Resource-aware Decentralization of a UKF-based Cooperative Localization for Networked Mobile Robots

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Abstract—Estimation techniques such as Unscented Kalman filter (UKF) are deployed for accurate joint location estimation in cooperative localization of cyber physical systems (CPS), e.g., to locate each robot in a cooperative mobile robotic network in GPS-denied environments. In order to avoid single point of failure in the centralized implementation of such estimation techniques, the decentralization of estimation algorithms has attracted considerable attention in the past two decades. However, the design of decentralized algorithms with reduced communication cost without loss in accuracy for compute-intensive estimation techniques such as UKF has been challenging. In the decentralized UKF, the tasks are partitioned and computed locally at robot nodes. Data communication overhead is overwhelming due to tight data dependency between the robots’ computations. In this paper, we present a CPS framework for UKF decentralization in which computation and communication are tightly intertwined and computation replication is deployed in order to reduce the communication overhead among cooperative mobile robots. We demonstrate and evaluate the performance of our proposed approach work in a wireless network of 15 Raspberry Pi 3 B, with quad-core 1.2GHz 64bit CPU, emulating a network of mobile robots with onboard computation and communication capabilities. Our experimental results show that the End-to-End execution time of decentralized UKF prediction and update steps with replication are faster by up to 12.29 and 3.57 times, respectively, compared to the partially decentralized UKF algorithm of [1].

I. INTRODUCTION

Cooperative cyber physical systems are driven by the tight coordination between computational components, physical sensors and the interaction with each other. Software development for such systems, because of the tight integration, heterogeneity of resources, as well as safety and application timing requirements can be very complex. Networked mobile robot operations are among the challenging cooperative CPS applications. An important component of the networked mobile robot operations is the localization of the robots. The fast and accurate localization is important because a delayed estimation will mislead navigation and other applications and may lead to a mission failure. Cooperative localization is a reliable scheme for localizing a team of communicating robots in GPS-denied environments (Figure 1). This localization technique uses relative measurements among the robots as a feedback signal to jointly estimate the location of robots.

Cooperative Localization (CL) algorithms deploy various estimation strategies such as Extended Kalman filters (EKF) [2], UKF [1], maximum likelihood [3], maximum a posteriori (MAP) [4], and particle filters [5], [6], [7], [8]. Among these techniques, EKF-based cooperative localization algorithms, due to their recursive nature and relative ease of implementation, have been studied extensively. However, the EKF, due to linearization approximation, is known to be inconsistent for highly nonlinear systems or when the filter has high initialization errors. UKF is an alternative recursive estimation filter, which is proven to work more consistently than the EKF for systems with nonlinear state and measurement models (c.f. [9]). For large systems, however, UKF is computationally more expensive than EKF. This paper focuses on UKF-based cooperative localization.

To increase the scalability and avoid a single point of failure, decentralization of localization has been promoted by robotics community. However, due to relative measurement updates, the local estimates of the cooperative robots are highly correlated. This creates a great challenge in the decentralization of cooperative localization algorithms. Decentralized cooperative localization algorithms normally come with a significant processing and communication requirements (see e.g., [10], [11], [12], [4], [13], [14]). Ignoring estimation correlations can lead to filter inconsistency and even divergence [15], [16].

Various decentralization schemes using EKF formulations have been proposed in the literature (see e.g., [10], [13], [14], [17]). In UKF, the estimation equations of the robots in the team are highly correlated. Thus, naive decentralization (or partitioning) of UKF algorithm can result in a large amount of data transfer among the robots. In [1], a partially decentralized UKF is proposed, in which the UKF equations are decoupled in a way that their computation is distributed among all the cooperative robots. However, a shared memory on a server is used to transfer data among robots, leading to
transfer of a large amount of data between the server and the robots. The communication time is usually much higher than the computation time, which makes this method inefficient.

The literature on the decentralized cooperative localization algorithms that is reviewed above is on the algorithmic decoupling of computations without loss in performance and accuracy (or with acceptable performance loss). The proposed methods lack rich knowledge on characteristics of the underlying network and computational resources. Wireless communication overhead, heterogeneity in embedded computation, and sensing features of mobile robots may incur significant delay and long execution time during localization. Therefore, decentralization is challenged by the tight coupling and integration of sensing, computation, and communication among the mobile agents. Without careful trade-off between computation and communication, a decentralized algorithm may suffer from unacceptably long End-to-End delay and may not be deployable for agents with high mobility. This mandates us to develop a CPS framework to overcome the complexity of decentralization of cooperative localization based on complex estimation techniques such as UKF without loss of of accuracy and computationally-identical to the centralized method.

We treat decentralization of UKF as a problem of $k$-way task partitioning problem. Each partition is a subgraph in UKF task graph that is computed in each robot. Unlike [1], the proposed method does not need a server for computation or sharing data. Applying the partitioning method on the target application, we observed that the majority of delay on critical path comes from communication delay. In order to reduce the data transfer further, we propose a communication graph to minimize the number of data links followed by computation replication based on Min-cut Max-flow Theorem. Replication has been researched extensively in VLSI CAD partitioning and design automation community (e.g., [18], [19]). The replication comes with the cost of an increase in local computation. In our target application, measurements show that the wireless communication delay is significantly higher than computational delay and hence, the computational delay overhead is negligible. The proposed min-cut replication results in significant reduction of UKF data transfer among the mobile nodes. The experimental results show that the End-to-End execution time of decentralized UKF prediction and UKF update steps are faster by up to 12.29 times and 3.57 times, respectively, compared to partially decentralized method [1].

II. BACKGROUND

A. Cooperative Localization

Localization is one of the basic applications in mobile robots that provides other applications such as navigation with location information. The fast and accurate localization is important because a delayed estimation will mislead navigation and other applications and might lead to mission failure. The time interval between execution of localization ($T$) should be small enough to be able to capture the motion of the robot and provide the higher level applications with accurate and almost real time location. In the GPS-denied environment, because of accumulated proprioceptive sensors error, the estimated location drifts from the actual location after a while.

Cooperative Localization (CL) is a scheme in which robots improve their localization accuracy by jointly processing inter-robot distance measurements obtained by exteroceptive sensors such as Kinect. In CL, a single measurement between two robots can improve the localization accuracy of other robots as well. For CL using Kalman filter and its variants such as EKF and UKF, as it is shown in Figure 1, robots share their control signals ($U$), state vectors ($X$), their uncertainty ($P$), and sensors measurements to obtain more accurate location ($Pred_X$) and uncertainty of estimated location ($Pred_P$) for further processing by higher level applications such as navigation. In the context of CL, the state vector ($X$) includes the global pose (position and orientation of the robots) as well as possibly other states potentially needed to model the dynamics of the robots (for example, steering angle and actuation). In a networked system of $N$ robots, state vector of each robot $i$ is represented by $n_i$ state variables, which adds up to $n$ states of the system ($X$).

B. UKF

UKF is a recursive filter for estimating the state of a system referred to as $X$. UKF is composed of prediction and update steps. Prediction step runs periodically, and update step runs whenever there is a measurement in the system. In the prediction step, the state estimation of the system will be predicted based on the control signals, self motion measurements, the previous state of the system, and covariance matrix. Update step runs whenever the system receives a measurement.

1) Prediction Step: The prediction step for the collective system with $n$ states starts with computing the square root matrix, as a triangular matrix, of matrix $P$ using Cholesky Decomposition (CD) method. After that, a set of $2n + 1$ sample points, called Sigma Points (SP), is generated by eq. 1b. In the equations, $c_{i}^{th}$ denotes the $i^{th}$ column of the matrix. The system model function will use $SPs$ and $U$ to
generate Transformed Sigma Points (eq. 1c). The predicted state is the weighted arithmetic mean of $T_{SP}$ (eq.1d). The prediction covariance matrix ($Pred_P$) will be obtained by eq. 1f using prediction error $E$ (eq.1e). In the equations, $l \in \{0, \cdots, 2n\}$, $c \in \{1, \cdots, n\}$, and $w$ are system defined constant.

$$CD = \text{CholeskyDecomposition}(P)$$ \hspace{1cm} (1a)  

$$SP_{(0)} = x, SP_{(c,c+n)} = x \pm (\sqrt{(n+c)}CD_{(c)})$$ \hspace{1cm} (1b)  

$$T_{SP(l)} = \text{SystemModel}(SP_{(l)}, U)$$ \hspace{1cm} (1c)  

$$Pred_X = \sum_{i=0}^{2n} w(i) T_{SP}(i),$$ \hspace{1cm} (1d)  

$$E_{(i)} = T_{SP}(i) - Pred_X$$ \hspace{1cm} (1e)  

$$Pred_P = \sum_{i=0}^{2n} w(i) E_{(i)}E_{(i)}^\top$$ \hspace{1cm} (1f)  

2) **Update Step:** Update step runs whenever the system receives a measurement ($Meas$). For example, when robot $A$ measures the relative distance from robot $B$, UKF update function fuses the measurement feedback with the predicted state, computed by prediction step, using Kalman gain. In update step, at first, innovation covariance ($S$), cross-covariance ($P_{xz}$), $E_{z}$, and innovation ($r$) will be generated by eq. 2 using the SPs.

$$T_{SP\_meas} = \text{MeasurementModel}(SP)$$ \hspace{1cm} (2a)  

$$Pred\_Meas = \sum_{i=0}^{2n} w(i) T_{SP\_meas}$$ \hspace{1cm} (2b)  

$$r = Meas - Pred\_Meas$$ \hspace{1cm} (2c)  

$$E_{z}(i) = T_{SP\_meas}(i) - Pred\_Meas$$ \hspace{1cm} (2d)  

$$S = \sum_{i=0}^{2n} w(i) E_{z}(i)E_{z}(i)^\top$$ \hspace{1cm} (2e)  

After that, the Kalman gain ($K$) will be computed using $S$ and $P_{xz}$ by eq. 3.

$$P_{xz} = \sum_{i=0}^{2n} w(i) E_{(i)}E_{z}(i)^\top$$ \hspace{1cm} (3a)  

$$K = P_{xz}S^{-1}$$ \hspace{1cm} (3b)  

Finally, the $Update\_X$ and $Update\_P$ will be computed by eq. 4.

$$Update\_X = Pred\_X + Kr$$ \hspace{1cm} (4a)  

$$Update\_P = Pred\_P - KS K^\top$$ \hspace{1cm} (4b)  

$Pred\_X$ and $Pred\_P$ (the output of prediction step), or $Update\_X$ and $Update\_P$ (the output of update step) will be used as $X$ and $P$ for the next iteration of UKF.

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C. **UKF in Cooperative Localization**

In this paper, we denote the components of the aggregated state vector $X$ of the team corresponding to robot $i$ by $X[i]$. In a similar way, the corresponding portion of $P$ and $E$ matrices for robot $i$ is denoted by $P[i]$ and $E[i]$. Note that each of these sub matrices has $n_i$ rows, which is the number of states of robot $i$. Figure 2 shows the shape and size of the main matrices$^1$.

Figure 3 shows the task graphs for UKF prediction and update steps, tagged with the equation numbers, where each robot $i$ sends its control signals ($U[i]$) to the UKF to compute the $T_{SP}$s using eq. 1c. At the end, the $Pred\_X$ and $Pred\_P$ might be used as the $X$ and $P$ of the next iteration of UKF. The navigation and other applications of robot $i$ will use $Pred\_X[i]$ and $Pred\_P[i]$. In the update step, the measured relative distance will be used to compute eq. 2. The outputs, which are $Update\_X$ and $Update\_P$, will be used as the $X$ and $P$ of the next iteration of UKF, and also will be sent to robots. The UKF update step runs whenever there is a relative distance measurement in the system.

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III. **PROPOSED FRAMEWORK FOR UKF DECENTRALIZATION**

To increase the scalability and avoid a single point of failure, it is necessary to decouple UKF equations in a way that the UKF computation is distributed and each robot $i$ computes $P[i]$ and $X[i]$, which are its own location parameters. Due to sequential data dependency between the robots to compute $CD$ and $Pred\_P$ (see Figure 4 and [20]), all the robots are actively involved in computing UKF and high amount of data transfer through wireless communication would increase the execution time of UKF. According to [1], all the computations of prediction step of robots are independent, except for the computation of $CD$ and $P$ (eq. 1a and 1f).

To fully decentralize the UKF, at design time, we partition the task graph into a set of subgraphs. Each subgraph is computed locally at the designated robot. In order to provide an effective decentralization, we need to perform optimizations to reduce the data transfer among the robots.

$^1$If $Pred\_X$ and $Update\_X$ are as the same size of $X$. Similarly, $|Pred\_P| = |Update\_X| = |P|$, and $|SP| = |T_{SP}| = |E|$.
In this paper, our objective is to obtain a fully decentralized UKF with reduced data links and transferred data size, while keeping the computation of UKF decentralized. In addition, we propose a method that further reduces the data transfer between robots by replicating the computations. Note that the computation in our method is identical to the computation in centralized UKF, so the accuracy of the filter will not be affected.

A. Row-based UKF Partitioning (R-UKF):

As mentioned in the previous section, each robot \( i \) needs \( X[i] \) and \( P[i] \) \(^2\). The control signals \((U[i])\) is generated locally. The UKF equations can be decoupled in a way that each robot just computes what it needs. For example, robot \( i \) just needs to compute \( SP[i] \). There is data dependency between computations of some sub matrices, hence each robot has to acquire them from other robots. Based on eq. 1 and [20], in the prediction step, all computations can be completely decoupled, except for \( CD \) and \( Pred_P \). \( CD[i] \) depends on \( CD[1:i-1] \) and \( Pred_P[i] \) depends on \( E[1:i-1] \), which has to be transferred from other robots.

\(^2\)Recall that we represent the corresponding \( n_i \) number of rows in \( X \), i.e. from row \( n_i \) to \( n_i+1 \), with \( X[i] \). To show the corresponding rows of \( X \) from \( i \)th robot to \( j \)th robot, we will use \( X[i:j] \). The same representation is used for other matrices as well.

Hence, each robot \( i \) after receiving \( CD[1:i-1] \) and \( E[1:i-1] \), computes the \( CD[i] \), \( SP[i] \), \( P[i] \), \( T_SP[i] \), \( Pred_X[i] \), \( Pred_P[i] \), and \( E[i] \). Figure 4 shows the proposed UKF prediction task graph. The task graph is divided into sub-graphs associated with each robot \( i \).

In our framework, this is referred to as Row-based UKF partitioning. This partitioning will reduce the size of data transfer among the robots. Note that the overall computation and result will be identical to the centralized UKF. Unlike [1], this partitioning does not require a server for computation or sharing data. The UKF prediction task graph shows that estimation of the location of robots is highly correlated and dependent on each other. As a result, the computation is sequential and parallelism is limited.

Figure 5 shows UKF update task graph. Each pair of robots, e.g. robot \( A \) and robot \( B \) can occasionally measure the relative distance between each other. Robot \( A \) that has the relative distance measurement will receive \( SP[B] \) from robot \( B \). Then, using \( SP[A] \), \( SP[B] \), and measurement data, robot \( A \) will compute \( S, r, \) and \( E_z \). After that, starting from robot 1, each robot \( j \) will compute \( K[j] \) using \( E[j] \), which is generated locally, and using \( S, r, \) and \( E_z \). In addition, Robot \( j \) receives \( K[1:j-1] \) from robot \( j-1 \) to compute \( Update_P[j] \) and \( Update_X[j] \). \( K[1:j] \) will be sent to robot \( j + 1 \) for computing the relative UKF update step. Note that R-UKF has exactly the same computation of a centralized UKF for CL.
B. Communication Graph Refinement

Based on eq. 1, submatrices such as $CD[i]$ and $E[i]$ that are computed in robot $i$ have to be transferred to all robots with label from $i + 1$ to $N$. We represent the $N$-way partitioned UKF task graph by $G = (V, L)$ where $V = V_1, V_2, \ldots, V_N$. $V_i$ is a set of UKF computation nodes in $V$ that belongs to robot $i$, and $L$ represents the data dependency between UKF computation nodes. Figure 6.a shows the data dependency between the nodes. Sending $CD[i]$ and $E[i]$ from $V_i$ directly to all other agents with greater label (from $i + 1$ to $N$) requires a complete communication graph, hence $N * (N - 1)/2$ communication links will be required which impose a large computation and communication overhead. Note that establishing a TCP/IP connection includes handshaking and several packet transmission, and broadcasting protocols such as UDP cannot be used because they are unreliable. Since each $V_i$ cannot finish the computation of $CD[i]$ before receiving $CD[1:i-1]$, it has to wait till $V_{i-1}$ finish the computation of $CD[i-1]$, and so on. Therefore the computation of $CD[i]$ is sequential, and its critical path starts from $V_1$ and ends in $V_i$. Considering the fact that $V_i$ has $CD[1:i-2]$ when it computes $CD[i-1]$, it is possible to send $CD[1:i-1]$ from $V_{i-1}$ to $V_i$. This way, communication links between $V_1$ to $V_{i-2}$ and $V_i$ are not needed (see Figure 6.b). Reducing communication links will reduce the communication overhead. Note that in the case of UKF, the computation delay is small compared to communication delay and the time overhead of establishing and maintaining a TCP/IP link. The number of communication links in the proposed communication model is $N - 1$ which is the minimum possible number of communication links for a connected network. Note that the data transfer between the nodes can be realized through ad-hoc or infrastructure network. This discussion is out of the scope of this paper.

The proper partitioning and communication graph reduce the communication cost, yet large matrices such as $CD$ and $E$ have to be transferred. Therefore, we propose to deploy a graph-based partitioning refinement technique called min-cut replication to further reduce the size of transferred data.

C. Communication Minimization by Computation Replication on R-UKF (RR-UKF):

The partitioning, stand alone, cannot reduce the data communication significantly. To further reduce the size of data transfer between partitions, we replicate some of the computations. Figure 7(a) shows data links between two nodes $V_1$ and $V_2$. The size of data that has to be transferred between $V_1$ and $V_2$ is 10. By replicating node $b$ in partition $V_2$ (Figure 7(c)), the size of data transfer will be reduce to 4 (2+2). To obtain UKF task graph with minimum data transfer, we propose a two phases method. First, Min-cut Max-flow replication is applied to prediction step of UKF, which generates a task graph with replicated nodes. Second, UKF update task graph is pruned given the changes in UKF prediction task graph after replication. Then Min-cut Max-flow replication is applied to the pruned UKF update task graph. The result is a fully decentralized UKF with minimum data transfer, and hence, minimum End-to-End delay.

**Replication technique:** In the proposed method, we present how to minimize the data transfer using min-cut replication technique on decentralized prediction and update steps of UKF for CL. We are given a $N$-way partition of $G = (V, L)$ represented by $V = V_1, V_2, \ldots, V_N$. $V_i$ is a set of nodes in $V$ that belongs to partition $i$, and $L$ represents the data dependency between nodes. The set of cut edges is the set of edges $C \subseteq L$ that connects the partitions. Lets consider two partitions $V_1$ and $V_2$ as shown in Figure 7. Lets assume that $I$ is the set of incoming edges to set $V_2$ from $V_1$. The objective is to find a subset of nodes in $V_1$ called replication set $R_1$ such that if it is replicated in $V_2$, the cutsize between $V_1$ and $V_2$ is minimized. After replication of $R_1$ in $V_2$, the cut edges between $R_1$ and $V_2$ are eliminated from the cut set and the input edges to $R_1$ are added to cutsize. $R_1$ is a minimum replication set respect to $V_2$ if the cutsize after replication is minimum among all possible $R_1$.

Figure 7 shows an example of $G$ (a), min-cut and replication set (b), and the replicated graph (c). To find the minimum replication set, we construct a network graph $G_f = (V_f, L_f)$, where $V_f$ includes nodes in $V_1$ that are reachable from $I$ (node a, b, and d) and the nodes in $V_2$...
that are adjacent to \( I \) (node \( i, h, \) and \( e \)). To apply min-cut max-flow theorem, two dummy nodes are added to \( G_f \) as source (node \( s \)) and sink (node \( t \)). The min-cut max-flow problem on this graph will separate the replication set from the rest of the network. The replication set \( R_1 \) is the subset of nodes in \( V_i \) starting from min-cut edges to the nodes adjacent to \( I \) (node \( b \)). After replication of \( R_1 \) in \( V_2 \), the cut edges between \( R_1 \) and \( V_2 \) (edge \( b \rightarrow h \), and \( b \rightarrow e \)) are eliminated from the cut set and the input edges to \( R_1 \) are added to cutsize (edge \( a \rightarrow b' \)). The replication in \( V_2 \) in respect to set \( V_1 \) does not affect the edges between \( V_i \) and other partitions. Hence, we can apply the min-cut replication between every two partitions in a \( N \)-way partitioned task graph independently.

To find the minimum weighted cut, min-cut max-flow replication considers weight or capacity for the edges. In \( N \)-way partition of UKF, the edges are weighted based on the data size and the wireless communication cost (e.g., number of hops). The pseudo code for min-cut replication is presented in algorithm 1. Next, we present replication algorithm for UKF prediction and UKF update step.

Algorithm 1: Min-cut Max-flow Replication

\[
\begin{align*}
1 & \text{input : } V_i, V_j, C_{ij} \\
2 & \text{output: } V^*_j \\
3 & \text{initialization: } V^*_j = V_j; \\
4 & \text{Construct Graph } G_f(V_i, V_j); \\
5 & |C| \leftarrow \text{min-cut max-flow on } G_f; \\
6 & \text{if } |C| < |C_{ij}| \text{ then} \\
7 & \quad |V^*_j = V_j \cup R_j| \\
8 & \text{end} \\
9 & \text{Return } V^*_j;
\end{align*}
\]

We present our algorithm based on aforementioned min-cut replication and show how the cut edge set between the \( N \)-way partitions of UKF is drastically reduced both in prediction and UKF update. The pseudo code of RR-UKF is presented in Algorithm 2. The input to the algorithm in sequential row-based \( N \)-way partition of UKF, known as R-UKF. First, Min-cut Max-flow replication is applied on each pair of partitions in decentralized UKF prediction (line [3-6]). Then UKF update with respect to measurement between \( V_a \) and \( V_b \) is considered. The refinement of decentralized UKF update is decomposed of two steps. The first step is referred to as Prune-cut-edges (line [7]). Due to replication in UKF prediction steps, some edges from cutsets between the partitions are moved inside the partitions. If those edges are deployed in the corresponding partition of update graph, the edges are local and hence, those edges from the cut set are removed. Visiting the cut edges in UKF update graph, this function checks if any of them is removed in \( V^*_p \) after replication. The second step is to apply Min-cut Max-flow replication between the \( N \)-way partitions of UKF update similar to UKF prediction (line [8-11]).

Algorithm 2: RR-UKF-Replication

\[
\begin{align*}
\text{input : } & \text{N-way partition of UKF prediction Task} \\
& \text{Graph } V_p = V_p1, V_p2, V_p3, \ldots, V_pk \text{ and} \\
& \text{Update(a,b) Task Graph } V_u(a, b) = V_u1, V_u2, \ldots, V_uk \\
\text{output: } & V^*_p \text{ and } V^*_u \\
1 & \text{//Initialization;} \\
2 & V^*_p = V_p, V^*_p = V_p, V^*_u = V_u \\
3 & \text{//====================== UKF prediction;} \\
4 & \text{for each } (V_p1, V_pj) \text{ where } C_{ij} \neq \emptyset \text{ do} \\
5 & \quad |V^*_p = \text{Min-cut Max-flow replication } (V_p1, V_pj) \\
6 & \text{end} \\
7 & V_u = \text{Prune-cut-edges } (V^*_p, V_u) \\
8 & \text{//====================== UKF update;} \\
9 & \text{for each } V_i \text{ and } V_l \text{ on } V_u \text{update do} \\
10 & \quad |V^*_u = \text{Min-cut Max-flow replication } (V_u1, V_uj) \\
11 & \text{end} \\
12 & \text{Return } (V^*_p \text{ and } V^*_u)
\end{align*}
\]

Figure 8 demonstrates the decentralized UKF prediction with min-cut replication. The minimum weight cut is shown in the figure. The result shows that by replicating \( SP[1], T_X[1], E[1], Pred_P[1], \ldots, SP[i-1], T_X[i-1], E[i-1], Pred_P[i-1], \) in \( V_i \), the cut size will be minimized. As a result, each partition \( v_i \), after receiving \( U[1 : i-1] \) from previous partition, will compute \( CD[1 : i] \). Instead of getting \( E[1 : i-1] \) and \( CD[1 : i-1] \) from previous partition, it will calculate them using \( U[1 : i-1] \) (new cut edge) and the \( Pred_P[1 : i] \) and \( Pred_X[1 : i] \) from previous step of UKF which are located in the same partition. The size of \( U[1 : i-1] \) is much smaller compared to \( E[1 : i-1] \) and \( CD[1 : i-1] \).

Figure 9 shows the partition between partitions. Partition \( V_A \) has the \( SP[B] \) and \( E[1 : A] \) as a result of replication in UKF prediction \( V^*_p \). Hence, \( S, r, \) and also \( K[1 : A] \) are generated (refer to eq. 2 and 3 ). Partition \( V_A \) will send \( K[1 : A-1] \) to partitions \( V_{ui} \) where \( i < A \). Partition \( A \) will send only relative distance measurement between robots \( A \) and \( B \) to partitions \( V_{ui} \) when \( i > A \). Since \( SP[A] \) and \( SP[B] \) already exist in \( V^*_p \), all other parameters are locally computed and hence those edges from cutset are removed.

IV. EXPERIMENTS

A. Experimental setup

To evaluate the proposed method, we implemented it in a network of single-board computers, called Raspberry Pi 3 B, with quad-core 1.2GHz Broadcom BCM2837 64bit CPU, 1 GB main memory, and a built-in WiFi module of 802.11 b/g/n. For these experiences, the CPU frequency
is fixed to 1.2GHz. Each board $i$ runs Linux kernel v4.9 and an application process that implements robot $i$’s task for decentralized UKF in C++ with an open-source linear algebra library EIGEN [21]. To create a realistic network environment, we set up an isolated WiFi network with one access point, Netgear N300 wireless router and $N$ boards. In the following, we report a median value for the End-to-End delay which is obtained from 500 rounds of UKF.

In our experimental studies, we considered robotic team scenarios with robots with 6 local states and with measurement and control signals size of 3. In our implementation, the main performance metric is the End-to-End delay from robot 1’s initiation to robot $N$’s completion of each execution cycle of UKF, consisting of the prediction and the update steps, which include both the computation and the communication delays on all the robots (or boards).

### B. Evaluation of UKF End-to-End Delay

To demonstrate the effectiveness of our decentralized UKF over the partially decentralized UKF [1], denoted by Partially Decentralized UKF, we compare them in Table I. Because of lack of space, in the tables, just the entries for the odd number of robots has shown. Recall that the Partially Decentralized UKF [1] assumes that there exists a server to store the whole data of UKF algorithm and that for each UKF cycle all robots takes part in the computation of UKF, like RR-UKF. The number of robots, $N$, is varied from 3 to 15. From the table, it is clear that on average, the End-to-End delay measured for RR-UKF is reduced by a factor of up to 12.29 for prediction step and by a factor of 3.57 for update step, compared to Partially Decentralized UKF [1].

To indicate the effect of reducing the size of transferred data using the replication technique in decentralized UKF, we compare our row-based decentralized UKF implementation without (R-UKF) and with computation replication (RR-UKF). Table II shows the total network traffic and the on board End-to-End delay measured for each UKF iteration in R-UKF and RR-UKF. Total data bytes transmitted by all robots at application level in RR-UKF have been drastically reduced up to 140 times for prediction step and 7.26 times for update step in comparison with R-UKF.

Computation replication eliminates transmission of $CD, E$ and $E_Z$ between robots and replaces with Measurement and $U$ in the transmission which are much smaller. Thus, it leads to significant reduction in the total packet frames transmitted by all robots at kernel level. The reduction ratio is up to 47.67 for prediction step and to 3.3 for update step.
The reason why this reduction ratio is smaller than that of the transmitted data bytes is that every packet frame, whose maximum size is 1500 bytes in our WiFi network, does not convey the same number of data bytes depending on the network I/O behavior of the application. In addition, the replicated computations increases the run time. As a result, the End-to-End delay from robot 1 to robot 70NANB17H192 is reduced from 92 ms (R-UKF) to 54 ms (RR-UKF) for prediction step when $N=15$. For update step, it is reduced from 92 ms (R-UKF) to 54 ms (RR-UKF). This significant reduction in the End-to-End delay allows each robot to react much faster in response to the resulting localization data of other robots.

V. CONCLUSIONS

Cooperative localization is a method for increasing the localization accuracy within a network of cooperative robots. Usually, probabilistic estimation algorithms such as EKF is used for processing the data shared by the group. UKF is a variant of Kalman filter, which is more accurate but has more computation, and decentralization of such filters with tight correlation among its tasks requires a large data transfer between agents. We showed that the size of the transferred data can be reduced by applying a replication technique. Our experimental results showed that the End-to-End execution time of the decentralized UKF prediction and update steps with replication are faster by up to 12.29 and 3.57 times compared to the partially decentralized UKF algorithm of [1]. In future, we plan to extend our work to include the delay and computational complexities of distance measurement sensors as well as to consider probabilistic estimation algorithms other than UKF.

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