Distributed collaborative 3D pose estimation of robots from heterogeneous relative measurements: an optimization on manifold approach Joseph Knuth* and Prabir Barooah

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SUMMARY

We propose a distributed algorithm for estimating the 3D pose (position and orientation) of multiple robots with respect to a common frame of reference when Global Positioning System is not available. This algorithm does not rely on the use of any maps, or the ability to recognize landmarks in the environment. Instead, we assume that noisy *relative measurements* between pairs of robots are intermittently available, which can be any one, or combination, of the following: relative pose, relative orientation, relative position, relative bearing, and relative distance. The additional information about each robot's pose provided by these measurements are used to improve over self-localization estimates. The proposed method is similar to a pose-graph optimization algorithm in spirit: pose estimates are obtained by solving an optimization problem in the underlying Riemannian manifold $(SO(3) \times \mathbb{R}^3)^{n(k)}$. The proposed algorithm is directly applicable to 3D pose estimation, can fuse heterogeneous measurement types, and can handle arbitrary time variation in the neighbor relationships among robots. Simulations show that the errors in the pose estimates obtained using this algorithm are significantly lower than what is achieved when robots estimate their pose without cooperation. Results from experiments with a pair of ground robots with vision-based sensors reinforce these findings. Further, simulations comparing the proposed algorithm with two state-of-the-art existing collaborative localization algorithms identify under what circumstances the proposed algorithm performs better than the existing methods. In addition, the question of trade-offs between cost (of obtaining a certain type of relative measurement) and benefit (improvement in localization accuracy) for various types of relative measurements is considered.

KEYWORDS: Robot localization; Multi-robot systems; Computer vision; Mobile robots; SLAM.

1. Introduction

In recent years, interest in utilizing teams of autonomous mobile robots has grown rapidly. Multi-robot teams are beneficial in many ways. Utilizing a group of low cost robots may be more economical than risking a single, more costly robot. In search and rescue operations, a group of robots can cover a larger area than a single robot. In hazardous conditions, the innate redundancy of a group of robots may be necessary to prevent catastrophic loss of mission capability. Regardless of the application, localization is a crucial task for any autonomous mobile robot team.

Localization, the estimation of an autonomous robot's global position and orientation, can be accomplished using a variety of sensors. Some of the more common sensors include Inertial Measurement Units (IMUs), vision-based sensors (cameras, LIDARs), and Global Positioning System (GPS). Of the three, GPS is the only senor capable of providing global measurements of a robot's position. However, in many situations, GPS measurements may not be available, or may only be intermittently available. For example, a group of unmanned aerial vehicles (UAVs) operating in an urban environment may temporarily lose GPS measurements when the signal is blocked by large buildings. In such a situation, the global pose (position and orientation) can be obtained by integrating

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over the relative pose measurements obtained using IMUs or vision-based sensors. This method of localization is referred to as *dead reckoning*.

Localization through dead reckoning can lead to a rapid growth in localization error.^{19,29} When a team of robots is deployed, relative measurements between pairs of robots may be available. For instance, a robot equipped with a monocular camera can measure another robot's bearing with respect to itself whenever the other robot is in its field of view. These measurements, although noisy, provide additional information on the robot's poses that can be used to improve localization accuracy.

In this paper, we propose a method for fusing measurements of various types to perform collaborative localization that improves localization accuracy over dead reckoning. We assume that all robots are equipped with proprioceptive sensors (cameras, IMUs, etc.) allowing each robot to measure its change in pose between time steps. We refer to these noisy measurements as *inter-time relative pose measurements*. In the absence of any other information, each robot can perform localization through dead reckoning using these noisy measurements. We further assume that each robot is equipped with exteroceptive sensors so that noisy relative measurements between robots may become available intermittently. These measurements, which we call *inter-robot relative measurements*, can be one (or a combination) of the following: relative pose, relative orientation, relative position, relative bearing, and relative distance, between a pairs of robots. We provide a method to perform collaborative localization by fusing the inter-time and inter-robot relative measurements to obtain an estimate of the absolute pose of every robot.

We provide both centralized and distributed algorithms for collaborative localization. In the centralized algorithm, all the measurements are assumed to be instantly available to a central processor. In the distributed algorithm, each robot performs localization using measurements from on-board sensors and information exchanged with neighboring robots. The complexity of the computations performed by a robot is only a function of the number of its neighbors at any given time, not of the total number of robots. This makes the distributed algorithm scalable to arbitrarily large groups of robots. In addition, the communication requirements are small. At every update, a pair of neighboring robots needs to exchange only (i) the relative measurement between them, and (ii) their current pose estimates.

Both simulations and experimental results (with a pair of P3-DX robots) are presented. In each case, the Monte Carlo method was used to empirically obtain the bias and standard deviation of estimation error by conducting multiple trials. Results show that the proposed collaborative algorithm substantially improves upon dead reckoning even when the team consists of a small number of robots. An examination of improvement versus measurement type provides a basis for deciding whether the cost of the sensors required for obtaining certain types of measurements is commensurate with the localization accuracy they provide.

1.1. Related work and contributions

Collaborative localization has been considered in the context of collaborative simultaneous localization and mapping (C-SLAM). Robots exchange local maps which are aligned and merged to improve robots' location estimates as well as to improve the maps (see refs. [2, 14, 28] and references therein). This requires the ability to identify common features in distinct maps generated by the robots. Recognizing common landmarks in distinct maps is often challenging. In addition, exchanging image data or maps between robots requires high bandwidth communication.

The second body of work therefore considers the collaborative localization problem as one in which only relative measurements (of pose, position, orientation, etc.) between pairs of robots are obtained and used to improve localization accuracy over self-localization. Our work falls into this category, which we term *collaborative localization from relative measurements*. Collaborative localization from relative measurements without map building is potentially useful to autonomous agents operating over large areas where collaborative mapping may be infeasible due to large computation, memory, and communication requirements.

Two features of the problem of collaborative localization from relative measurements make it challenging: (i) The nonlinear relationship between the variables to be estimated (absolute positions and orientations) and measurements (relative position, orientation, bearing, distance, etc.), and (ii) the non-commutativity of rotations in 3D.

Two distinct approaches to localization of single or multi-robot teams are relevant to the discussion of prior art: (i) pose-graph optimization methods, and (ii) filtering approaches.

The pose graph-based methods (also known as network or view-based representations) have been popular and successful in SLAM (see ref. [20] and references therein) and bundle adjustment.³⁴ A pose graph is a graph consisting of nodes that correspond to unknown poses of robots at various time instants and landmarks, and edges that correspond to observations between robot-landmark pairs. The best set of pose estimates are estimated by minimizing a cost function that quantifies how well a given set of poses predict the measurements. Under certain Gaussian assumptions on the measurement noise, the least square solution obtained is the maximum likelihood (ML) estimate of the poses. With additional statistical information on prior belief of the poses, again involving Gaussianity, the least square solution turns out to be the maximum *a posteriori* estimate (MAP) as well. Pose-graph optimization methods used in bundle adjustment and single robot SLAM can be directly applied to the multi-robot case. A centralized algorithm for ML collaborative localization through pose-graph optimization is presented in ref. [13]. Distributing the computations among the robots presents additional challenges;¹⁵ addresses them by using relative pose graphs. The MAP estimator for 2D collaborative localization proposed in ref. [27] also belongs to this category. In ref. [17] a pose graph-based algorithm for distributed collaborative localization in 3D is proposed that uses a linearization of the relation between the pose measurements involving rotations.

When robots' absolute orientations are known, the problem of cooperative localization through pose-graph optimization becomes a linear estimation problem.^{4,33} The problem we consider is, however, nonlinear since orientations are not known. In ref. [3], the authors use a two-phase approach to solve the 2D collaborative localization problem, where each phase is solved using linear estimation techniques: First, the absolute orientations are estimated using relative orientation measurements, then these are used along with relative position measurements to obtain absolute position estimates, and finally an improved estimate of the absolute orientations are obtained from the position estimates. This method is restricted to 2D localization.

The method proposed here also belongs to the pose graph-based approach: the estimation problem is formulated as an optimization defined by a graph, where nodes represent robot poses at various times and edges represent inter-time and inter-robot measurements. There is a subtle – but major – distinction between existing pose-graph optimization methods, which we call the Euclidean Pose-Graph Optimization (EPGO), and our proposed method. While existing methods use a vector space parameterization of orientation (such as the complex part of the unit quaternion representation) and then apply vector-space minimization techniques to search for the minima, we perform optimization directly on the product Riemannian manifold in which the problem is naturally posed without relying on a specific parameterization. A gradient-descent method on the Riemannian manifold is then used for searching for the minima. The gradient descent algorithm is independent of parameterization as well; any parameterization of the orientations can be used for numerical implementation without affecting the solution obtained. The advantages of doing so are discussed in detail in Section 3 (see Remark 1). Simulations show that the proposed method based on the Riemannian Manifold Pose-Graph Optimization, which we call RPGO, outperforms EPGO, the traditional pose-graph optimization performed by vector-space-based methods, when the measurement noise is large.

For EPGO in 3D, various tricks have to be employed to ensure that the result from vectorspace optimization leads to accurate results (see ref. [10] and references therein). The Kalman filter (KF)-based approaches require, in addition to a specific vector-space parameterization of rotations, linearization of nonlinear relationships between variables and measurements. Indeed, most of the papers on collaborative localization consider the 2D case: a robot's pose is described by three scalars: x, y-coordinates and an angle θ . Rotations in 2D commute and a vector-space parameterization does not lead to serious difficulties. Although many of these 2D algorithms can be extended to the 3D case in principle, the extensions are far from straightforward. In fact, no details are provided for the extension to 3D case; all simulation and experimental results in refs. [3, 8, 13, 21, 32, 33] are for the 2D case. The belief propagation approaches of refs. [8, 13, 21], although developed without a specific parameterization, are feasible only if a vector-space parameterization of the orientations is used. Otherwise belief propagation of orientation estimates requires representation and computation of conditional probability density functions in SO(3), a particularly challenging problem in itself. Some progress in fusion and propagation of pdfs in SO(3) has been reported in refs. [22, 36].

There is a large body of work on collaborative localization from relative measurements using a filtering-based approach in both centralized and distributed settings. Extended Kalman Filter (EKF)-based formulations have been adopted in refs. [25, 30, 32]. Collaborative localization of two aircrafts

from *indirect* relative measurement (observing common feature points on the ground) using the EKF has been proposed in ref. [26]. The papers^{8, 12, 31} propose particle-filter-based methods for collaborative localization. Leung *et al.*²¹ propose the Bayesian framework such that many filtering-based methods can be applied within it. They propose decentralized algorithms to update a robot's belief of its own pose that is equivalent to the centralized belief whenever a certain Markovian property holds.

In comparison with the KF-based methods described above, our method offers a distinct advantage. The linearization involved in the EKF requires a small angle approximation to hold at all times. Unless the time interval between two successive inter-vehicle measurements is extremely small, which is unlikely in many practical situations, the small angle approximation is violated, which is likely to lead to poor covariance updates and poor pose estimates. This is verified through simulations presented here.

This work, as compared with much of the earlier work on collaborative localization from relative measurements, differs in a number of novel ways. The proposed approach is (i) directly applicable to 3D pose estimation, (ii) able to utilize heterogeneous measurement types (of the relative position, orientation, bearing, distance, or any combination thereof) between pairs of robots, and (iii) able to handle a time-varying neighbor relationship in the distributed setting. To the best of our knowledge, although earlier works have some of these features, no existing method has all three features.

Ability to fuse heterogeneous relative measurements provides an important functionality. In particular, many of the papers mentioned above^{3, 8, 10, 13, 32} require measurements of relative pose, i.e., position and orientation, between robot pairs. While relative position can be obtained through stereo vision or laser range finders, obtaining relative orientation is quite challenging. It requires robots to be equipped with recognizable targets with known geometry, but even then orientation measurements can be quite noisy. In fact, obtaining relative position measurements with stereo vision is also challenging due to small baseline enforced by the size of robots. Laser range finders have limited range and therefore are not suitable to, say, UAVs, where the distance between two UAVs may be large. However, even if only one camera of a robot sees a neighboring robot, a bearing measurement can be obtained. Or, radio frequency (RF)-based techniques, such as time of arrival (TOA) measurements, can be used to measure distance during wireless communication. With the ability to fuse all types of relative measurements, all available measurements can be utilized for improvement in localization accuracy. Although Martinelli *et al.*²⁵ and Rekleitis *et al.*³¹ consider the case of heterogeneous measurements, their algorithms are limited to the 2D case. Our work extends to 3D, the comparison between various measurement types initiated in refs. [25, 31].

The third contribution of the paper is the low communication requirements of the proposed algorithm. A crucial requirement in some of the prior works^{3, 7, 13, 25, 27, 32} on collaborative localization is that of constant all-to-all communication among robots. In contrast, the proposed distributed algorithm allows the number of neighbors of a robot (those it can exchange information with) to vary in an arbitrary manner with time. In the extreme case when the number of neighbors is always 0, the proposed algorithm simply degenerates into dead reckoning.

Some preliminary aspects of the proposed method were reported in ref. [18]. Compared with that paper, we make several novel contributions here. While the method in ref. [18] required all inter-robot measurements to be of the relative pose, here we extend the methodology to enable fusion of many types of measurements. As discussed above, this represents a significant extension. We also present detailed proofs of all the mathematical results, which were not present in ref. [18]. Finally, here we include substantially more simulation and experimental results than in ref. [18].

The rest of the paper is organized as follows. The problem is precisely stated in Section 2. A centralized estimation scheme is described in Section 3, and a distributed algorithm that is inspired by the centralized scheme is described in Section 4. In Section 5, two existing state-of-the-art algorithms are reviewed briefly. Simulation and experimental results with the proposed algorithms are presented in Sections 6 and 7, respectively. The paper concludes with a discussion in Section 8.

2. Problem Statement

2.1. The collaborative localization problem

Consider a group of r mobile robots indexed by i = 1, ..., r. Time is measured by a discrete counter k = 0, 1, 2, ... Each robot is equipped with a local, rigidly attached frame of reference, that is, a

coordinate system defined in the robot's local reference frame. We call the frame of reference attached to robot *i* at time *k* frame *i*(*k*). Between any two frames of reference, say frames *u* and *v*, we denote the Euclidean transformation from *v* to *u* by \mathbf{T}_{uv} , where \mathbf{T}_{uv} is an element of the special Euclidean group SE(3). Specifically, if p_u is a point expressed in frame *u* and p_v is the same point expressed in frame *v*, then $p_u = \mathbf{T}_{uv}p_v$. We call \mathbf{T}_{uv} the *relative pose* of frame *v* with respect to frame *u*.

Let frame 0 denote some fixed frame of reference that is common to all robots. The *absolute pose* of frame *u* is then given by the transformation \mathbf{T}_{0u} . We will often denote the absolute pose simply as \mathbf{T}_u . Robot *i* is said to be *localized* at time *k* when an estimate is known for the absolute pose of frame *i*(*k*). We denote such an estimate by $\hat{\mathbf{T}}_{i(k)}$.

In this work, we consider the case when measurements of a robot's absolute pose, perhaps from the measurements of a GPS and compass, are either not available or only rarely available. Instead, each robot is equipped with proprioceptive sensors such that, at every time k, the robot is able to obtain a relative pose measurement with respect to its previous pose. That is, a robot i at time k is able to measure the relative pose $\mathbf{T}_{i(k-1)i(k)}$. We refer to these measurements as *inter-time relative pose measurements*. Such measurements can be obtained with inertial sensors or vision-based sensors. In addition, they need not be obtained from a sensor alone. Instead, a measurement could also be the estimate computed by fusing sensor measurements with predictions of the robot's motion from a dynamic/kinematic model.

In addition to the aforementioned proprioceptive sensors, each robot is equipped with exteroceptive sensors so that occasionally robot i is able to obtain a relative measurement of one or more other robots. We call these inter-robot relative measurements. If robot i collects a measurement of robot j at time k, it can be one of the following:

- **Relative pose**: The Euclidean transformation from frame *j*(*k*) to frame *i*(*k*); denoted by the symbol **T**.
- **Relative orientation**: The element of *SO*(3) that describes change in orientation from frame *j*(*k*) to frame *i*(*k*); denoted by the symbol **R**.
- **Relative position**: The vector in \mathbb{R}^3 that describes change in position between the frames i(k) and j(k), expressed in frame i(k); denoted by the symbol **t**.
- **Relative bearing**: The vector of unit length that points from frame *i*(*k*) to frame *j*(*k*), expressed in frame *i*(*k*); denoted by the symbol *τ*.
- **Relative distance**: The distance between frames i(k) and j(k); denoted by the symbol δ .

Which pairs of robots will be able to obtain an inter-robot relative measurement will depend on many factors, including the kind of sensors they have on board, the range of sensors, etc. An implicit assumption here is that a robot is able to uniquely identify another robot of which it obtains a relative measurement, so that there is no ambiguity on which pair of robots a relative measurement corresponds to.

The *collaborative localization problem* is the problem of estimating the pose of every robot at the current time k with respect to the common frame of reference by utilizing the inter-time and inter-robot measurements collected up to time k.

The above situation is best described in terms of a directed, time-varying, fully labeled graph $\mathcal{G}(k) = (\mathcal{V}_0(k), \mathcal{E}(k), \ell(k))$, where nodes $\mathcal{V}_0(k)$ correspond to variables, and edges $\mathcal{E}(k)$ correspond to measurements, and shows how the noisy relative measurements relate to the absolute pose of each robot at every time step. The graph is defined as follows. For each robot $i \in \{1, \ldots, r\}$ and each time $k' \leq k$, a unique index (call it u) is assigned to the pair (i, k'). How this indexing is done is immaterial. The indices $\{1, \ldots, rk\}$ define a set $\mathcal{V}(k)$ that is a subset of the *node set* $\mathcal{V}_0(k)$ of the graph. We refer to the frame of reference attached to robot i at time k' as frame u, where u is the node assigned to the pair (i, '). We introduce another node, denoted by 0, that corresponds to the common frame of reference in which every robot's pose is to be determined. Node 0 is called the grounded node. The node set of the graph is then defined as $\mathcal{V}_0(k) := \mathcal{V}(k) \cup \{0\}$. The relative pose of frame u with respect to frame 0 is denoted simply by \mathbf{T}_u . We call the poses $\{\mathbf{T}_u\}_{u \in V(k)}$ the node variables of $\mathcal{G}(k)$.

The set of *directed edges* at time k, denoted as $\mathcal{E}(k)$, corresponds to the noisy inter-time and inter-robot measurements collected up to time k. That is, suppose robot i is able to measure robot j's relative pose at time k', and let u, v be the nodes corresponding to robots i, j at time k', respectively. Then for all $k \ge k'$, there exists a directed edge $e \in \mathcal{E}(k)$ corresponding to this measurement. Since robot i measures robot j, the edge e leaves node u and arrives at v, we denote this by $e \triangleright (u, v)$.

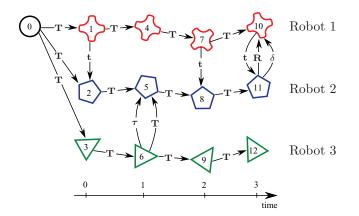


Fig. 1. A time history of three robots up to time k = 3 with inter-time and inter-robot relative measurements. Each (robot, time) pair is labeled with the corresponding node index from $V_0(3)$. Arrows indicate edges in $\mathcal{E}(3)$, i.e., relative measurements. Each edge is labeled to indicate the type of measurement. Each robot had GPS and compass measurements at the initial time k = 0. Thereafter, no other GPS or compass measurements were available.

Similarly, each inter-time relative pose measurements of a robot also creates an edge in the graph. To delineate the type of measurement, a label from the set { **T** (pose), **R** (orientation), **t** (position), τ (bearing), δ (distance) } is attached to each edge. The map from the set of edges to the set of labels is denoted by $\ell(k)$. For each edge, $e \in \mathcal{E}(k)$, where $e \triangleright (u, v)$, if $\ell(k)(e) = s$, which means there is a measurement of type *s* for frame *v* with respect to frame *u*, where $s \in \{\mathbf{T}, \mathbf{R}, \mathbf{t}, \tau, \delta\}$. The noisy relative measurement associated with edge $e \triangleright (u, v)$ is denoted by $\hat{\mathbf{T}}_{uv}, \hat{\mathbf{T}}_{uv}, \hat{\mathbf{T}}_{uv}$, or $\hat{\delta}_{uv}$ for $\ell(k)(e) = \mathbf{T}$ (pose), **R** (orientation), **t** (position), τ (bearing), or δ (distance), respectively. A measurement of type **T** is really two measurements, one of type **R** and another of **t**. We still use the nomenclature "measurement of type **T**" for ease of comparison with prior work, since relative pose measurements are commonly considered in the existing literature.

If the absolute pose in global GPS coordinates of at least one robot is known at time 0 through the use of GPS and compass, then node 0 can be associated with a terrestrial reference frame. When such measurements are not available, node 0 could correspond to the initial frame of reference of one of the robots. In either case, estimating the node variables is equivalent to determining the robots' poses with respect to the frame of the grounded node 0. Without the grounded node, the problem of localization from relative measurements is indeterminate up to a rotation and a translation.

The graph $\mathcal{G}(k)$ is called the *measurement graph* at time k. Figure 1 shows an example of the graph corresponding to the measurements collected by three robots up to time index 3. Because each robot may be equipped with more than one sensor, multiple distinct edges may exist between a pair of nodes.

To ensure that at least one estimate exists for every robot at each time k, we make the following assumption.

Assumption 1. Each robot has access to an estimate of its absolute pose at time 0.

Due to Assumption 1, an estimate of the pose of robot *i* at any time *k* (equivalently, the node variable T_u , where node *u* corresponds to the pair (i, k)) can be computed by composing the intertime relative pose measurements obtained by robot *i* up to time *k*. This estimate is equivalent to robot *i* performing dead-reckoning. In practice, Assumption 1 holds if all robots have a GPS and compass measurement at time 0. If no GPS measurement is available, but each robot can obtain a relative pose measurement with respect to, say, robot 1, then again the assumption holds.

Often many more edges are present in the graph $\mathcal{G}(k)$ than those necessary to form a single estimate of the node variables; robots can benefit from collaborative localization in such a scenario. As an illustrative example, consider the scenario shown in Fig. 1. The path (0, 1, 4, 7, 10) provides the dead-reckoning estimate of robot 1 at time 3, or equivalently an estimate of \mathbf{T}_{10} . Similarly, the path (0, 2, 5, 8, 11) provides as estimate for the node variable \mathbf{T}_{11} . In addition, the three edges between nodes 10 and 11 corresponding to relative position, orientation, and distance measurements

all provide additional information about how T_{10} and T_{11} relate. Because there is noise in each relative measurement, by incorporating the information found in these three edges, we expect to get a better estimate of both T_{10} and T_{11} . The goal of collaborative localization is to utilize all edges in the graph $\mathcal{G}(k)$ to improve over any estimates of the node variables that are obtained using only a single path. When the measurements are linearly related to the node variables, this can be accomplished by using the Best Linear Unbiased Estimator, as done in refs. [3, 4]. In our case, the relationship between the measurements and the node variables is nonlinear.

The discussion above also shows that Assumption 1 is not necessary, but merely sufficient for localization. For robot *i* to be localized, all that is needed is that there exists a time *k* so that there is an undirected path from node *u* (where *u* corresponds to *i*, *k*) to node 0 such that the edges along this path are of type **T**. In that case, an estimate of \mathbf{t}_u can be obtained by concatenating the relative pose measurements along the path from 0 to *u*, and hence robot *i* is localized at time *k*. After that, even if no inter-robot relative measurements are available, robot *i* can perform dead reckoning. Assumption 1 is taken in order to simplify the presentation.

2.2. The distributed collaborative localization problem

The problem stated above does not put any restriction on access to the measurements collected up to time k. In particular, a method that assumes that all measurements collected by all the robots up to time k are instantly available to a central computer is allowed. However, when a large number of robots are involved, with communication accomplished through the use of low bandwidth wireless links with limited range, such a centralized scheme is not feasible. In addition, retaining past measurements indefinitely will quickly exhaust both available memory and available processing capabilities of a centralized computing unit.

We now modify the problem by including constraints on communication and computation. In particular, a robot is now required to localize itself by using information that is available from onboard sensors and data it can collect from its *neighbors*. Two robots i and j are said to be neighbors at time k if they can communicate at time k.

The *distributed* collaborative localization problem is the following: Each robot is to localize itself by using measurements collected by on-board sensors and information it can obtain by communicating with its neighbors. To simplify the development of the distributed algorithm, the following assumption is made.

Assumption 2. If robot i can obtain a relative measurement of robot j at time k, then i and j can communicate at that time.

Although this assumption may seem strict, it can, in fact, always be satisfied: robot i simply drops any measurements involving j if it cannot communicate with j.

3. Centralized Collaborative Localization Algorithm

In this section, we present a solution to the collaborative localization problem where all the relative measurements are instantly available to a central processor at each time k. To distinguish this algorithm from other pose-graph methods discussed in the sequel, we will refer to it as the RPGOalgorithm. The centralized solution naturally leads to a distributed scheme, which will be described in the next section.

3.1. Localization through optimization over a graph

Instead of addressing the problem of estimating the robots' current poses at time k, we examine the more general problem of estimating all the node variables $\{\mathbf{T}_u\}_{u\in\mathcal{V}(k)}$ of the measurement graph $\mathcal{G}(k)$ using the noisy relative measurements, $\{\hat{\mathbf{M}}_e\}_{e\in\mathcal{I}(k)}$, where for each $(u, v) \triangleleft e \in E(k)$, $\hat{\mathbf{M}}_e$ is a measurement from node u to node v of one of the following: relative pose, orientation, position bearing, or distance.

In the following, we will interchangeably refer to both the absolute pose \mathbf{T}_u and its corresponding orientation and position, \mathbf{R}_u and \mathbf{t}_u respectively, as node variables. Similarly, a pose measurement $\hat{\mathbf{T}}_{uv}$ will be identified with its corresponding orientation and position measurements, $\hat{\mathbf{R}}_{uv}$ and $\hat{\mathbf{t}}_{uv}$.

We pose the collaborative localization problem as an optimization of a cost function over the set of node variables, where the cost function measures how well a given set of absolute poses explain the noisy measurements collected up to time k. The initial condition for each node variable $\mathbf{T}_u = (\mathbf{R}_u, \mathbf{t}_u), u \in \mathcal{V}(k)$ is given by the dead-reckoning estimate, whose existence is assured by Assumption 1.

Given a measurement $\hat{\mathbf{M}}_e$ between nodes u and v, let $c_e(\mathbf{R}_u, \mathbf{R}_v, \mathbf{t}_u, \mathbf{t}_v | \hat{\mathbf{M}}_e) \in \mathbb{R}_+$ denote the cost of the measurement $\hat{\mathbf{M}}_e$ as a function of the node variables. A suitable cost function must be chosen for each of the four measurement types (excluding pose measurements and instead considering a orientation/position pair) such that a higher cost indicates more disagreement between the observed measurements and the estimated node variables. This is done by considering the following fact. When the node variables and measurements are know exactly, each measurement can be given as a function of the node variables as follows:

$$\hat{\mathbf{R}}_{u\,v} = \mathbf{R}_{u}^{T} \mathbf{R}_{v}, \qquad \hat{\boldsymbol{\tau}}_{u\,v} = \mathbf{R}_{u}^{T} (\mathbf{t}_{v} - \mathbf{t}_{u}) / \|\mathbf{t}_{v} - \mathbf{t}_{u}\|, \\
\hat{\mathbf{t}}_{u\,v} = \mathbf{R}_{u}^{T} (\mathbf{t}_{v} - \mathbf{t}_{u}), \qquad \hat{\delta}_{u\,v} = \|\mathbf{t}_{v} - \mathbf{t}_{u}\|.$$
(1)

Since noise is present in the measurements, how much a noisy measurement differs from its value given as a function of the node variables provides a suitable cost.

In (1), $\hat{\mathbf{t}}_{uv}$, $\hat{\boldsymbol{\tau}}_{uv}$, and $\hat{\delta}_{uv}$ are all elements of a real vector space (\mathbb{R}^3 or \mathbb{R}) and so the standard Euclidean norms can be used to measure the distance between the noisy measurements and the equivalent node variable representations. In contrast, both $\hat{\mathbf{R}}_{uv}$ and $\mathbf{R}_u^T \mathbf{R}_v$ are elements of the manifold SO(3) and are thus considered to be abstract operators and not equated to any particular parameterization (such as rotation matrices). To this end, a suitable distance function on the manifold SO(3) must be chosen to measure their relative distance. Such a function is given by the Riemannian distance,

$$d: SO(3) \times SO(3) \to \mathbb{R}_+, \quad (p,q) \mapsto \sqrt{-\frac{1}{2} \operatorname{Tr}\left(\log^2(p^T q)\right)}, \tag{2}$$

where Tr denotes the trace function. More details on this distance function can be found in Appendix A. The cost function at each time k is chosen as a sum of edge costs over all edges (measurements):

$$f\left(\left\{\mathbf{T}_{u}\right\}_{u\in\mathcal{V}(k)}\right) := \sum_{(u,v)\lhd e\in\mathcal{E}(k)} c_{e}(\mathbf{R}_{u}, \mathbf{t}_{v}, \mathbf{R}_{v}, \mathbf{t}_{v} \mid \hat{\mathbf{M}}_{e}),$$
(3)

where $c_e(\mathbf{R}_u, \mathbf{t}_u, \mathbf{R}_v, \mathbf{t}_v | \hat{\mathbf{M}}_e)$, the cost for edge $e \triangleright (u, v)$, is given by

$$c_{e}(\mathbf{R}_{u}, \mathbf{t}_{u}, \mathbf{R}_{v}, \mathbf{t}_{v}) := \frac{1}{2} \\ \times \begin{cases} \frac{d^{2}(\hat{\mathbf{R}}_{u\,v}, \mathbf{R}_{u}^{T}\mathbf{R}_{v}) + \|\hat{\mathbf{t}}_{u\,v} - \mathbf{R}_{u}^{T}(\mathbf{t}_{v} - \mathbf{t}_{u})\|^{2} & \text{if } \ell(k)(e) = \mathbf{T} \\ \frac{d^{2}(\hat{\mathbf{R}}_{u\,v}, \mathbf{R}_{u}^{T}\mathbf{R}_{v}) & \text{if } \ell(k)(e) = \mathbf{R} \\ \frac{1}{\|\hat{\mathbf{t}}_{u\,v} - \mathbf{R}_{u}^{T}(\mathbf{t}_{v} - \mathbf{t}_{u})\|^{2} & \text{if } \ell(k)(e) = \mathbf{t} \\ \frac{1}{\|(\hat{\mathbf{t}}_{u\,v}\|\mathbf{t}_{v} - \mathbf{t}_{u}\|) - \mathbf{R}_{u}^{T}(\mathbf{t}_{v} - \mathbf{t}_{u})\|^{2} & \text{if } \ell(k)(e) = \mathbf{\tau} \\ \frac{1}{\|(\hat{\boldsymbol{\delta}}_{u\,v} - \|\mathbf{t}_{v} - \mathbf{t}_{u}\|)\|^{2}} & \text{if } \ell(k)(e) = \delta \end{cases}$$

$$(4)$$

and $\|\cdot\|$ indicates the standard Euclidean norm on \mathbb{R}^3 .

If the relative measurements were completely error-free, the minimum value of the cost function would be 0. By minimizing the cost function, we expect to find an improved estimate for the absolute pose of each robot over what can be determined through dead reckoning alone. The cost function in (3) is similar to one proposed in ref. [35] for a static camera network; here the cost function changes with time.

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3.2. Optimization over the product manifold

Finding the minimum of a function defined over a vector space has been studied extensively. However, the function $f(\cdot)$ in (3) is defined on a curved surface, specifically, the product Riemannian manifold $(SO(3) \times \mathbb{R}^3)^{n(k)}$, where $n(k) = |\mathcal{V}(k)|$, the cardinality of the set $\mathcal{V}(k)$. Although it is possible to parameterize the product manifold and thus utilize standard vector space optimization techniques, such a parameterization will inevitably introduce either an increase in dimensionality, and thus introduce constraints, or the parameterization will fail to be bijective over the entire domain. See ref. [37] for a discussion of the relevant parameterizations of SO(3) and their associated problems.

In contrast, we desire a method that allows us to solve an unconstrained optimization problem in which the solution remains on the manifold during all intermediate steps. We accomplish this through the use of a gradient descent algorithm on product manifold.

Gradient descent in a Riemannian manifold is analogous to gradient descent in a vector space in the following sense. Given a smooth real-valued function f defined on a manifold M, the gradient of f at $p \in M$, denoted as grad f(p), is a vector in the tangent space of M at p, which is denoted by T_pM . Just as in Euclidean space, grad f(p) points in the direction of the greatest rate of increase of f. An explicit expression for the gradient of the cost function (3) is provided in the next proposition; the proof is in Appendix C.

Proposition 1. The gradient of the cost function shown in (3) at $p = (\mathbf{R}_1, \mathbf{t}_1, \dots, \mathbf{R}_{n(k)}, \mathbf{t}_{n(k)}) \in (SO(3) \times \mathbb{R}^3)^{n(k)}$ is

$$grad f(p) = (grad f(\mathbf{R}_1), grad f(\mathbf{t}_1), \dots, grad f(\mathbf{R}_{n(k)}), grad f(\mathbf{t}_{n(k)})),$$

where for u = 1, ..., n(k),

$$gradf(\mathbf{R}_u) = \sum_{e \in \mathcal{E}(k)} grad c_e(\mathbf{R}_u) \qquad gradf(\mathbf{t}_u) = \sum_{e \in \mathcal{E}(k)} grad c_e(\mathbf{t}_u).$$

The gradients of the edge cost c_e with respect to each manifold, $grad c_e(\mathbf{R}_u)$ and $grad c_e(\mathbf{t}_u)$ for all $u \in \mathcal{V}(k)$, are provided in Appendix D.

Consider a function $f : S \to \mathbb{R}$ for some set *S*. When *S* is a vector space, the function *S* is minimized using gradient descent by iteratively updating an estimate $p_t \in S$ such that $p_{t+1} = p_t - \eta_t \operatorname{grad} f(p_t)$, where $\eta_t \in \mathbb{R}_+$ is chosen so that $f(p_{t+1}) < f(p_t)$.⁵ An analogous update law can be defined when *S* is instead a Riemannian manifold. The key difference is that, in general, the addition operator is not defined on the manifold. Rather, the *parallel transport* map can be used to traverse the manifold in the direction of $-\operatorname{grad} f$.

The parallel transport map on the product manifold $(SO(3) \times \mathbb{R}^3)^n(k)$ at a point

$$p := (\mathbf{R}_1, \mathbf{t}_1, \dots, \mathbf{R}_{n(k)}, \mathbf{t}_{n(k)}) \in (SO(3) \times \mathbb{R}^3)^{n(k)},$$
(5)

denoted by \exp_n , is given by

$$\exp_p(\xi) = \left(\mathbf{R}_1 \exp(\mathbf{R}_1^T \xi_{\mathbf{R}_1}), \mathbf{t}_1 + \xi_{\mathbf{t}_1}, \dots, \mathbf{R}_{n(k)} \exp(\mathbf{R}_{n(k)}^T \xi_{\mathbf{R}_{n(k)}}), \mathbf{t}_{n(k)} + \xi_{\mathbf{t}_{n(k)}}\right),$$
(6)

where $\xi = (\xi_{\mathbf{R}_1}, \xi_{\mathbf{t}_1}, \dots, \xi_{\mathbf{R}_{n(k)}}, \xi_{\mathbf{t}_{n(k)}})$ is an element of the *tangent space* $T_p[(SO(3) \times \mathbb{R}^3)^{n(k)}] = T_{\mathbf{R}_1}SO(3) \times \dots \times T_{\mathbf{t}_{n(k)}}\mathbb{R}^3$, and the exp(·) function appearing on the right hand side of (6) is the map exp : $\mathbb{L}(\mathbb{R}^3) \to \mathbb{L}(\mathbb{R}^3)$ defined by $exp(X) = \sum_{k=0}^{\infty} \frac{X^k}{k!}$ for $x \in \mathbb{L}(\mathbb{R}^3)$. The derivation of (6) is provided in Appendix B. The gradient descent law is

$$p_{t+1} = \exp_{p_t}(-\eta_t grad f(p_t)), \quad t = 0, 1, \dots,$$

where $\eta_t \ge 0$ is the step size for iteration *t*. The parameter η_t is chosen as the Armijo step size $\eta_t^{(A)} = \beta^{N_t} \alpha$, where N_t is the smallest non-negative integer such that

$$f(p_t) - f(exp_{p_t}(\beta^{N_t}\alpha \, grad f(p_t))) \ge \sigma \beta^{N_t} \alpha \|grad f(p_t)\|$$
(7)

for scalar tuning parameters $\alpha > 0$, $\beta, \sigma \in (0, 1)$. The norm $\|\cdot\|$ defined in the vector space $T_p[(SO(3) \times \mathbb{R}^3)^n]$ is given by

$$\|(\xi_{\mathbf{R}_{1}},\xi_{\mathbf{t}_{1}},\ldots,\xi_{\mathbf{R}_{n(k)}},\xi_{\mathbf{t}_{n(k)}})\|^{2} = \sum_{u=1}^{n(k)} \left(\frac{1}{2}\operatorname{Tr}\left(\xi_{\mathbf{R}_{u}}^{T}\xi_{\mathbf{R}_{u}}\right) + \xi_{\mathbf{t}_{u}}^{T}\xi_{\mathbf{t}_{u}}\right),$$

which comes from the *Riemannian metric*. More details on Riemannian metric can be found in Appendix A. Theorem 4.3.1 in ref. [1] guarantees that the iterates p_t converges to a critical point of the cost function f defined in (3) as $t \to \infty$.

The pseudo-code of the centralized RPGO algorithm is given in Algorithm 1, where $\epsilon > 0$ is a user-specified accuracy threshold. Correctness of the RPGO algorithm (convergence to a critical point of the cost function f in (3)) follows from Theorem 4.3.1.¹

It should be noted that a choice of parameterization is necessary to explicitly compute updates in the proposed gradient descent algorithm. However, because the gradient descent is carried out on the manifold, this choice does not impact the particular value of the node variables found at any step.

4. Distributed Collaborative Localization Algorithm

In this section, we propose an algorithm for solving the distributed localization problem. The algorithm requires limited memory, processor power, and communication bandwidth for its execution. To distinguish this algorithm from the centralized RPGO algorithm presented in the previous section, we will refer to it as the Distributed Riemannian Pose Graph Optimization algorithm, or D-RPGO.

For each robot *i*, let $N_i^{(+)}(k)$ denote the set of all robots $j \in \{1, ..., r\}$ such that, at time *k*, robot *i* can obtain a relative measurement with respect to *j*. Similarly, let $N_i^{(-)}(k)$ denote the set of all robots $j \in \{1, ..., r\}$ such that, at time *k*, robot *j* can obtain a relative measurement with respect to *i*. The *neighbors* of robot *i* at time *k* are then given by the set $N_i(k) = N_i^{(+)}(k) \cup N_i^{(-)}(k)$. Due to Assumption 2, robot *i* can communicate with its neighbors $N_i(k)$ during time *k*.

Consider the local measurement graph $\mathcal{G}_i(k) = (\mathcal{V}_i(k), \mathcal{E}_i(k), \ell_i(k))$ of robot *i*, whose node set is simply the neighbors of *i* at time *k* along with the grounded node 0 and *i* itself: $\mathcal{V}_i(k) = N_i(k) \cup \{0, i\}$. The edges of $\mathcal{G}_i(k)$ correspond to the inter-robot measurements at time *k* between *i* and its neighbors along with an edge $e \triangleright (0, j)$ for each $j \in \mathcal{V}_i(k)$ (see Fig. 2 for an example). Each node in the local measurement graph $\mathcal{G}_i(k)$ is associated with an absolute pose of a robot at time *k*. No past poses belong to this graph. An edge $(p, q) \triangleleft e \in \mathcal{G}_i(k)$ (where i = p or q) corresponds to an inter-robot relative measurement between robots *p* and *q* at time *k*. The additional edges $e \triangleright (0, j), j \in \mathcal{V}_i(k)$ correspond to the "initial" estimate robot *j*'s absolute pose at time *k*, denoted by $\hat{\mathbf{T}}_{0,j}(k)$. Each robot *j* obtains the estimate $\hat{\mathbf{T}}_{0,j}(k)$ at time *k* by concatenating its pose estimate obtained at time k - 1 with the noisy inter-time relative pose measurement describing its motion from k - 1 to *k*. The estimate $\hat{\mathbf{T}}_{0,j}(k)$ is then used as the measurement associated with edge $e \triangleright (0, j)$. The graph $\mathcal{G}_i(k)$ is now a measurement graph since each edge has an associated noisy relative measurement. The edge $(0, j) \triangleleft e \in \mathcal{E}_i(k)$ for $j \in \mathcal{V}_i(k)$ ensures that Assumption 1 is satisfied for the local measurement graph $\mathcal{G}_i(k)$ for each *i*.

The distributed algorithm works as follows. At each time k, every robot $i \in \{1, ..., r\}$ forms an initial estimate $\hat{\mathbf{T}}_{0i}(k)$ of its absolute pose as described above and obtains inter-robot relative measurements of each of its neighbors $j \in N_i^{(+)}(k)$. Robot *i* then transmits to each $j \in \mathcal{V}_i(k)$ its initial estimate of its absolute pose $\hat{\mathbf{T}}_{0i}(k)$ along with all inter-robot relative measurements between itself and *j* that it obtained at time *k*. Robot *i* receives in turn robot *j*'s estimate of its current absolute pose $\hat{\mathbf{T}}_{0j}(k)$ along with relative measurements involving itself that *j* collected at that time. Robot *i* then executes Algorithm 1 on the local measurement graph $\mathcal{G}_i(k)$. The unknown node variables on this graph consist of its own pose and the poses of its neighbors (all at time *k*). After the computation, only the estimate of *i*'s own current pose is retained in its local memory; all other computed values **Input:** $\mathcal{G}(k)$, all noisy measurements on $\mathcal{E}(k)$, an initial guess for each node variable $(\mathbf{R}_u, \mathbf{t}_u)$, $u \in \mathcal{V}(k)$. **Output:** $\{(\hat{\mathbf{R}}_u, \hat{\mathbf{t}}_u)\}_{u \in \mathcal{V}(k)}$ $\hat{p} \rightarrow \text{ initial guess } (p \text{ defined in (5)});$ **repeat foreach** $u \in \mathcal{V}(k)$ **do** $\mid \text{ Compute } grad f(\hat{\mathbf{R}}_u) \text{ and } grad f(\hat{\mathbf{t}}_u) \text{ for the cost } f \text{ in (3) (as shown in Appendix D);}$ **end** Determine $\eta^{(A)}$, the Armijo step size from (7) (with \hat{p} for p_t); **foreach** $u \in \mathcal{V}(k)$ **do** $\mid \hat{\mathbf{R}}_u \rightarrow \hat{\mathbf{R}}_u \exp(-\eta^{(A)} \hat{\mathbf{R}}_u^T grad f(\hat{\mathbf{R}}_u));$ $\hat{\mathbf{t}}_u \rightarrow \hat{\mathbf{t}}_u - \eta^{(A)} grad f(\hat{\mathbf{t}}_u);$ **end until** $\|grad f\left(\{(\hat{\mathbf{R}}_u, \hat{\mathbf{t}}_u)\}_{u \in \mathcal{V}(k)}\right)\| > \varepsilon;$

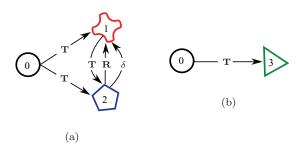


Fig. 2. The local measurement graphs for the robots of Fig. 1: (a) $\mathcal{G}_1(3)$ and $\mathcal{G}_2(3)$ (in this example they are the same), and (b) $\mathcal{G}_3(3)$.

are discarded. Since the distributed algorithm is simply Algorithm 1 applied to a local measurement graph, it inherits the correctness property of Algorithm 1. Note that if robot i has no neighbors at time k, the distributed collaborative localization algorithm is equivalent to performing self-localization from inter-time relative measurements.

The local measurement graphs $\mathcal{G}_1(3)$, $\mathcal{G}_2(3)$, and $\mathcal{G}_3(3)$ corresponding to the example of Fig. 1 are shown in Fig. 2. At time k = 3, robot 3 can see no other robots, so it will update its absolute pose using the inter-time relative pose measurement alone, without the aid of any inter-robot relative measurements. Robots 2 and 1, in contrast, will use the relative measurements between them obtained at time k = 3 to update their pose estimates.

4.1. Communication requirements

In the distributed algorithm, all the past data are truncated and replaced by the current estimate. As a result, the size of the problem data for each robot *i* at time *k* (encoded in the local subgraph $G_i(k)$), is bounded – and small – for all *i* and *k*. Specifically, for a robot *i* with N_i neighbors, the communication requirement of the distributed algorithm is $O(N_i)$. This can be seen as follows. A robot has to store current estimates of the node variables in its local measurement graph, and the number of nodes in that graph is $N_i + 1$. Each node variable can be represented by 6 numbers (3 for position and 3 for orientation). The robot also has to store the relative measurements between itself and its neighbors. The maximum memory requirement is when a robot is able to obtain measurements of all four types of relative measurements in the same time instant (relative position, orientation, bearing, and distance). In that case, each neighbor causes a requirement of 3 + 3 + 2 + 1 = 9 numbers to represent the interrobot measurements (2 for orientation in 3D, and 1 for distance). Thus, the total memory (storage) requirement for robot *i* is at most $6(N_i + 1) + 9N_i$. At every time step, a pair of neighboring robots has to exchange (i) the relative measurement(s) between them, and (ii) their current pose estimates:

at most 15 numbers. If information is broadcast to all neighbors, a likely scenario with autonomous robots communicating wirelessly, a robot only has to transmit $15N_i$ numbers.

4.2. Computational complexity

Just as the communication requirements scaled with the number of neighbors for a given robot *i*, so too does the computational complexity for a given robot's implementation of the D-RPGO algorithm. Consider a robot *i* at time *k*, and let $N_i(k)$ indicate the number of neighbors for robot *i* at time *k*. The dimensionality of the optimization problem is then equal to $N_i + 1$. In addition, the number of operations necessary to find the gradient for robot *i* is $O(N_i(k)^3)$. This follows from the fact that the gradient is a $N_i(k) + 1$ dimensional vector found by summing over the number of edges between neighbors, at most $N_i(k)^2$.

5. Existing Methods of Collaborative Localization

In this section, we briefly introduce two existing state-of-the-art methods for collaborative localization. The first algorithm we consider is based on standard pose-graph optimization methods (see ref. [20] and references therein). The second algorithm utilizes an Indirect Extended Kalman filter (IEKF) to perform collaborative localization, similar to the method used in ref. [26]. These two algorithms will be compared with the proposed algorithm through simulations in Section 6.3.

5.1. Pose-graph optimization

Unlike the algorithm presented in Section 3, standard pose-graph optimization techniques require a suitable parameterization of SO(3) to be used. For use in this paper, unit quaternions were chosen for this parameterization.⁶ For each node $u \in \mathcal{V}(k)$, let \mathbf{q}_u denote the unit quaternion corresponding to the node variable \mathbf{R}_u . Similarly, given a orientation measurement $\hat{\mathbf{R}}_{uv}$, let $\hat{\mathbf{q}}_{uv}$ denote the corresponding unit quaternion. Estimates for the node variables at time *k* are determined by minimizing the cost function,

$$f(\{\mathbf{q}_u, \mathbf{t}_u\}_{u \in \mathcal{V}(k)}) := \sum_{(u,v) \triangleleft e \in \mathcal{E}(k)} g_e(\mathbf{q}_u, \mathbf{t}_u, \mathbf{q}_v, \mathbf{t}_v)^T P_e g_e(\mathbf{q}_u, \mathbf{t}_u, \mathbf{q}_v, \mathbf{t}_v),$$
(8)

where $P_e > 0$ is a scaling matrix and $g_e(\mathbf{q}_u, \mathbf{t}_u, \mathbf{q}_v, \mathbf{t}_v)$ is a suitable vector edge error defined for each measurement type. Although standard pose-graph optimization is able to handle all measurement types considered thus far, for ease of exposition, only measurements of the relative position and orientation will be considered in the following discussion.

No canonical vector edge error exists for parameterization; many choices are possible. One suitable choice, which is the vector edge error used in the subsequent comparisons, is given as follows:

$$g_e(\mathbf{q}_u, \mathbf{t}_u, \mathbf{q}_v, \mathbf{t}_v) = \begin{cases} \mathbf{q}_u^{-1} \otimes \mathbf{q}_v \otimes \hat{\mathbf{q}}_{uv}^{-1} - 1 & \text{if } \ell(k)(e) = \mathbf{R} \\ C(\mathbf{q}_u)^T (\mathbf{t}_v - \mathbf{t}_u) - \hat{\mathbf{t}}_{uv} & \text{if } \ell(k)(e) = \mathbf{t} \end{cases},$$
(9)

where $e \triangleright (u, v)$, \otimes denotes the quaternion multiplication,⁶ and *C* denotes the map that takes a unit quaternion to its corresponding 3×3 rotation matrix representation. Many vector-space optimization algorithms can be used to search for (9). One common choice is to use the Levenberg–Marquadt algorithm, as was done in ref. [20]. To implement the Levenberg–Marquadt algorithm, a *minimal parameterization* is utilized, e.g., the first three components of the four-element unit quaternion. More details of the EPGO algorithm used for comparison with the RPGO algorithm can be found in ref. [16].

Remark 1. The orientations $(\mathbf{R}_{(\cdot)})$ that appear in our problem formations and throughout the development of the RPGO algorithm are abstract rotation operators, or elements of SO(3). No particular parameterization (rotation matrices, quaternions, etc.) is assumed during the development. In contrast, the cost function for the EPGO algorithm, (9) utilizes a specific parameterization in the form of unit quaternions. A different parameterization would necessarily lead to a different cost function, and possibly a different estimate delivered by the corresponding algorithm. The

estimates found using the EPGO algorithm are functions of both the measurements and the chosen parameterization, rather than a function of the measurements alone.

5.2. Indirect extended Kalman filter

We next consider an algorithm based on IEKF, similar to that used in ref. [26]. The two primary differences between the algorithm used in ref. [26] and the IEKF algorithm given below are as follows. (i) Inter-time relative measurements are of the relative pose, rather than measurements of linear and angular velocities. This is motivated by the idea that measurements are captured by a camera, rather than an inertial measurement unit. (ii) The IEKF algorithm presented here utilizes measurements of the relative position between robots, where as the algorithm presented in ref. [26] uses measurements of the bearing with respect to a time-varying set of feature points. Under these two changes, the IEKF algorithm is able to function under the same circumstances as the RPGO algorithm presented in Section 3.

As with the EPGO algorithm, and in contrast to the RPGO algorithm, the IEKF algorithm requires a suitable parameterization of SO(3) to be used. Unit quaternions are again chosen for this parameterization. A complete description of the resulting filtering equations is available in ref. [16]; here we omit the details due to lack of space.

6. Simulation Results

In this section, we present simulations comparing the centralized and distributed algorithms and study the effect of collaboration on localization accuracy using each of the inter-robot measurement types.

First, in Section 6.1, we consider the case that all inter-robot relative measurements are of the relative pose. We examine the difference in localization accuracy between the centralized and distributed algorithms, as well as the effect of increasing number of robots on localization accuracy. In Section 6.2, we consider the heterogeneous measurement case in which the inter-time relative measurements are of relative pose, but the inter-robot relative measurements may be of relative pose, orientation, position, bearing, or distance. We examine the effects these various measurement types have on the accuracy of location estimates.

To study the change in localization accuracy, we consider the following metrics. The *position* (*estimation*) *error* of robot *i* is $\mathbf{e}_i(k) := \hat{\mathbf{t}}_i(k) - \mathbf{t}_i(k)$, where $\mathbf{t}_i(k)$ is its global position at time *k* and $\hat{\mathbf{t}}_i(k)$ is the estimate of this position. The *bias* in the position estimation error of robot *i* is defined as $\|\mathbf{E}[\mathbf{e}_i(k)]\|$, where $\mathbf{E}[\cdot]$ denotes expectation and $\|\cdot\|$ refers to 2-norm, and the *standard deviation* as $\sqrt{\operatorname{Tr}(Cov(\mathbf{e}_i(k), \mathbf{e}_i(k)))}$, where $Cov(\cdot, \cdot)$ denotes the covariance.

6.1. Centralized RPGO versus DRPGO

To compare the centralized and distributed algorithms, we examine the localization of a group of five robots. All inter-robot relative measurements were of the relative pose. Each of the five robots travels along a distinct zig-zag path in 3D, as shown in Fig. 3(a). Two robots could obtain relative pose measurements at time k if the Euclidean distance between them at that time was less than 7 m. Due to Assumption 2, communication between robots is possible between at least those pairs with a distance of less than 7 m. Furthermore, 25% of these potential measurements were dropped to simulate random failure. A plot of the number of neighbors of robot 1 over time is shown in Fig. 3(b). The orientation measurements for each relative pose (both inter-robot and inter-time) were corrupted by independent identically distributed (i.i.d.) unit quaternions drawn from the Von Mises–Fisher distribution²⁴ centered around the zero-rotation quaternion and with a concentration parameter of 10, 000. Noise in the relative translation measurements was simulated by adding i.i.d. zero-mean normal random variables with a covariance matrix $I_{3\times3} \times 10^{-6}$.

Figure 4 shows the bias and standard deviation in the position error of a single robot acting in a five-robot team, estimated using a 100-iteration Monte Carlo simulation. The group of robots performed localizing using either the centralized or distributed collaborative localization algorithms with all inter-robot measurements being of the relative pose. The plots indicate that improvement in localization accuracy with the distributed algorithm is quite close to that with the centralized algorithm. This is promising since the distributed algorithm is applicable to large teams of robots in highly dynamic scenarios that can lead to arbitrary time variation in neighbor relationships. The

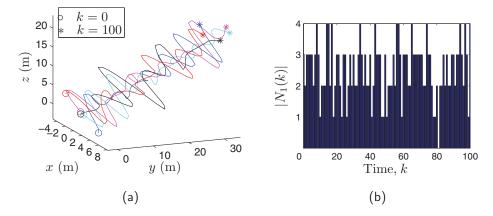


Fig. 3. (a) The 3D trajectories for robots 1 through 5, used in all simulations, and (b) the number of neighbors of robot 1 as a function of time.

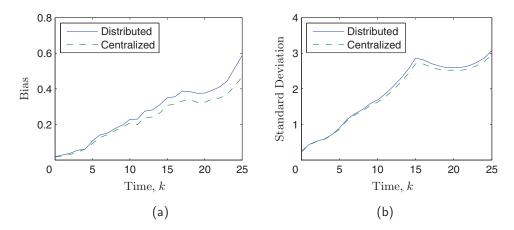


Fig. 4. The (a) bias and (b) standard deviation in position error for a robot acting in a group of five localizing using either the centralized or distributed algorithm. This estimate was obtained using a 100-iteration Monte Carlo experiment.

centralized algorithm is not applicable in a realistic setting, but it provides a measure of the best possible performance. From this point forward, we will only study the distributed algorithm.

We next conduct a series of Monte Carlo simulations with the distributed algorithm and compute means and variances of the resulting localization errors. One goal of this study is to verify that the accuracy improvements observed above are not a chance occurrence. Another goal is to examine how the localization accuracy varies with the number of robots in the team.

In each of the sample runs of the Monte Carlo simulation, the robots traveled the same paths as shown in Fig. 3(a). Measurement noise was introduced in the same manner as in the previous simulations, and varied randomly from run to run. Neighbor relations were again determined as described earlier, and kept the same from run to run to preclude that from being an additional source of randomness. Simulations for robot teams of size 1, 2, 3, 4, and 5 were carried out. In each case 1000 iterations were performed. When only one robot is present in the team, collaborative localization is equivalent to self-localization without the aid of any inter-robot relative pose measurement. As the number of robots in the team increases, the number of neighbors for a robot at any given time will tend to increase and so greater improvement in localization accuracy is expected.

The bias and standard deviation in the position estimation error $\mathbf{e}_i(k)$ for robot 1 (i = 1) are shown in Fig. 5. Both bias and standard deviation show significant improvement with distributed collaborative localization over self-localization. This is evident even for a team of only two robots. As the number of robots in the team increases, the localization error of robot 1 decreases. The improvement in accuracy, however, shows a diminishing return with increasing team size.

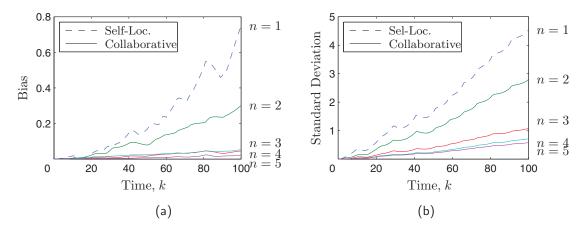


Fig. 5. The (a) bias and (b) standard deviation in the position estimation error for robot 1 utilizing the distributed algorithm. The curves correspond to the number of robots in the group: 1, 2, 3, 4, or 5.

6.2. Effect of measurement type

We now perform simulations for the case when inter-robot relative measurements are of various types (pose, orientation, position, bearing, and distance). To examine the effect that measurement type has on localization accuracy, in each experiment, we let all inter-robot relative measurements be of the same type. The Monte Carlo simulations were conducted, in which we again consider five robots traveling in zig-zag paths shown in Fig. 3(a). Errors in the pose, orientation, and position measurements are introduced as described in Section 6.1. Noise in the bearing measurements is induced by rotating the true bearing through the application of a unit quaternion generated from the i.i.d. Von Mises–Fisher distribution. The noise in the distance measurements is normally distributed. The bias and standard deviation of position error for robot 1 are shown in Fig. 6.

We see from the plots that improvement over self-localization occurs for all measurement types, with the exception of distance measurements. While distance measurements improve the standard deviation of the position estimates, they have little or no effect on the bias. That distance has little effect on the accuracy is consistent with the conclusions in the study²⁵ for 2D localization. The fact that bearing measurements lead to higher improvement over distance or position measurements was also observed in ref. [25] for the 2D case. While the conclusions in ref. [25] were based on single simulations, ours are based on Monte Carlo simulations.

Trade-offs between cost of sensors and the resulting benefit in localization can be analyzed from these empirically observed trends. Although full pose provides the most benefit to localization accuracy, it is clear that any of the considered inter-robot measurement types can be used to improve localization accuracy over dead reckoning. In particular, after relative pose, relative position seems to be the most valuable types of inter-robot measurements, leading to significant reduction in both bias and variance over dead reckoning. This means that the cost of having sensors capable of measuring relative position (stereo vision, laser range finder, or monocular camera-based bearing sensor along with an RF-based distance measurement) may very well be justified by the localization accuracy they lead to. On the other hand, it is also apparent that the improvement due to inter-robot bearing measurements is quite comparable to that due to inter-robot relative position measurements. Yet only a single camera is necessary to measure the bearing, where as (in general) stereo vision is necessary to measure the full relative position, which is still quite prone to large errors unless large baseline stereo is used. Thus, given cost, payload, and reliability constraints, monocular cameras might be a better choice than stereo vision. These conclusions come with the caveat that they have been drawn from one set of Monte Carlo simulations; more extensive studies are needed to establish how general these trends are.

6.3. Comparison with alternate methods of distributed collaborative localization

We now present simulations that provide some insight into how the D-RPGO algorithm performs when compared with two state-of-the-art collaborative localization algorithms. All inter-robot relative measurements are of the relative pose for these simulations.

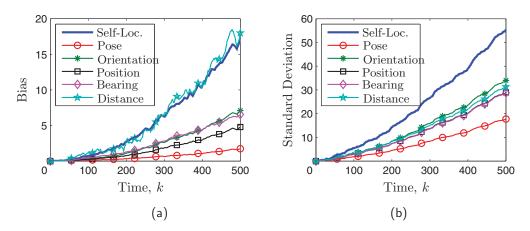


Fig. 6. Simulation: The (a) bias and (b) standard deviation in position error for robot 1 when the distributed algorithm is applied to a group of five robots utilizing inter-robot relative measurements. The labels "pose," "orientation," "position," "bearing," and "distance" indicate the group of robots used inter-robot measurements of the respective types to perform collaborative location using the distributed algorithm.

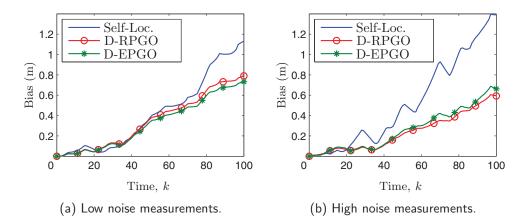


Fig. 7. Simulation: Comparison between D-RPGO and D-EPGO algorithms when (a) inter-robot orientation measurements are less noisy (K = 10,000) and (b) inter-robot orientation measurements are more noisy (K = 100), where K is the concentration parameter for the Von Mises–Fisher distribution from which the noisy measurements are drawn. The bias in position estimation error of robot 1 (in a group of five robots utilizing noisy inter-robot relative pose measurements), computed with both algorithms, averaging over the same graph, is shown.

The first alternative that we consider is the standard pose-graph implementation presented in Section 5.1, referred to as the EPGO algorithm. To maintain comparability, we provide the same local measurement graph to both D-RPGO algorithm and EPGO algorithm, which we call the D-EPGO algorithm.

A group of five robots are simulated to move along the 3D path described above. Errors in the pose measurements were induced as in simulations in Section 6.2. Simulations were performed varying the concentration parameter K in the Von Mises–Fisher distribution, from which the noisy rotations (quaternions) used to corrupt the inter-robot orientation measurements are drawn. These simulations show that when K is very large, that is, the variance is very low, D-EPGO does very well, even outperforming the D-RPGO algorithm. However, when K is small, that is, the noise variance is large, D-RPGO outperforms D-EPGO. Due to lack of space, only the results for K = 10,000 and 100 are shown; see Fig. 7. It is clear that the proposed D-RPGO algorithm outperforms the D-EPGO algorithm when K = 100 (more noise), while the D-EPGO algorithm outperforms the D-RPGO algorithm when K = 10,000 (less noise). The standard deviation is shown in Fig. 8. In the case of standard deviation, difference between the two algorithms is small.

Next, we consider the IEKF algorithm presented in Section 5.2. A pair of robots is simulated traveling along distinct sinusoidal paths in 3D space. Measurements are generated as described

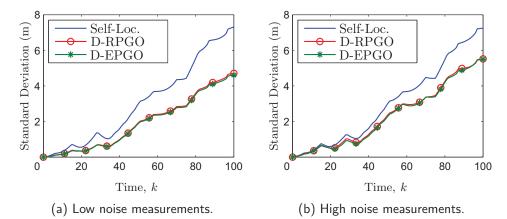


Fig. 8. Simulation: Comparison between D-RPGO and D-EPGO algorithms when (a) inter-robot orientation measurements are less noisy (K = 10,000) and (b) inter-robot orientation measurements are more noisy (K = 100), where K is the concentration parameter for the Von Mises–Fisher distribution from which the noisy measurements are drawn. The standard deviation in position estimation error of robot 1 (in a group of five robots utilizing noisy inter-robot relative pose measurements), computed with both algorithms, averaging over the same graph, is shown.

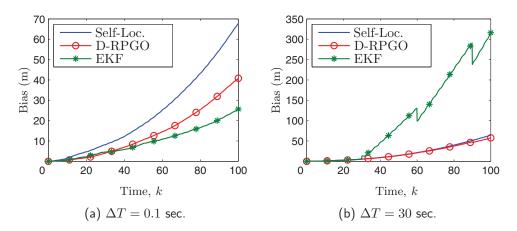


Fig. 9. Simulation: Comparison between D-RPGO and IEKF algorithms with two different sampling periods for inter-robot measurements. The bias in position estimation error of robot 1 (in a group of two robots utilizing noisy inter-robot relative pose measurements) computed with both algorithms.

earlier. The trends observed from extensive simulations can be summarized as follows. When the time interval between successive inter-robot measurements, call it ΔT , is small, the IEKF performs well as, or better than, the D-RPGO algorithm. However, when the time between measurements is large, the D-RPGO algorithm provides significantly better estimates of the robots' poses compared with the IEKF algorithm. Figure 9 provides numerical results for the case of a large ΔT (30 sec in this example), when IEKF performs poorly, and a small ΔT (0.1 sec), when the IEKF performs well. How small ΔT has to be for IEKF to perform well depends on many factors, including the motion of the robots, noise in the measurements, etc. For the parameters used in the simulations mentioned above, ΔT has to be smaller than 0.1 sec for the IEKF to perform as well as the D-RPGO algorithm.

We believe the reason for this behavior of the IEKF is the error introduced by the linearization involved in covariance propagation. The linearized state equations rely on the assumption that the angle between the true and estimated orientation is very small. When the time interval between inter-robot measurements is sufficiently small, this approximation holds. In that case the error in the covariance matrix due to linearization is small enough that it does not outweigh the added benefit of using covariance information. However, the small angle approximation is violated for large time intervals, leading to quite poor covariance estimates, which in turn lead to poor pose estimates.

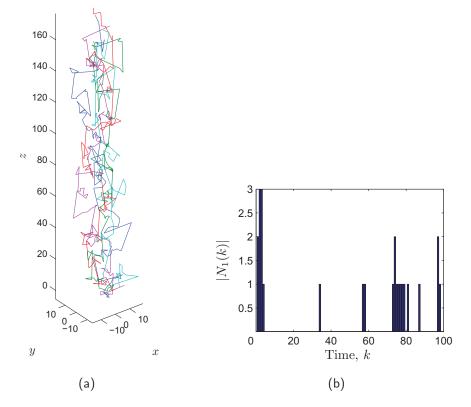


Fig. 10. (a) The 3D trajectories for robots 1 through 5 corresponding to one particular realization of the random path, and (b) the number of neighbors of robot 1 as a function of time for the given paths.

6.4. Application to randomized paths

Thus far the D-RPGO algorithm has only been applied to a single deterministic set of paths, which is necessary for the careful evaluation of algorithm's performance. To address any concerns about possible dependency on paths, we now present a simulation in which the paths of the robots are random. The random paths were generated such that each robot tended to follow the same moving set point. The probability of moving toward that point was scaled by the distance from it. One particular realization of such a path for each of the five robots is found in Fig. 10(a). The number of neighbors for robot 1 during the given path is shown in Fig. 10(b).

We again consider the case when all inter-robot measurements are of the relative pose. A 500iteration Monte Carlo simulation was conducted and the root-mean-square localization error was estimated for both self-localization (dead reckoning) and D-RPGO algorithm cases. The root-meansquare position estimation error is shown in Fig. 11.

Comparing these results with those presented in Fig. 5, we observe that improvement in localization accuracy by using the D-RPGO algorithm appears to be largely independent of the path. In addition, the results seen in Fig. 11 are indicative of an observation made early in this work: that the average number of neighbors, not the total number of robots, is the dominant factor determining localization improvement in the application of the D-RPGO algorithm. In contrast to our previous simulations, the average number of neighbors here proves to be $\ll 1$ and so we see improvements more akin to the r = 1 case in Section 6.1, rather than the r = 5 case.

7. Experimental Results

In this section, we present results for experiments conducted with two Pioneer P3-DX robots; these are shown in Fig. 12. Each robot was equipped with a calibrated monocular Prosillica EC 1020 camera and wheel odometers. Measurements from these sensors were fused to obtain the noisy inter-time relative pose measurements. Each robot is additionally equipped with a target allowing the on-bard

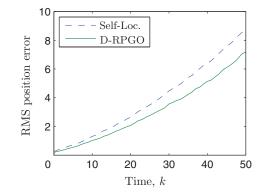


Fig. 11. Simulation: The root-mean-square position error for robot 1 when the distributed algorithm is applied to a group of five robots utilizing inter-robot relative measurements over random paths.



Fig. 12. Two Pioneer P3-DX robots equipped with cameras and targets. Robot 1 is shown on the left, while robot 2 is on the right.

cameras to measure the inter-robot relative pose by exploiting the known geometry of each target. The true pose of each robot was determined using an overhead camera capable of tracking each robot's target. The sensor suite was polled every 0.2 sec with the noisy inter-robot relative pose measurements available at most of, but not all, times.

All robots moved in straight lines with their paths approximately parallel. Six different pose estimates of the robots were obtained at each time. The first was a dead-reckoning estimate, obtained from the inter-time relative pose measurements alone. The remaining five estimates were obtained by using the distributed collaborative localization algorithm with the inter-robot noisy relative measurements being of full pose, orientation only, position only, bearing only, or distance only respectively. These measurements were obtained by projecting the relative pose measurements. The resulting global position estimates, along with the true positions, for robot 1 are reported in Fig. 13. Simulations presented in Section 6 indicate that we should see significant improvement in localization accuracy even in this small team, and the experimental results are consistent with that conclusion. As in the simulations, distinct improvement in localization accuracy is seen when collaborative localization is performed irrespective of the type of inter-robot measurement, although the improvement varied depending on the measurement type.

To show that the improvements seen in the previous experiments are not just a random occurrence, the previous experiment was repeated 24 times to produce an empirical estimate of the bias and standard deviation in position estimation error (by taking appropriate averages) for each type of interrobot measurement. The results are reported in Fig. 14. The experimental results again show that all measurement types lead to an increase in localization accuracy, with relative pose measurements leading to maximum improvement, as expected. This is also the same trend as was observed in the simulations. In contrast to the simulations, here we see that distance measurements do lead to a non-negligible improvement in the bias of localization accuracy. As in the simulations, we see that both bearing and position measurements lead to similar improvement in the bias. However, in contrast to the simulations, in the experiments measurements seem to improve the bias

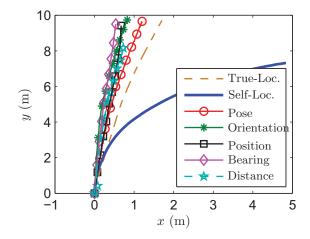


Fig. 13. Experimental: A plot of the location of robot 1 in the overhead camera frame of reference when both robots move in a straight line. The true path (determined using the overhead camera), estimated path using self localization, and estimated path using the distributed collaborative localization algorithm are all reported. The various curves correspond to the type of inter-robot relative measurement used.

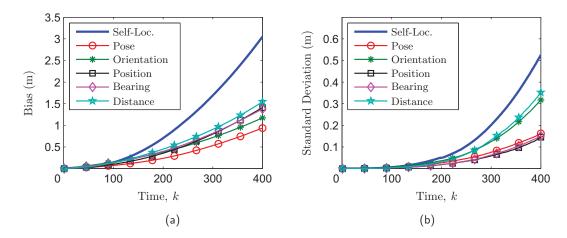


Fig. 14. Experimental: The (a) bias and (b) standard deviation in position error for robot 1 with collaborative localization when both robots move in a straight line. The bias and standard deviation were empirically estimated by averaging over 24 repeated trials.

more than bearing or position measurements. The trend of the standard deviation improvement with orientation, position, and orientation measurements are similar in both simulation and experiments. The most significant difference between the trends observed in simulations and experiments are in the standard deviation improvement between pose, orientation, and position measurements: they are much closer in the experiments than in the simulations. In fact, the standard deviation with position or orientation measurements seem to be a little smaller than that with pose measurements. We believe this is an artifact of the small number of experimental samples averaged to estimate the bias and standard deviation empirically, which limits the accuracy of the estimates.

In short, the experiments verify that the proposed algorithm for distributed collaborative localization leads to statistically significant improvement in localization accuracy even with a small number of robots. Several trends seen in the experiments about the relative merits of the different types of measurements are consistent with those in the simulations. However, there are a few noticeable differences as well. Although the exact cause of these differences is not clear, one should note that the simulations and experiments differ in a number of ways. The biggest difference is that the bias and standard deviations estimated from the experimental samples, which in turn is due to the difficulty associated with conducting repeated experiments. The other significant difference comes from the measurement noise distributions. In the simulations, measurement noise was drawn from distributions

that were somewhat arbitrarily chosen, while the noise distributions in the experiments are unknown. The third potential source of difference is the paths of the robots. In ref. [19], which examined the growth of dead-reckoning error, we showed that the path traversed by a robot plays a crucial role in the bias and standard deviation of the localization error. In particular, the bias grows without bound if the robot moves in a straight line, but stays uniformly bounded by a constant if the robot stays inside a bounded region. Similar effects are likely in collaborative localization. Therefore, the difference between the paths used in the simulation and experiments might be another source of difference in localization accuracy observed.

8. Summary

In this paper, we introduce a novel distributed algorithm (D-RPGO) for estimating the 3D pose of multiple robots when noisy inter-robot measurements of various types (relative pose, orientation, position, bearing, or distance) between pairs of robots are intermittently available. The distributed algorithm is inspired by a centralized algorithm for solving a least square-type problem. The cost function is chosen to measure how well the estimates explain the relative measurements. A gradient-descent in a product Riemannian manifold is used to solve the optimization problem.

The proposed method is close in spirit to the pose-graph optimization problems commonly encountered in mapping and localization. The primary distinction is that our cost function is defined on a Riemannian manifold and is minimized on that surface without converting the problem to a vector space optimization problem. The estimator is thus naturally independent of the choice of parameterization of SO(3) utilized to carry out numerical computation. While Euclidean pose-graph algorithms may require careful choice of design variables for a specific rotation parameterization for it to perform well, the Riemannian pose-graph algorithm does not. In addition, the Riemannian version can achieve better performance in some cases. In particular, the D-RPGO algorithm outperforms a distributed version of the traditional pose-graph optimization (that uses a specific Euclidean parameterization of robot orientations) in terms of accuracy when the noise in the inter-robot measurements is large.

It was found that the computation time of the D-RPGO algorithm is comparable with that of the D-EPGO algorithm in the simulations we performed. However, there are many ways to speed up the computations involved in Euclidean optimization that are not currently available for Riemannian optimizations. We suspect, when applied to large graphs, Riemannian optimization will be slower than Euclidean optimization. Fortunately, the measurement graphs that appear in distributed computation are small, since their size scales with the number of neighbors of a robot.

The IEKF has been a popular tool in past work on collaborative localization as well as mapping. Simulation comparison of the proposed D-RPGO algorithm with the IEKF showed that the IEKF performs poorly compared with the proposed D-RPGO algorithm unless the time interval between successive inter-robot relative measurements is quite small. This has important implications for many practical applications, since it is more likely that inter-robot measurements will be available infrequently. However, in those special cases when inter-robot measurements arrive frequently, the IEKF performs better then the D-RPGO algorithm due to the consideration of the estimation covariance. One such scenario is the one considered in ref. [26], where robots observe common feature points on the ground. In such a scenario the IEKF maybe preferred over the proposed method.

The focus of this paper has been on relative localization with an arbitrary origin or with GPS fixes available only in the beginning. The algorithm is readily applicable to scenarios where global position or orientation measurements may become available over time, such as intermittent GPS or landmarks with known position/orientation detected by robots. For instance, when a landmark with a known position or orientation is detected, a new node is introduced in the graph corresponding to the landmark whose node variable(s) enter the optimization as known values instead of variables to optimize over. The landmarks are effectively grounded nodes.

A limitation of the proposed method over the IEKF (and other filtering methods) is that the latter also provides a covariance estimate, while the proposed method does not. A method to estimate the covariance, or some measure of estimation error, is an important future task. Although simulations reported here provide an indication of when the proposed method performs over the existing stateof-the art methods, a more thorough study needs to be conducted to obtain a better understanding of the relative merits of the proposed method over alternatives. Future studies will also address issues such as identification of the neighboring robots, and rejection of outliers in relative measurements.

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Appendix A: Riemannian Manifolds

In this section we provide greater detail on the building blocks of Algorithm 1, including Riemannian manifolds. A full introduction to the study of Riemannian geometry is outside the scope of this paper; the interested reader is referred to refs. [1, 9].

A manifold *M* of dimension *d* is a topological space that can locally be represented by \mathbb{R}^d . The set of all linear operators on the vector space \mathbb{R}^3 is denoted by $\mathbb{L}(\mathbb{R}^3)$. Elements of $\mathbb{L}(\mathbb{R}^3)$ can be represented by 3×3 matrices. We can then describe the manifold of 3D rotations, denoted as SO(3), by the set $\{R \in \mathbb{L}(\mathbb{R}^3) : R^T R = id, det(R) = 1\}$. The symbol *id* denotes the identity operator. Given a point $p \in SO(3)$, the tangent space of SO(3) at *p*, denoted as $T_pSO(3)$, is given by

$$T_p SO(3) = \left\{ p\hat{v} : \hat{v} \in \mathbb{L}(\mathbb{R}^3), \, \hat{v}^T = -\hat{v} \right\}.$$

A *Riemannian metric* on a manifold *M* is given by defining $\forall p \in M \ g_p : T_pM \times T_pM \to \mathbb{R}$ such that

- $\forall p \in M, g_p$ is an inner product;
- given two vector fields X and Y on M, the map $g: M \to \mathbb{R}, p \mapsto g_p(X_p, Y_p)$ is smooth.

The Riemannian metric g gives rise to an inner-product norm on the tangent space T_pM for each $p \in M$. That is, for $\xi \in T_pM$, $\|\xi\|^2 = g_p(\xi, \xi)$. A *Riemannian manifold* is a manifold equipped with a Riemannian metric. The first Riemannian manifold that we consider is (SO(3), g), where g is the Riemannian metric given by

$$g_p(A, B) = \frac{1}{2} \operatorname{Tr} \left(A^T B \right)$$
(A1)

for all $p \in SO(3)$, $A, B \in T_pSO(3)$. From this point on, when we refer to SO(3), we mean this Riemannian manifold. To further simplify the notation, when the argument $p \in SO(3)$ is clear, we will denote $g_p(X_p, Y_p)$ by $g(X_p, Y_p)$.

For a Riemannian manifold, the analog to a straight line in Euclidean space is given by a *geodesic*. For two points p and q on a manifold M, a geodesic from p to q, denoted by γ_{pq} , is the "shortest" path from p to q. More precisely, if we consider the set of all parameterized paths from p to q, given by $\Gamma = \{\gamma : [0, 1] \rightarrow M : \gamma(0) = p, \gamma(1) = q\}$, then a geodesic γ_{pq} is given by a path that minimizes

$$\int_0^1 \|\gamma'(t)\| dt$$

over all paths in Γ , and $\gamma'(t)$ is the derivative of $\gamma(t)$ with respect to t, which is defined by the following relationship:

$$\gamma'(t_0)f = \frac{d(f(\gamma(t)))}{dt}|_{t_0}$$

for all $f: M \to \mathbb{R}$. Note that the derivative $\gamma'(0)$ is an element of the tangent space $T_{\gamma(0)}M$. For the manifold SO(3), the geodesic from p to $q \in SO(3)$ is given by

$$\gamma_{pq}(t) = p \exp(tv), \ t \in [0, 1],$$

where $v = \log(p^{-1}q)$, the map exp : $\mathbb{L}(\mathbb{R}^3) \to \mathbb{L}(\mathbb{R}^3)$ is given by $\exp(X) = \sum_{k=0}^{\infty} \frac{X^k}{k!}$ and log is the inverse map of exp.¹

For an arbitrary Riemannian manifold, the Riemannian metric defines a corresponding distance function on the manifold as follows. For a Riemannian manifold M, the distance between two points $p, q \in M$ is given by

$$d_M(p,q) = \int_0^1 \|\gamma_{pq}(t)\| dt$$

In particular, for the manifold SO(3),

$$d_{SO(3)}(p,q) = \sqrt{-\frac{1}{2}\operatorname{Tr}\left(\log^2(p^T q)\right)}$$

The subscript SO(3) on the distance function will be omitted when the arguments make it clear what manifold the distance refers to. The second manifold that we consider is $(\mathbb{R}^3, < \cdot, \cdot >)$, standard Euclidean 3-space with the Riemannian metric given by the inner product on \mathbb{R}^3 . Given points $p, q \in \mathbb{R}^3$ the (unique) geodesic connecting p and q is the straight line from p to q and $d(p,q) = ||p-q|| := < p-q, p-q >^{1/2}$, the standard Euclidean inner product norm.

In addition to the distance function, the Riemmanian metric also specifies a parallel transport function. Let M be an arbitrary Riemmanian manifold. For $p \in M$ and $\xi \in T_p M$, the *parallel transport* function, denoted as \exp_p , finds a new point $q \in M$ given by moving along a geodesic beginning at p and with initial velocity ξ . More precisely,

$$q = \exp_p(\xi) = \gamma(1),$$

where γ is the parameterized geodesic with $\gamma(0) = p$ and ξ being the tangent to $\gamma(t)$ at t = 0. In the case of SO(3), $\exp_p(\xi) = p \exp(p^T \xi)$.

Finally, we consider the gradient for real-value functions defined on the manifold. Given a Riemannian manifold M and a scalar, real-valued function $f: M \to \mathbb{R}$, the gradient of f at $p \in M$,

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¹ This use of $\log(p)$ for $p \in SO(3)$ is only valid on the region where exp is a diffeomorphism, which it is at least on the open ball about $I \in SO(3)$ of radius π (Proposition 1.4).¹¹

denoted by grad f(p), can be defined by the following relation:

$$g_p(grad f(p), \xi_p) = (f \circ \gamma)'(0) \quad \forall p \in M,$$
(A2)

where $\gamma : [0, 1] \to M$ is any curve with $\gamma(0) = p$, and $(f \circ \gamma)'(0) := \frac{d(f \circ \gamma)(t)}{dt}|_0$. For an example of a gradient calculation that will be useful later, consider the following function:

$$f_1: SO(3) \to \mathbb{R}, \ p \mapsto \frac{1}{2}d^2(p,q)$$
 (A3)

for some fixed $q \in SO(3)$. For an arbitrary $p \in SO(3)$, we consider the geodesic $\gamma_{pq}(t) = p \exp(tv)$, where $v = log(p^{-1}q)$. It can be shown that

$$(f_1 \circ \gamma)(t) = \frac{1}{2}(1-t)^2 d^2(p,q)$$

and thus

$$(f_1 \circ \gamma)'(0) = -d^2(p,q) = \frac{1}{2} \operatorname{Tr} (v^2).$$

The corresponding tangent vector to $\gamma(t)$ at p is given by $\xi_p := \gamma'(0) = pv$. It follows from (A1) that

$$g_p(grad f_l(p), \xi_p) = \frac{l}{2} \operatorname{Tr} \left((grad f_l(p))^T p v \right).$$

Applying (A2) we have

$$grad f_{I}(p) = -pv = -p\log(p^{-1}q) \quad \forall p \in SO(3).$$
(A4)

For a function $f : SO(3) \to \mathbb{R}$, the gradient at $\mathbf{R} \in SO(3)$ is given by²³:

$$gradf(\mathbf{R}) = f_{\mathbf{R}} - \mathbf{R} f_{\mathbf{R}}^T \mathbf{R},\tag{A5}$$

where $f_{\mathbf{R}} \in L(\mathbb{R}^3)$ is the linear operator whose matrix representation (using the canonical basis vectors for \mathbb{R}^3) is given by: $(f_{\mathbf{R}})_{ij} = \frac{\partial f}{\partial \mathbf{R}_{ij}}$, where \mathbf{R}_{ij} represents the (i, j)th entry of the 3 × 3 matrix representation of **R**. Gradient calculation for a function $\mathbf{f} : \mathbb{R}^3 \to \mathbb{R}$ is straightforward.

Apart from SO(3) and \mathbb{R}^3 , the other Riemannian manifold that is useful in this study is SE(3). Instead of SE(3), however, we consider the equivalent manifold $SO(3) \times \mathbb{R}^3$. There exists a natural way to define the Riemannian metric on this manifold, based on the Riemannian metrics in SO(3) and \mathbb{R}^3 so that we can deal with geodesics and gradients. This has to do with the fact that $SO(3) \times \mathbb{R}^3$ is a product manifold, which is the topic of the next section.

Appendix B: Product Manifold

Let $\{(M_i, g^{(i)})_{i=1}^n\}$ be a set of *n* Riemannian manifolds. We define the *product Riemannian manifold* (M, g) as follows:

$$M := M_1 \times \dots \times M_n,$$

$$g_p(\xi, \zeta) := \sum_{i=1}^n g_{p_i}^{(i)}(\xi_i, \zeta_i)$$
(B1)

for all $p = (p_1, ..., p_n) \in M$ and all $\xi = (\xi_1, ..., \xi_n)$, $\zeta = (\zeta_1, ..., \zeta_n) \in T_p M$. We first restrict our considerations to product Riemannian manifolds of the form $M = M_1 \times M_2$ (i.e. n = 2). The

extension to any finite combination is straightforward. The following lemma will be useful in the sequel. The proof of the lemma can be found in ref. [23].

Lemma 1. $[^{23}]$ Let $M = M_1 \times M_2$ be a product Riemannian manifold as defined in (B1) and consider a smooth function $f : M \to \mathbb{R}$. Then for any parameterized path $\gamma = (\gamma_1, \gamma_2) : [0, 1] \to M$,

$$\frac{d}{dt}f(\gamma(t))|_{t_0} = \frac{d}{dt}f(\gamma_1(t), \gamma_2(t_0))\Big|_{t_0} + \frac{d}{dt}f(\gamma_1(t_0), \gamma_2(t))\Big|_{t_0}$$

where $t_0 \in [0.1]$ *.*

Theorem 1. Let $M = M_1 \times M_2$ be a product Riemannian manifold as defined in (B1). Consider a smooth function $f : M \to \mathbb{R}$ and for all $(p, q) \in M$ $(p \in M_1, q \in M_2)$, define

$$f_1^q : M_1 \to \mathbb{R}, \ a \mapsto f(a,q),$$

 $f_2^p : M_2 \to \mathbb{R}, \ b \mapsto f(p,b).$

Then for all $(p,q) \in M$

$$grad f(p,q) = \left(grad f_1^q(p), grad f_2^p(q)\right)$$

Proof. Fix $(p, q) \in M$ and consider an arbitrary tangent vector $\xi = (\xi_1, \xi_2) \in T_{(p,q)}M$. Choose a parameterized path $(\gamma_1, \gamma_2) = \gamma : [0, 1] \to M$ such that ξ_1 is tangent to $\gamma_1(t)$ at $p = \gamma_1(t_0)$ and ξ_2 is tangent to $\gamma_2(t)$ at $q = \gamma_2(t_0)$ (and thus ξ is tangent to $\gamma(t)$ at $(p, q) = \gamma(t_0)$). Using Lemma 1, we have

$$(f \circ \gamma)'(t_0) = \frac{d}{dt}(f_1^q \circ \gamma_1)|_{t_0} + \frac{d}{dt}(f_2^p \circ \gamma_2)|_{t_0}.$$

From this relation and (A2), we find that

$$g(grad f(p, q), \xi) = g^{(1)}(grad f_1^q(p), \xi_1) + g^{(2)}(grad f_2^p(q), \xi_2).$$
(B2)

Let $grad f(p, q) = (A, B) \in T_{(p,q)M}$ for some unknown $A \in T_pM$, $B \in T_qM$. Using the definition of g given in (B1), we can rewrite (B2) as

$$g^{(1)}(A,\xi_1) + g^{(2)}(B,\xi_2) = g^{(1)}(gradf_1^q(p),\xi_1) + g^{(2)}(gradf_2^p(q),\xi_2).$$

Since this equality holds for all $\xi_1 \in T_p M$ and all $\xi_2 \in T_q M$, we have $A = grad f_1^q(p)$, $B = grad f_2^p(q)$, which completes the proof.

As the number of manifolds increases, the notation $grad f_{I}^{q}(p)$ becomes more cumbersome. We will therefore simplify the notation by writing grad f(p) (to mean $grad f_{I}^{q}(p)$) whenever the manifold the gradient is found with respect to is clear from the choice of p and q.

The following corollary is a direct consequence of Theorem 1.

Corollary 1. Given a set of n Riemannian manifolds $\{(M_i, g_i)\}_{i=1}^n$, if the product Riemannian metric on the product manifold (M, g), where $M = M_1 \times \cdots \times M_n$, is defined as

$$g_p(X, Y) = g_1(X_1, Y_1) + \dots + g_n(X_n, Y_n)$$

for $p = (p_1, ..., p_n) \in M$ and $X = (X_1, ..., X_n), Y \in T_p M = T_{p_1} M_1 \times ... T_{p_n} M_n$, then

$$grad f(p) = (grad f(p_1), \ldots, grad f(p_n))$$
.

We next consider what geodesics look like on the product manifold.

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Theorem 2. If *M* is a product Riemannian manifold as defined in (B1), then a geodesic between $p = (p_1, p_2)$ and $q = (q_1, q_2) \in M$ can be obtained by the product of geodesics between p_1 and q_1 , and p_2 and q_2 , on M_1 and M_2 respectively.

Proof. Let γ_{pq} , $\gamma_{p_1q_1}$, $\gamma_{p_2q_2}$ denote the geodesics on M, M_1 , and M_2 respectively. By definition, $\gamma_{pq} =: (\gamma^{(1)}, \gamma^{(2)})$ is the path that minimizes

$$\gamma_{pq} = \operatorname*{argmin}_{\gamma \in \Gamma_{pq}} \int_0^1 \|\gamma'(t)\|^2 dt,$$

where Γ_{pq} denotes the set of all paths on M from p to q, and for $\gamma \in \Gamma_{pg}$, $\gamma(t)$ denotes a parameterization of path γ such that $\gamma(0) = p$ and $\gamma(1) = q$. Using the definition of the innerproduct norm on the product manifold, we see that

$$\begin{split} \gamma_{pq} &= \underset{(\gamma^{(1)},\gamma^{(2)})\in\Gamma_{pq}}{\arg\min} \int_{0}^{1} (\|\gamma^{(1)'}(t)\|^{2} + \|\gamma^{(2)'}(t)\|^{2}) dt \\ &= \left(\underset{\gamma\in\Gamma_{p_{1}q_{1}}}{\arg\min} \int_{0}^{1} \|\gamma'(t)\|^{2} dt, \ \underset{\gamma\in\Gamma_{p_{2}q_{2}}}{\arg\min} \int_{0}^{1} \|\gamma'(t)\|^{2} dt \right) \\ &= (\gamma_{p_{1}q_{1}},\gamma_{p_{2}q_{2}}). \end{split}$$

That parallel transport on M is given by parallel transport on each individual M_i immediately follows. Thus, proving (6).

Appendix C Proof of Theorem 1

Proof. The cost function (3) is a function from \mathcal{M} to \mathbb{R} , where \mathcal{M} is the product manifold

$$\mathcal{M} := (SO(3) \times \mathbb{R}^3)^n$$
, where $n := |\mathcal{V}(k)|$.

We define the Riemannian metric on this product manifold as follows. For $p = (\mathbf{R}_1, \mathbf{t}_1, \dots, \mathbf{R}_n, \mathbf{t}_n) \in \mathcal{M}$ and $X, Y \in T_p \mathcal{M}$, where $X \in T_p \mathcal{M}$ is expanded as $(X_{\mathbf{R}_1}, X_{\mathbf{t}_1}, \dots, X_{\mathbf{R}_n}, X_{\mathbf{t}_n})$,

$$g_p(X, Y) := \sum_{i \in n} \left(g_{\mathbf{R}_i}(X_{\mathbf{R}_i}, Y_{\mathbf{R}_i}) + \langle X_{\mathbf{t}_i}, Y_{\mathbf{t}_i} \rangle \right).$$

This meets the conditions of Corollary 1 and so for $p = (\mathbf{R}_1, \mathbf{t}_1, \dots, \mathbf{R}_n, \mathbf{t}_n) \in \mathcal{M}$, we have

$$grad f(p) = (grad f(\mathbf{R}_{l}), grad f(\mathbf{t}_{l}), \dots, grad f(\mathbf{R}_{h}), grad f(\mathbf{t}_{h}))$$

All that remains is to determine $grad f(\mathbf{R}_u)$ and $grad f(\mathbf{t}_u)$ for u = 1, ..., n. Using linearity of the gradient operator, we have

grad
$$f(\mathbf{R}_u) = \sum_{e \in \mathcal{I}(k)} grad c_e(\mathbf{R}_u)$$
, and $grad f(\mathbf{t}_u) = \sum_{e \in \mathcal{I}(k)} grad c_e(\mathbf{t}_u)$.

First define $f_2 := \frac{1}{2}d^2(\mathbf{R}_u^T\mathbf{R}_v, \hat{\mathbf{R}}_{u\,v}), f_3 := \frac{1}{2}\|\hat{\mathbf{t}}_{u\,v} - \mathbf{R}_u^T(\mathbf{t}_v - \mathbf{t}_u)\|^2$, so that for $e \triangleright e(u, v), c_e = f_2 + f_3$. For h = u or h = v, we therefore have

grad
$$c_e(\mathbf{R}_h) = grad f_2(\mathbf{R}_h) + grad f_3(\mathbf{R}_h),$$
 (C1)

$$grad \ c_e(\mathbf{t}_h) = 0 + grad \ f_3(\mathbf{t}_h). \tag{C2}$$

The Riemannian distance d(p,q) is bi-invariant, meaning that for arbitrary $R, \bar{R} \in SO(3)$, $d(p,q) = d(Rp\bar{R}, Rq\bar{R})$. Using bi-invariance, we obtain $d^2(\mathbf{R}_u^T \mathbf{R}_v, \hat{\mathbf{R}}_{uv}) = d^2(\mathbf{R}_v, \mathbf{R}_u \hat{\mathbf{R}}_{uv})$, which implies grad $f_2(\mathbf{R}_v) = grad f_1(\mathbf{R}_v|q)$, where $f_1 : SO(3) \to \mathbb{R}$ is given by (A3) and $q = \mathbf{R}_u \hat{\mathbf{R}}_{uv}$. Applying formula (A4), we obtain

grad
$$f_2(\mathbf{R}_v) = -\mathbf{R}_v \log(\mathbf{R}_v^T \mathbf{R}_u \hat{\mathbf{R}}_{uv}).$$

The expression for grad $f_2(\mathbf{R}_u)$ is similarly obtained by thinking of f_2 as a function of \mathbf{R}_u and keeping \mathbf{R}_v , $\hat{\mathbf{R}}_{uv}$ fixed. These gradients are compactly represented as:

$$grad f_2(\mathbf{R}_h) = \begin{cases} -\mathbf{R}_h \log \left(\mathbf{R}_h^T \, \mathbf{R}_v \, \hat{\mathbf{R}}_{u \, v}^T \right) & \text{if } h = u \\ -\mathbf{R}_h \log \left(\mathbf{R}_h^T \, \mathbf{R}_u \, \hat{\mathbf{R}}_{u \, v} \right) & \text{if } h = v \\ 0 & \text{otherwise} \end{cases}$$

The gradient of the function $f_3 : SO(3) \to \mathbb{R}$ (thinking of f_3 only as a function of \mathbf{R}_u while $\mathbf{t}_u, \mathbf{t}_v, \hat{\mathbf{t}}_{uv}$ are fixed) can be obtained using (A5) with straightforward but tedious calculations; which turn out to be $gradf_3(\mathbf{R}_u) = -(\mathbf{t}_v - \mathbf{t}_u)\hat{\mathbf{t}}_{uv}^T + \mathbf{R}_u\hat{\mathbf{t}}_{uv}(\mathbf{t}_v - \mathbf{t}_u)^T\mathbf{R}_u$. The gradient of $f_3 : \mathbb{R}^3 \to \mathbb{R}$ is simply $(\frac{\partial f_3(x)}{\partial x})^T$, which is

$$grad f_{3}(\mathbf{t}_{h}) = \begin{cases} \left(\mathbf{t}_{h} + \mathbf{R}_{h} \, \hat{\mathbf{t}}_{u \, v} - \mathbf{t}_{v}\right) & \text{if } h = u \\ \left(\mathbf{t}_{h} - \mathbf{R}_{u} \, \hat{\mathbf{t}}_{u \, v} - \mathbf{t}_{u}\right) & \text{if } h = v \\ 0 & \text{o.w.} \end{cases}$$

These formulae completely specify the gradient of the edge cost, $grad c_e(\cdot)$ in (C1). The expressions for $grad f(\mathbf{R}_u)$ and $grad f(\mathbf{t}_u)$ for u = 1, ..., n that are provided in the theorem are obtained simply by adding the components of the gradients that are derived above.

Appendix D: Gradient of the Cost Function (3) for Heterogeneous Measurements

As in the proof of Proposition 1, the gradient of function f in (3) at a point $p = (\mathbf{R}_1, \mathbf{t}_1, \dots, \mathbf{R}_n, \mathbf{t}_n) \in (SO(3) \times \mathbb{R}^3)^n$ is given by

$$gradf(p) = \sum_{e \in \mathcal{E}(k)} grad c_e(p) =: (gradf(\mathbf{R}_1), \dots, gradf(\mathbf{t}_n)),$$
(D1)

where $grad g_e(p)$ is the gradient of the edge cost function for edge e = (u, n). Finding the gradient of the cost function (3) then reduces to finding the gradients of the edge costs c_e (specified in (4)) for each edge $e \in \mathcal{E}(k)$.

It follows from Corollary 1 that the gradient of the cost function c_e in (4) for the edge $e = (u, v) \in \mathcal{E}(k)$ is

grad
$$c_e(p) = (grad \ c_e(\mathbf{R}_1), grad \ c_e(\mathbf{t}_1), \dots, grad \ c_e(\mathbf{R}_n), grad \ c_e(\mathbf{t}_n)).$$

For $\ell(k)(e) = \mathbf{T}$ (pose), the gradients have already been computed in the previous section, which can be compactly represented as $(e \triangleright (u, v))$

$$grad c_{e}(\mathbf{R}_{h}) = \begin{cases} -2\mathbf{R}_{h} \Big(\log(\mathbf{R}_{h}^{T}\mathbf{R}_{v}\hat{\mathbf{R}}_{uv}^{T}) \\ +\mathbf{R}_{h}^{T}(\mathbf{t}_{v}-\mathbf{t}_{u})\hat{\mathbf{t}}_{uv}^{T} - \hat{\mathbf{t}}_{uv}(\mathbf{t}_{v}-\mathbf{t}_{u})^{T}\mathbf{R}_{h} \Big)^{\text{if } h = u} \\ -2\mathbf{R}_{h} \log(\mathbf{R}_{h}^{T}\mathbf{R}_{u}\hat{\mathbf{R}}_{uv}) & \text{if } h = v \\ 0 & \text{o.w.} \end{cases}$$

grad $c_e(\mathbf{t}_h) = 2I_{uv}(h)(\mathbf{t}_u + \mathbf{R}_u \hat{\mathbf{t}}_{uv} - \mathbf{t}_j),$

where $I_{uv}(h) = 1$ if h = u, -1 if h = v, and 0 otherwise. If $\ell(k)(e) = \mathbf{R}$ (orientation) or $\ell(k)(e) = \mathbf{t}$ (position), we have the following expressions for the gradient from the previous section. If $\ell(k)(e) = \mathbf{R}$,

$$grad c_e(\mathbf{R}_h) = \begin{cases} -2\mathbf{R}_h \Big(\log(\mathbf{R}_h^T \mathbf{R}_v \hat{\mathbf{R}}_{uv}^T) & \text{if } h = u \\ -2\mathbf{R}_h \Big(\log(\mathbf{R}_h^T \mathbf{R}_u \hat{\mathbf{R}}_{uv}) & \text{if } h = v \\ 0 & \text{otherwise} \end{cases},\\ grad c_e(\mathbf{t}_h) = 0, \end{cases}$$

and if $\ell(k)(e) = \mathbf{t}$, then

$$grad c_e(\mathbf{R}_h) = \begin{cases} -2\mathbf{R}_h \Big(\mathbf{R}_h^T (\mathbf{t}_v - \mathbf{t}_u) \hat{\mathbf{t}}_{uv}^T - \hat{\mathbf{t}}_{uv} (\mathbf{t}_v - \mathbf{t}_u)^T \mathbf{R}_h \Big) & \text{if } h = u \\ 0 & \text{otherwise} \end{cases},\\ grad c_e(\mathbf{t}_h) = 2I_{uv}(h)(\mathbf{t}_u + \mathbf{R}_u \hat{\mathbf{t}}_{uv} - \mathbf{t}_v). \end{cases}$$

If $\ell(k)(e) = \tau$ (bearing) or $\ell(k)(e) = \delta$ (distance), the gradient grad $c_e(\mathbf{R}_h)$ can be computed by using formula (A5). The gradient grad $c_e(\mathbf{t}_h)$ is obtained by differentiation. We obtain, for $\ell(k)(e) = \tau$ (bearing)

grad
$$c_e(\mathbf{R}_h) = \begin{cases} -2\mathbf{R}_h \Big(\mathbf{R}_h^T (\mathbf{t}_v - \mathbf{t}_h) \hat{\boldsymbol{\tau}}_{uv}^T \| \mathbf{t}_u - \mathbf{t}_v \| - \hat{\boldsymbol{\tau}}_{uv} \| \mathbf{t}_v - \mathbf{t}_u \| (\mathbf{t}_v - \mathbf{t}_u)^T \mathbf{R}_h \Big) & \text{if } h = u \\ 0 & \text{otherwise} \end{cases},$$

grad $c_e(\mathbf{t}_h) = -4I_{uv}(h)[(\mathbf{t}_v - \mathbf{t}_u) - \|\mathbf{t}_v - \mathbf{t}_u\|\mathbf{R}_u\hat{\boldsymbol{\tau}}_{uv}],$

and if $\ell(k)(e) = \delta$ (dist),

$$grad c_e(\mathbf{R}_h) = 0,$$

$$grad c_e(\mathbf{t}_h) = -2I_{uv}(h) \frac{(\hat{\delta}_{uv} - \|\mathbf{t}_v - \mathbf{t}_u\|)}{\|\mathbf{t}_v - \mathbf{t}_u\|} (\mathbf{t}_v - \mathbf{t}_u).$$

The gradient grad f(p) can now be computed by using these formulas.