Data Sharing and Assimilation in Multi-robot Systems for Environment Mapping

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Abstract: We consider scenarios where a mobile multi-robot system is used for mapping a spatial field. Gaussian processes are a widely employed regression model for this type of tasks. For the sake of generality, scalability, and robustness, we assume that planning and control are fully distributed and that robots can only communicate via range-limited channels. In such scenarios, one core challenge is how to let the robots efficiently coordinate in order to maintain a shared view of the mapping process, and, accordingly, make plans minimizing overlaps and optimizing joint information gain from obtained measurements. A simple approach of sharing and utilizing all the sampled data would not scale to large teams, neither for computation nor for communication (assuming a general ad hoc robot network). Building on previous work where robots adaptively plan where to sample data by selecting convex containment regions, we propose a data sharing and assimilation strategy which aims to minimize the impact on communication and computation while minimizing the loss on accuracy in map estimation. The strategy exploits convexity of the regions to create compact meta-data that are locally shared. Submodularity of information processes and properties of GPs are used by the robots to create highly informative summaries of the sampled regions, that are shared on-demand based on the meta-data. In turn, a received summary is assimilated by a robot into its local GP only if/when needed. We perform a number of studies in simulation using real data from bathymetric maps to show the efficacy of the strategy for supporting scalability of computations and communications while guaranteeing learning accurate maps.

SCIENCE AND TECHNOLOGY PUBLICATIONS

1 INTRODUCTION

The availability of *spatial maps describing the state* of an environment with respect to values of interest is essential for the timely modeling monitoring of its physical, chemical, and biological properties. For instance, in aquatic environments¹ there's a clear interest in having accurate maps describing values such as depth, salinity, pH, pollution, temperature, presence of flora and fauna, etc. Typically, these quantities of interest have a continuous distribution over the spatial field, which can extend over several kilometers. Given the extension of the regions of interest, the use of autonomous mobile robots, and in particular of *mobile multi-robot systems* (MRSs), is strongly envisaged to effectively tackle the tasks of *exploring* an environment and/or constructing data maps (Stachniss, 2009). In a given time budget, a well-designed MRS can perform measures in multiple sites in parallel, potentially providing a large speed-up in map construction and boosting map accuracy. In this work, we precisely start from the use of MRSs for environment mapping missions: in a limited time budget, an accurate map of the given region and attributes of interest must be constructed by exploiting robot mobility and robots sensors providing localized measures (i.e., the range of the sensor is very narrow, fundamentally point-like, which is typical, for instance, for measuring things such as salinity, pH, concentration of chemicals). Moreover, for the sake of generality, we assume the use of an ad hoc network for robot communications.

In such scenarios, measures of interest are defined over a continuous field, while robots can only take samples at discrete points, asking for building *regression models*. Moreover, accounting for sensor noise and errors, as well as for the limited time budget available to cover a vast region, the number of

514

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¹In this work we are precisely motivated by performing *automatic monitoring and surveying in marine environments*. Therefore, we mostly refer to this type of scenarios. However the methods presented in the rest of the paper are general and are not restricted to mapping tasks in aquatic environments.

sampling points is necessarily limited. Therefore, a main challenge is to *plan in the MRS where to sample* and how to combine the discrete and noisy measures into a regression model providing a reliable, continuous map estimator at the end of mission. Planning is expected to be adaptive, to account for incrementally gathered evidence measures. Moreover, localized sensors and *submodular* properties of information environments (Krause and Guestrin, 2011) partially rule out the use of approaches of intelligent partitioning of the environment (e.g., (Cortes et al., 2004)), asking, in general, for complex and dynamic schemes for planning.

In the described scenario, robots need to coordinate to issue plans that can jointly perform an optimized sampling process. Ruling out, for the sake of generality and scalability, the use of a centralized controller, the realization of an optimized and distributed scheme requires exchanging sampled data in order to maintain a *shared map* and plan accordingly (e.g., to direct sampling to unexplored regions, or to those regions where estimation uncertainty is higher, or, equivalently, where predicted information gain is higher (Ma et al., 2018)). What data to share and how to use (i.e., assimilate) the shared data in a distributed MRS for planning purposes is the main focus of this work: we will introduce a novel strategy for the selection of data to share and for data assimilation in distributed MRSs. Note that data sharing and assimilation must ensure scalability of communications (i.e., make a parsimonious use of available bandwidth) and of on-board computations (i.e., allow near real-time computations), while guaranteeing high accuracy in final map estimation. In general, use and control of data sharing is a central issue in MRSs, being involved in virtually any task of interest (e.g. (Roumeliotis and Bekey, 2002; Jager and Nebel, 2001)).

Since the importance of each piece of data that can be shared depends in turn on how data is used to build a regression model, we consider here Gaussian Processes (GPs) (Rasmussen, 2003) as the reference model to learn the target mapping. This is a specific yet general choice since GPs are widely employed in the relevant robotics literature for environmental data mapping (Ma et al., 2018; Li et al., 2020; Popović et al., 2020; Ghassemi and Chowdhury, 2019). While providing a number of advantages (e.g., in terms of capturing the statistical properties of spatial correlations), a main drawback when using GPs is related to the $O(n^3)$ complexity for each model updating, with *n* being the size of the available data measures. This makes it computationally expensive to run in real-time on embedded platforms such as mobile robots, as the data size grows. Moreover,

being non-parametric models, GPs put a major burden on data sharing, given that the model cannot be necessarily compacted in a few parameters.

In this work, using GPs as a target model, we provide a *strategy to perform an intelligent selection and assimilation of the sharing data in an MRS*. The goal is to limit the exposure of communicated and processed information for the sake of scalability, while still allowing the robots to effectively plan their sampling actions. In order to frame the data sharing and assimilation strategy in a concrete scheme for GPbased planning, we build on our previous work (Di Caro and Yousaf, 2021b), where robots adaptively and iteratively identify convex containment regions where to sample at selected locations.

The proposed strategy makes the robots exploiting the convexity of the regions to create compact meta-data descriptions, that are locally shared consuming little bandwidth. Submodularity in contiguous regions and statistical properties of GPs are exploited to let each robot independently create compact and informative data summaries of sampled regions, that are shared with other robots on-demand based on meta-data and estimated utility. In turn, a received data summary is assimilated by a robot into its local GP only if/when needed. Overall, the use of compact descriptions and data summaries, as well as the use of opportunistic communications and computations fully support scalability of the processes and ensure a minimal, strategically tunable, loss in final map accuracy. We show these properties through a number of simulation studies.

2 RELATED WORK

A central aim of our work is to define a strategy to let the robots select and share a 'minimal enough' amount of information, in the form of GP data, to enable informed and coordinated planning in the MRS while ensuring scalability. In the following we review a few works that similarly aim to mitigate the burden of data either in single robot or multi-robot systems.

(Ma et al., 2018) utilizes the concept of Sparse Online Gaussian Process (SOGP), to limit the size of the training data for GP. Their approach is well suited for single robotic exploration case where each data point is tested online to determine if GP could still be modelled with existing data point i.e. without increasing the size of variance-covariance matrix but with re-normalized kernel parameters induced by new data point. This is a cheaper alternative as it limits the growth of GP, it is however not suited in the case of multi-robots as there is no notion of 'localized summaries' that enable other robots to form decisions to either consider the batch of data or reject it based on their current budgets and location of the summaries.

In (Choudhury et al., 2002) the burden of data is mitigated by forming multiple clusters to model multiple GP's. Different sets of hyperparameters are used to model the compactly supported covariance structures of different clusters. The clustering mechanism is based on maximizing the log likelihood for each cluster, and using a predetermined partitioning of size *m*. Although the approach is well suited for addressing data's computational load, the data size still remains constant within each cluster, and the clustering process is done after the data collection process, hence not suitable for online exploration.

(Chen et al., 2013) defines the *p*PIC algorithm where each machine (e.g., individual robot) compute a local summary and provide it to master, while master computes the global summary and distribute over the machines. The machine makes use of local data, local summary and global summary to produce predictions over unobserved locations. They test their algorithm for computation, time and communication complexity (Ghassemi and Chowdhury, 2019) provides a method of decentralized data sharing where observations of individual robot is downsampled to fixed 400 observation if the size of data grows more then 400 observations by each robot. Although this can be assumed as naive approach, the exact downsampling technique isn't clear.

The approaches mentioned above aim to either decentralize the GP such that it that can be distributed and run over several machines in parallel or downsample the data to mitigate the burden of GP computation, or both. Our aim is different, as we seek to summarize data not specifically for highly accurate predictions, but only to sufficiently produce accurate predictions that can enable robots to plan to sample in informative regions: summaries are only accurate enough to "mimic" similar behavior among the robots as if all the data would be shared among them.

3 REGION-BASED PLANNING

As mentioned in the Introduction, we build on our previous work (Di Caro and Yousaf, 2021b) where we have introduced an adaptive sampling approach for environmental mapping tasks using MRSs and GPs. The methodology is based on an *iterated multi-stage process*: (i) each robot, or group of robots, uses Monte Carlo to perform Bayesian inference on the currently estimated GP model with the goal of identifying the *best region where to sample new data in the next time*

slice, where best here means the region maximizing the expected information gain for GP learning; (ii) inside the region, Bayesian optimization is employed to select a set of candidate locations where to go sampling; (iii) an *orienteering model* is used to select a subset of the locations such that the overall expected utility is maximized and a path that satisfies given time constraints is computed through them; (iv) each robot then follows the computed *plan*: it goes to its selected region and performs sampling over the selected path. Process (i-iv) is *iterated* and *fully distributed*.

For computational reasons, as well as to support safe and reliable robot navigation, the sampling regions are chosen to be convex, elliptical more precisely (in 2D). Note that an ellipse can effectively stretch along any desired orientation and width and is compactly represented by a vector of only 5 parameters, $\mathbf{c} = (c_x, c_y, a_l, a_w, \theta)$, defining respectively the coordinate of the center, lengths of major and minor axes, and rotation angle in a given reference frame. In the following, describing our strategy for data sharing and assimilation, we refer to sampled regions as clusters (of sampled locations), to convey the fact that we are not necessarily bounded to use elliptical regions: our strategy can be applied to any planning scheme producing a 'cluster' of points, as long as it can be compactly described by a vector of parameters such as c defining a convenient embedding region.

In the process described above, robots can adopt different schemes of *local communication* in the *ad hoc robot network*, ranging from no communication, to full, sustained communication (an ideal configuration). This implements a full spectrum of coordination for maintaining a shared data map and letting the individual robots planning being mutually aware of others' plans and status of the map learning process. Hereafter, the two extreme cases are referred to as *No Data Sharing (NDS)* and *Full Data Sharing (FDS)*.

While data sharing might be essential to achieve a coordinate and optimized planning in the distributed MRS, this comes with a cost, both in terms of communications and computations, that could negatively impact scalability of performance. We therefore seek for a good balance, that can let the robots share only the truly useful data, to ensure scalability while only minimally affecting final map accuracy.

4 DATA SHARING AND ASSIMILATION

Our approach for data sharing and assimilating is hinged to the *submodularity* property of information gathering processes (Krause and Guestrin, 2011), which is addressed by tracking the expected changes in the GP model with the addition or deletion of measurement points. The intuitive understanding of submodularity is that if we obtain more measurements in the region where there is already significant data recorded, any additional measurements will only marginally increase our information of the location and will not 'influence' our environment model significantly. Leveraging submodularity, we can filter the data inside the region (cluster) by forming a sum*mary*. The objective of the summary is to subsample the data already sampled in the region, by selecting a subset of data which is much smaller than the original set but informative enough to minimize information loss with respect to the GP. This subset makes a cluster summary which can used for data sharing. Note that, by exploiting the the fact that a cluster's embedding region is represented by a vector of parameters such as c for the case of ellipses, both the data summary and coordinates and geometric description of the cluster (which we call its *meta-data*), are expressed in compact forms suitable to be effectively shared in the ad hoc network.

Moreover, in our strategy, we account for robots' time budget for traveling, such that a robot, *on-demand*, only requests and assimilates (i.e., data summaries) that are expected to be used and that can carry a relevant information gain for its GP. E.g., data summaries of regions that are far away and beyond the current traveling budget will not be considered.

4.1 Summaries of Cluster Data

As mentioned before, we consider the scenario where robots sample in 2D elliptical regions that define a data cluster. Based on the described iterative adaptive planning process, at a time t, a robot can have sampled in $m \ge 1$ different cluster regions, with the measured data for locations and values being assimilated in robot's GP and stored in internal data structures, where the k-th cluster is represented by its parameters $c^k = (c_x, c_y, a_l, a_w, \theta)$, which make the metadata of the cluster. At the same time, in the internal data structures, these parameters uniquely identify the set of data points inside the cluster: $c^k \rightarrow \mathbf{D}^k = [\mathbf{x}_1, y_1], ..., [\mathbf{x}_i, y_i]$, where the location of a specific data point is given by a position vector \mathbf{x} and its value y is the measured attribute of interest to map.

In a GP, a data point contributes to its 'local' outcome and extends its influence to its local neighbourhood based upon optimized GP's kernel hyperparameters. Based on this, here we define the notion of *influence points*: subset *s* of points of the full data set in the cluster, $s \subset \mathbf{D}$, whose presence inside

GP's structure affects GP's prediction outcomes the most. Influence points form the *cluster summary*. A data point that is not an influence point can be removed, such that other data points assimilated inside the GP could accurately predict the outcome value of the removed data point by predicting its corresponding mean value conditioned on the summary.

We can assess the quality of a summary set *s* by measuring the prediction error that it induces on data points x_k that are not part of the summary set compared to the prediction that the full GP would give:

$$error(x_k|s) = \mathcal{GP}(\mu_{x_k}|s) - y_k \tag{1}$$

Here the error is computed as the difference between GP's prediction outcome $\mathcal{GP}(\mu_{x_k}|s)$ for point x_k conditioned on the summary set *s*, and the value y_k , which is the prediction from the the GP including all the data from the cluster, not only the the summary set. In the notation $\mathcal{GP}(\mu_{x_k}|s)$, μ_{x_k} indicates the mean value of GP's prediction at the location of that data point x_k .

Computing the prediction error in this way, we are making the *naive* assumption that to predict the output value at certain location inside the cluster we don't need the GP being conditioned on data outside the cluster, hence, the summary is localized to a single cluster. This naive assumption forms the basis for our *summarization* process and the result of which will be tested in our computation experiments.

For each sampled cluster c^j comprising of a set \mathbf{D}^j of n_j data points, in order to define the *best* subset of influence points (i.e., the *summary*), a robot needs to select the subset $s^j \subset \mathbf{D}^j$ that can provide a minimal expected error over all points in the cluster. At this aim, we define a *combinatorial optimization problem* to find the subset $s^j \subset \mathbf{D}^j$ such that if the GP is purely conditioned on s^j , its cumulative prediction error for all the data points in the cluster is minimized (indeed, only the points in the set $\mathbf{D}^j \setminus s^j$ give a contribution):

$$\min_{s^{j} \in \mathcal{P}(\mathbf{D}^{j})} \sum_{k=1}^{n_{j}} \left\| y_{k} - \mathcal{GP}(\mu_{x_{k}} | s^{j}) \right\| w_{j}^{2} \qquad (2)$$

where $\mathcal{P}(\mathbf{D}^{j})$ is the power set of the data points \mathbf{D}^{j} inside the cluster c^{j} and w_{j} is equal to the cardinality of the subset s^{j} . The role of w_{j} is to avoid that the whole set \mathbf{D}^{j} is selected, and, more in general, balancing prediction accuracy and size of the summary set, acting as a penalty for large sets. It should be noticed that the minimization problem (2) is not limited to regression algorithms such as GP, but it can be utilized to form summaries using any regression algorithm.

The optimization problem is NP-hard, hence we tackle it with a custom *heuristic* to ease its solution in online scenarios. We form adjacency matrix using the euclidean distance between the data points. For each

data point in the adjacency matrix we compute its average distance to all the other data points as well as determine its closest neighbour. We then compare its closest neighbour to check if both the data points are a mutually closest pair, i.e no other points are closest to either of the pair. If this condition is satisfied, we select the point from the pair that has greater average distance to all the other data points. The rationale is that spatially distant points are more informative compared to neighbouring data points, due to their relatively high uncertainty. The process is repeated until all pairs are accounted for and an initial summary is formed by the points that are selected.

We then form a GP based on this summary and test if the prediction of the GP is accurate enough by randomly comparing GP's prediction on the un-selected data points output value. More specifically, the k-th data point is added to the summary and assimilated into the GP if the following condition on the relative prediction error is satisfied:

$$\frac{|\mathcal{GP}(\mu_{x_k}|s_{k-1}^j) - y_k|}{|y_k|} > \varepsilon \tag{3}$$

where s_{k-1}^{J} is the current summary set (i.e., *before* checking the *k*-th point). ε is a strategically-defined error bound, that was set to 0.3 in the experiments. The choice has been motivated by the observation that the aim of summarizing is to balance accuracy vs. size, with the goal of preparing a summary set for sharing that can make a receiving robot sufficiently informed for efficient future planning. Thus, an error bound of 30% seemed sufficient for the purpose.

The process of adding data points to the summary set is terminated when the moving average of the last *m* consecutive data points goes below the error bound ε . In the experiment *m* was heuristically chosen equal to 5. The points that the new GP had assimilated after the termination form the summary of the given cluster.

4.2 Sharing Summary Data

Robots do not locally broadcast their summaries directly, but broadcast first their way more compact *meta-data*. This is to let the nearby robots know about the regions that are available for summaries. A robot R hearing a meta-data message of another robot Schecks if the listed regions are already in its memory. If this is the case, R doesn't need the summary data from S, preventing from obtaining duplicate data.

If the meta-data message indicated the presence of new regions, the data of which is not in R's memory, the summary associated with those regions could be useful for R's future planning. However, before requesting the summaries, R uses the meta-data of each cluster c^k available from *S* to verify that the cluster is reachable in the future, based on its currently available discounted travel budget. We discount the robots current travel budget to account for the fact that the robot may travel to that region sometime in future. In practice, if a cluster c^k is too far, it would be of no use to get its summary, since *R* will never go to or traverse that region during the rest of the mission. Therefore, only the clusters that are reachable and correspond to unexplored regions are explicitly requested from *S*. Once received, summaries and associated meta-data are stored in memory, but are not assimilated yet in *R*'s GP model, which is explained in the next section.

4.3 Data Assimilation

The assimilation of received data summary into the local GP is a two fold process. First, the robot determines the next location / region where to sample based on the inputs given by its local planner. Let c^{next} be the cluster region to visit. Its parameters are compared with the meta-data of all the clusters stored in the robot's memory, that have been received from other robots. If the distance of any cluster c^k in memory falls in some close proximity of c^{next} , then the robot *assimilates the summary s*^k associated to cluster s^k into its current GP, thereby augmenting its knowledge without actually visiting the locations.

The process then repeats, where the robot identifies the new cluster to visit with the augmented knowledge and compares the parameters of the new region from the planner, with all the other meta-data of clusters stored in its memory. For the sake of optimization, a robot assigns flags to clusters in its memory whose summary has already been assimilated, and removes them from future computation for assimilation.

This sequential process ensures that size of GP is kept minimum for fast computation and sufficiently rich for informed planning and survey. Note that the described process is general, and not really bounded to any specific planning strategy.

5 COMPUTATIONAL EXPERIMENTS

We have empirically validated the proposed approach for controlled data sharing and assimilation in a series of *simulation experiments* conducted in a custom environment developed in Python. In particular, we have addressed *accuracy* in learning the ground-truth and *scalability* of performance and computation with increasing number of robots in the team. We have compared the proposed strategy with the two 'extreme cases' NDS and FDS mentioned in Section 3 where, respectively, no data sharing and full data sharing happen among the robots. The NDS and FDS cases represent upper and lower bounds in terms of performance for efficient sampling. In the NDS case each robot is on its own such that the iterative planning for the next region where to sample data is only based on local G data. In the FDS case, at each instant all robots have global knowledge of mission status (i.e., of the estimated GP map) such that individual planning decisions get implicitly coordinated.

In the following, we refer to the data sharing and assimilation approach proposed in this paper as *SDS*, standing for *Selective Data Sharing*.

We are motivated by surveying marine environments, so the test scenarios utilize actual bathymetry maps from the GEBCO database (www.gebco.net). Robot's mission is to collectively learn these maps by forming a global GP at the end of the experiment (i.e., at the end of the mission, all individual GP maps are merged to create one single, suitably aggregated map). To ensure that computational experiments are close to realistic physical experiments, we have considered a time budget of $T_{max} = 4h$ for the mission, with the robots moving at a constant velocity v = 1 m/s (modeled after the LutraProp robot bots, senseplatypus.com/lutra-prop), and a large area to map. To equalize the experiments, we scaled all bathymetry data to match $6000 \times 6000 \text{ m}^2$ area and discretized the data into a grid of 10000×10000 possible sampling points. Robots' depth sensor returns the ground-truth value corrupted by a Gaussian noise equal to 5% of the expected value. Communication range is set to 90m, after typical ranges in ad hoc wi-fi networks.² To initialize the robot's GP, a prior sparse random sampling of the area is conducted on the given dataset, with a total of 10 data points, (e.g., similar to (Li et al., 2020; Mishra et al., 2018)).

5.1 Baseline Algorithm, Metrics for Accuracy and Scalability

As we have pointed out, our goal with the SDS strategy is to strike an effective and efficient balance between accuracy in map reconstruction and overall scalability. Therefore, we need first to have a baseline for assessing goodness in map reconstruction accuracy. At this aim, we compare the accuracy performance of the three considered approaches, SDS, FDS, and NDS, against a *random-waypoint* (*RWP*) (Bettstetter et al., 2003) model for robot mobility. In RWP a robot iteratively identify the next location where to sample according to a random selection of direction and distance for traveling (in the experiments we have selected angle ranges that make a robot to have a tendency to move forward, and distances that are relatively long to attenuate the submodularity effect). In practice, RWP is a *random and uninformed strategy for planning*. The use of random coverage approaches as baseline is common practice in the literature (Popović et al., 2020; Li et al., 2020).

In the experiments, we use the *Root Mean Square Error* (RMSE) of GP's predictions vs. bathymetry ground truth computed over all the sampling points in the map. Based on the RMSE, we report the percentage of improvement of each approach vs. RWP according to the following formula:

$$\frac{\text{RMSE}(S, \text{ Ground Truth})}{\text{RMSE}(RWP, \text{ Ground Truth})} \times 100$$
(4)

where RMSE(*S*, Ground Truth) is the root mean square error between the predictions given by the final GP constructed by algorithm *S* and the ground truth values. *S* is a substitute for SDS, FDS, NDS.

5.2 Performance in Mapping Accuracy

The results from the experiments addressing map accuracy are reported in Table 1 for a team size of = 5robots. We have considered seven different marine rectangular areas from different parts of the world, latitude and longitude coordinates of the corners are shown. Each area has a different bathymetric profile, which can be more or less rugged. E.g., in a profile featuring smooth and highly correlated changes, a fully informed approach such as FDS is expected to strongly outperform less informed approaches. Instead, if the profile is rugged, showing rapid and sudden changes, the advantage of FDS might be comparatively less. Also in a mostly flat profile with a few abrupt changes the use of an informed strategy might not necessarily bring an advantage even over RWP, since it's mostly a matter of luck to hit the right spot.

Results show that FDS outperforms NDS as well as SDS in all the instances, as expected. However, the relative advantage differs from area to area precisely based on the discussion above. Unfortunately, as it will be shown in the next section, there's a price to pay for keeping all robots fully informed: FDS isn't really scalable. Moreover, in remote locations where there is not a global communication infrastructure, it might simply be unfeasible to assume global communications. The performance of SDS is, rightfully, between that of FDS and NDS. This shows that the controlled data sharing and assimilation does well its job by providing a mapping accuracy which is closer to FSD than NDS indeed, but with a significant saving in

²Experiments were conducted in Intel(R) Xeon(R) CPUX5690 @ 3.47GHz hardware with 16GB RAM.

Table 1: Mapping accuracy of the different approaches on various bathymetry datasets. Reported percentages for FDS, NDS, SDS are improvements over RWP according to (4).

Dataset	lon0	lon1	lat0	lat1	FDS%	NDS%	SDS%
Oman Coast	58.389	23.66	58.914	24.197	41.56	2.34	28.93
North Pacific Ocean A	168.347	172.856	39.831	43.647	60.49	14.60	39.77
North Pacific Ocean B	145.814	152.731	31.177	35.993	9.67	5.83	12.82
North Pacific Ocean C	165.352	170.806	22.956	26.514	35.43	24.02	29.76
South Pacific Ocean	168.03	173.65	-28.15	-22.63	9.74	14.31	8.38
Philippine Sea	129.681	134.743	17.289	22.218	8.88	1.41	8.67
Genoa	8.306	9.314	43.085	44.022	69.73	64.28	50.32
Average Standard Deviation					36.26 20.93	18.99 21.11	26.81 15.42

Parameters: n = 5, Time budget = 4hr, Max replanning period = 1h, v = 1 m/s, $\varepsilon = 0.3$.

computation and communications, as it will be shown in the next section. Moreover, being inherently based on local, range-limited communications, SDS is always a feasible approach in real scenarios.

Note that all the approaches systematically perform better than the RWP baseline. In SDS, data shared in opportunistic communications enables little less than 30% improvement in performance compared to no communication (RWP and NDS).

Interestingly, f the cases of Genoa, South Pacific Ocean and Philippines Sea datasets, SDS performed nearly as good as optimistic FDS approach, without the burden of computation that the FDS approach suffers from. An interesting observation from the tested data sets is that the standard deviation of all three approaches are more or less the same and it is attributed to exactly the same experiment environment and survey strategy that is provided to test all three approaches. The standard deviation in the reported improvements is the result of how survey strategy performs at particular dataset. If the dataset is more or less flat, then informed strategies generally perform as good as uninformed strategies and hence, the improvement on such datasets is usually lower.

5.3 Scalability of Computations

In order to perform a *scalability analysis* we study how computations for local GP updating and communications for data sharing grow with team sizes for the different algorithms. In particular, we measure the *average data size which is assimilated into robot's GPs at the end of the experiment*. The size *M* of a GP (in terms of number of used data points) directly impacts on the time for each GP updating, which goes as $O(M^3)$. The more the robots engage in data sharing, the more the GP inside robots grows, and the more computation time is required for individual updates, and the more communication bandwidth is used. In the case of NDS, since there's no data sharing, GP's data size is fairly independent of team size, it thus serves as a lower bound. Instead, FDS serves as upper bound, given that robots engage in global communications. We will therefore refer the measures to the FDS baseline, measuring the improvement (i.e., the reduction in size) with respect to FDS.

We consider teams of size 5, 10 and 15 robots, and for each algorithm we measure: average GP data size at the end of the experiment; standard deviation of the data size; % of decrease of data size with respect the data size in FDS; accuracy loss with respect FDS. Results are reported in Table 2 and refer to the scenario North Pacific Ocean A of Table 1.

It can be noted from the results that the average size of the GP data assimilated by robots in FDS grows significantly compared to the average sizes in SDS and NDS, which shows the poor scalability of the FDS strategy for large team sizes. The average data size for NDS remains more or less the same, as expected, while the average data size in SDS seems to increase only marginally compared to FDS. The relatively large standard deviation of the data size in SDS is the result to the stochastic nature of whole planning, sampling, data sharing and assimilation process. On the account for accuracy, we observe that even though both SDS and NDS suffer from some accuracy losses, SDS' loss in accuracy in SDS is minimal compared to FDS. Also we can notice a general decreasing trend in the magnitude of loss with the increase of team size, both for SDS and NDS. It is plausible that as the region gets saturated with more robots for mapping, the effects of full data sharing and utilization for efficient planning of sampling actions tends to diminish. Overall, these results well support the scalability of the strategies with reduced or even no data sharing.

Moreover, since the computation time for obtaining a new plan is proportional to the size of the GP

# Robots	Strategy	Average GP size	Standar dev.	% Decrease of avg. GP size	% Decrease in accuracy
5	FDS	318.5	-	-	-
5	NDS	67.4	7.1	79.8	40.0
5	SDS	197.2	10.2	38.1	14.8
10	FDS	870	-	-	-
10	NDS	90.5	13.9	89.5	26.5
10	SDS	259.8	39.2	70.1	7.1
15	FDS	1021	-	-	-
15	NDS	90.7	33.4	91.1	10.2
15	SDS	242.1	336.2	76.2	5.4

Table 2: Scalability in GP data size and mapping accuracy for various team sizes and data sharing strategies. All percentage decreases are referred to the baseline FDS strategy. Accordingly, FDS' percentages are reported.

data³, we also analyze how *planning computation time* is affected by GP's data size. This is a central metric to assess the capability of the robot team to *effectively operate in a real-world scenario*. In fact, if even a single planning operation would consume a relevant fraction of the available time budget, the robot team will not really be used to perform extensive data sampling, as needed. In other words, the previous scalability of GP size needs to be accompanied by equivalently good scalability results for computing a plan in order to truly assess strategy scalability.



Planning iteration

Figure 1: Computation time for planning over the iterations.

In our model, a new plan is issued at more or less regular time intervals. Therefore, we study how *computation time varies across planning iterations*. Note that since GP's size grows over time as a result of sampling actions and data sharing, the planning computation time for a robot is expected to grow accordingly. Results are shown in Figure 1, which refers to a 10-robot team for North Pacific A scenario of Table 1. Note that the *y*-axis in in logarithmic scale. We can observe that the computation times for FDS are by far the highest and raise up rapidly, showing the non-scalability of the strategy. SDS and NDS have similar in performance to in the first three planning iterations. This is due to the fact that robots only share the data if they are in close proximity and only assimilate the data if needed in upcoming planning iteration, so initially SDS behaves similar to NDS as far as plan computation time is concerned. After the third iteration, SDS shows an increase in computation times, but still stays well below FDS. The same trend is shown for larger teams (not shown). However in these cases, SDS' times become of order of minutes, such that a strategic adjustment of parameter ε of Eq. (3) would be required for efficient online execution.

From the results reported in Figure 1 and Table 1, we can conclude that in the proposed SDS, the utilization of data summaries, sharing them with other robots on demand, and assimilating them at the right time enable significant a retention of the accuracy compared to the ideal FDS strategy, while keeping computational load relatively low and therefore supporting the *good scalability of the SDS approach*.

6 CONCLUSIONS

We have considered scenarios where a mobile and distributed multi-robot system is used for mapping a spatial field in a given limited time budget. Building on our previous work (Di Caro and Yousaf, 2021b; Di Caro and Yousaf, 2021a) we have focused on the case where mapping is modeled as a regression task and tackled using a GP model. Robots collectively learn the GP by taking measures at discrete locations. In turn, robots use the GP to iteratively and adaptively compute plans aimed at identifying the most informative regions where to go and take measures. While the use of multiple robots can bring clear benefits, at the same time a number of challenges arise regarding coordination and joint action planning. Central to these challenges is the use of communications to share data and therefore maintain a shared view of the GP map to let the robots coordinate their planning actions.

In this paper we have precisely addressed the is-

³As described in (Di Caro and Yousaf, 2021b), for computing a plan (i.e., a region and a path where to sample) a robot makes use of a utility function based on a measure of expected information gain for GP learning.

sue of selectively filtering and compacting the data that each robot could share, and define efficient strategies to let nearby robots only exchange data that are expected to be informative for a robot conditioned to its local GP and available time budget. In addition, we have defined a strategy for the selective assimilation of the data in a robot's GP only if and when truly needed. All these strategies are aimed at endowing a multi-robot system with data sharing and assimilation mechanisms that would guarantee effective mapping accuracy while ensuring overall scalability in computations and communications.

More specifically, we have described an orchestrated approach (termed SDS in the text) where robots form summaries of the regions they sample, organize the summaries into clusters with meta-data, share their summaries with nearby robots on-demand, assimilate the summaries into their local GP at the right time. We tested the approach in simulation using actual bathymetry datasets and showed how the proposed summarization, sharing, and assimilation strategy compares against strategies where either all data is shared and assimilated by all the robots or no data is shared among the robots for planning. Results have shown that the proposed strategy can balance very well accuracy and scalability, automatically providing highly accurate maps without incurring in excessive computational and communication load, overall showing a very promising scalability.

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