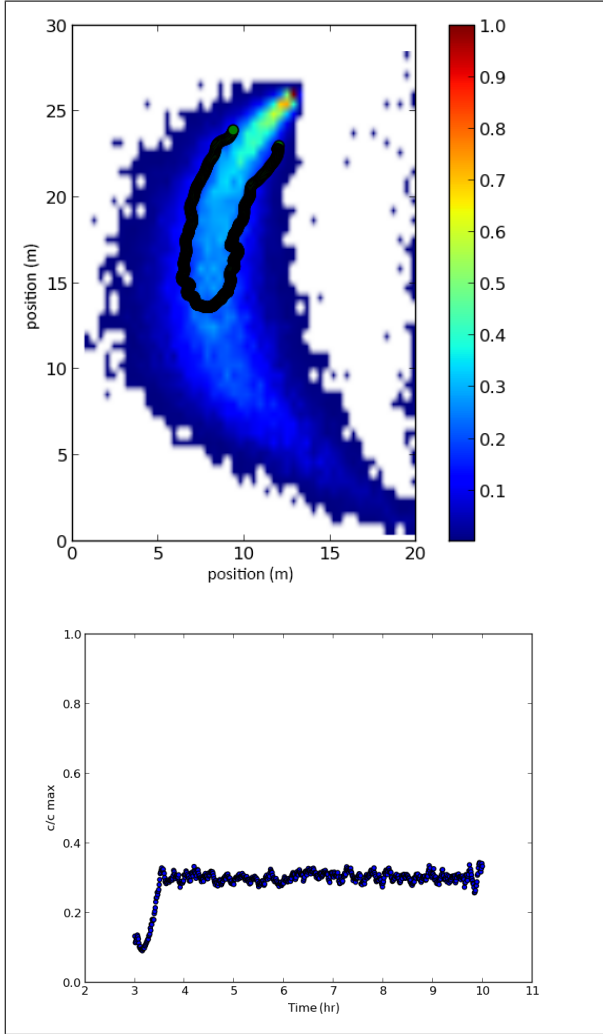

Fig. 7: Results for the low variability test case for single robot tracking the plume front. The upper image shows the chemical concentration of a pollution plume at the end of the simulation run. The control objective for the mobile robot is to follow the 30% contour. The robot track is shown superimposed on the plume as a black line. The lower graph shows the time history of the concentration observations made by the mobile platform.

For the second case we introduce variability that we might anticipate when we transition the control algorithm to field implementation, i.e., testing the algorithms on mobile robots operating in a coastal ocean environment using on-board sensors to observe pollution concentration. Based on previous experimental work we can expect significant variability in these instantaneous measurements. To prepare for this transition from simulation to experiment it is necessary to evaluate how the control approach performs in a more authentic environmental scenario, not just in

the idealized Gaussian plume case. To accomplish this we reduce the concentration sampling radius to $0.05\,\mathrm{m}$ which still captures the underlying advection and diffusion, but also emulates the intermittency and peak/mean ratio described by Jones and Crimaldi. The results are shown in Fig. 8, which illustrates that the increased variability prevents the algorithm from guiding the robot to follow the plume front. The robot track shows that the robot moves very little from its original position, suggesting that the reliance on gradient measurements is brittle with respect to the localized, quickly varying concentration measurements typical of experimental implementation. The concentration readings illustrate the intermittency we might expect from an on-board sensor where the concentration at any given point is likely to be zero because of the packetized pollution. While this result shows that the original control algorithm is not robust to this intermittency, it does illustrate the capability of the environmental model we propose to capture this effect. Furthermore we can vary a single parameter, the concentration sampling radius, to smoothly transition from the idealized Gaussian plume case to a more authentic case that captures the additional complexity of an intermittent plume environment observed in ocean experiments.

V. CONCLUSIONS

We present an environmental model of pollution plumes in an ocean environment. In addition to the underlying advection and diffusion dynamics, the model explicitly represents variability and intermittency in pollution concentration measurements to enable evaluation of robotic control algorithms. The parametric representation allows the evaluation of control algorithm robustness with respect to variability. To successfully transition algorithms developed in simulation it is critical to understand how these algorithms will perform with the intermittent sensor measurements to be expected in field deployments.

We illustrated the utility of this approach through a case-study. In this case-study we evaluated the performance of a previously proposed control algorithm using the proposed environmental model. We showed that with the variability set sufficiently low, by increasing the size of the concentration sampling radius, the model approximates the Gaussian plume model. As variability is increased the performance of the control algorithm degrades. Understanding the relationship between variability and performance is critical to transitioning these algorithms into the field.

Field experiments in marine environments are costly and time consuming, therefore it is critical that the simulations used to evaluate their efficacy capture the salient aspects of ocean operations. Previous research has shown that the real-time sensor observations of pollution plumes in such environments are highly variable; to be experimentally successful robotic control algorithms must address concentration intermittency. The model we propose provides a method for emulating this intermittency to quantify performance of candidate robotic algorithms tasked with supporting ocean
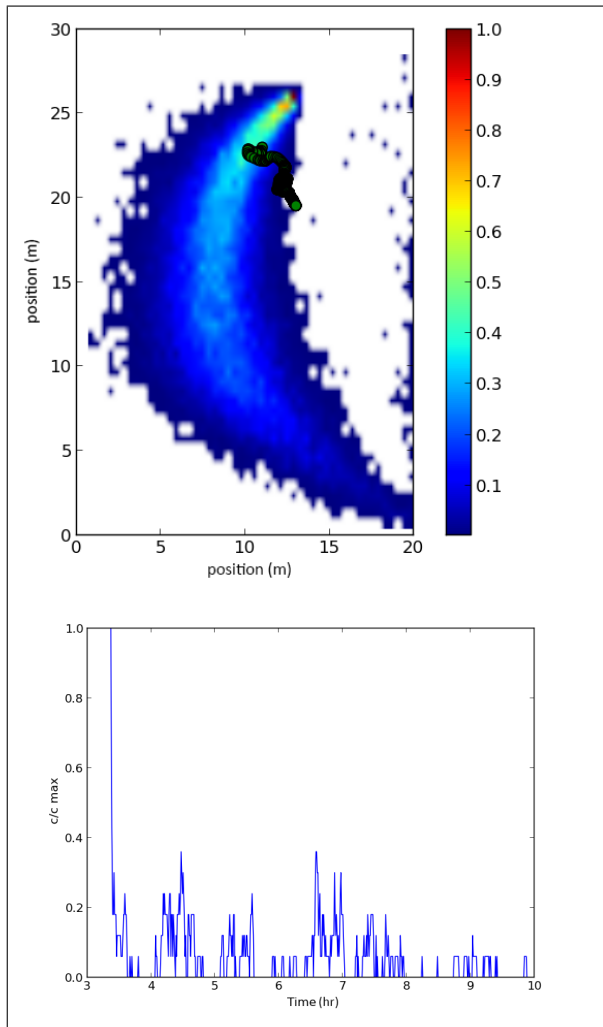
Fig. 8: Results for the high variability test case. In the upper image the robot location is shown superimposed on the plume image. The lower graph shows the time history the concentration observations made by the mobile platform, illustrating the variability in instantaneous observations.

pollution remediation. In application ocean pollution dynamics have a wide variety of characteristics. As researchers transition robotic algorithms to work in such environments it is important to predict how these strategies will perform when implemented on operational hardware. The proposed model allows developers to evaluate the performance in a wide variety of test cases to better quantify sensitivity to variability in sensor measurements before incurring the costs associated with experimental deployment.

VI. Acknowledgments

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