

# Synchronization Problems in Computer Vision with Closed-Form Solutions

Federica Arrigoni<sup>1,2</sup> · Andrea Fusiello<sup>1</sup>

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#### Abstract

In this paper we survey and put in a common framework several works that have been developed in different contexts, all dealing with the same abstract problem, called *synchronization* by some authors, or averaging, or graph optimization by others. The problem consists in recovering some variables from a set of pairwise relation measurements. In particular, we concentrate on instances where the variables and the measures belong to a (*semi-)group* and the measures are their mutual differences (or ratios, depending on how the group operation is called). The groups we deal with have a matrix representation, which leads to an elegant theory and closed-form solutions.

**Keywords** Synchronization  $\cdot$  Averaging  $\cdot$  Graph optimization  $\cdot$  Multiple point-set registration  $\cdot$  Structure from motion  $\cdot$  Multi-view matching

# 1 Introduction

Consider a network of nodes where each node is characterized by an unknown state, and suppose that pairs of nodes can measure the ratio (or difference) between their states. The goal is to infer the unknown states from the pairwise measures. This is a general statement of the *synchronization* problem (Karp et al. 2003; Giridhar and Kumar 2006; Singer 2011). States are represented by elements of a group  $\Sigma$ , that is why the problem is actually referred to as *group synchronization*. The groups of interest are defined in Table 1, while Fig. 1 represents their inclusion.

The problem can be usefully modelled by introducing a graph  $\mathcal{G} = (\mathcal{V}, \mathcal{E})$ , which is referred to as the *measurement graph*, where nodes correspond to the unknown states and edges correspond to the pairwise measures, and it is well-

 Andrea Fusiello andrea.fusiello@uniud.it
 Federica Arrigoni

arrigoni.federica@spes.uniud.it

<sup>1</sup> DPIA, University of Udine, Via Delle Scienze 208, Udine, Italy

<sup>2</sup> Czech Institute of Informatics, Robotics and Cybernetics, Czech Technical University in Prague, Jugoslavskych partyzanu 3, Prague, Czech Republic posed only if such a graph is connected. Synchronization can be seen as upgrading from relative (pairwise) information, which involves two nodes at a time, onto absolute (global) information, which involves all the nodes simultaneously. In the literature, the same problem is also referred to as *averaging* (Govindu 2004; Wilson et al. 2016; Hartley et al. 2013) or *graph optimization* (Carlone et al. 2015) in some cases.

As an example, consider the graph in Fig. 2, where nodes and edges are labelled with integer numbers: the task is to recover the unknown numbers in the nodes by measuring their differences (on the edges). Two things can be immediately observed: a solution exists only if the sum of the differences along any cycle is zero, and, when it exists, the solution is not unique, for adding a constant to the nodes does not change the differences.

Measures are typically corrupted by errors, which can be gross errors (outliers) and/or a diffuse noise with small variance. If  $\mathcal{G}$  is a tree then these errors will creep in the solution, however, as soon as redundant measures are considered (i.e. the graph has at least one cycle), they are exploited by synchronization to globally compensate the errors. The solution minimizes a suitable cost function which evaluates the coherence between the unknown states and the pairwise measures. By construction synchronization enforces *cycle consistency* (Enqvist et al. 2011; Zach et al. 2010), namely the property that the composition of relative measures along any cycle in the measurement graph should return the identity.

Table 1         Matrix groups/semigroups considered in this paper				
GL(d)	The General Linear Group is the set of invertible matrices: $GL(d) = \{M \in \mathbb{R}^{d \times d} \text{ s.t. } \det(M) \neq 0\}$			
SL(d)	The Special Linear Group is the set of matrices with unit determinant: $SL(d) = \{M \in \mathbb{R}^{d \times d} \text{ s.t. } det(M) = 1\}$			
O(d)	The Orthogonal Group is the set of rotations and reflections: $O(d) = \{M \in \mathbb{R}^{d \times d} s.t. M^{T}M = MM^{T} = I_d\}$			
SO(d)	The Special Orthogonal Group is the set of rotations: $SO(d) = \{M \in O(d) \ s.t. \ \det(M) = 1\}$			
GA(d)	The General Affine Group is the set of affine maps: $GA(d) = \left\{ \begin{bmatrix} M & \mathbf{t} \\ 0^{T} & 1 \end{bmatrix}, s.t. \ M \in GL(d), \mathbf{t} \in \mathbb{R}^d \right\}$			
SE(d)	The Special Euclidean Group is the set of direct isometries: $SE(d) = \left\{ \begin{bmatrix} M & \mathbf{t} \\ 0^{T} & 1 \end{bmatrix}, s.t. M \in SO(d), \mathbf{t} \in \mathbb{R}^d \right\}$			
$\mathcal{S}_d$	The Symmetric Group is the set of total permutations: $S_d = \{M \in \{0, 1\}^{d \times d} \text{ s.t. } M1 = 1, 1M = 1\}$			
$\mathcal{I}_d$	The Symmetric Inverse Semigroup is the set of partial permutations: $\mathcal{I}_d = \{M \in \{0, 1\}^{d \times d} \text{ s.t. } M1 \leq 1, \ 1M \leq 1\}$			



**Fig. 1** Subgroups of GL(d). The identity is in the intersection of all of them



**Fig. 2** Synchronization over  $(\mathbb{Z}, +)$ 

Several instances of synchronization have been studied in the literature.  $\Sigma = \mathbb{R}$  yields *clock synchronization* (with offset) (Karp et al. 2003; Giridhar and Kumar 2006), from which the term *synchronization* originates, where all the nodes in a network are synchronized to a common clock. The same problem is studied under the name *levelling* in topography and surveying (Bjerhammar 1973). If the clock model includes also a drift (besides the offset), the problem can be addressed in  $\Sigma = GA(1)$  (Solis et al. 2006). Synchronization over GA(3) has also been applied to RGB colour matching in image mosaicking (Santellani et al. 2018).  $\Sigma = \mathbb{Z}_2$  gives sign synchronization (Cucuringu 2015), which was used to identify communities in a network where the interaction between the nodes is described by two dichotomous values, e.g., agreement/disagreement, and the network has a natural partition into two communities.  $\Sigma = \mathbb{R}^d$  is a *translation* synchronization (Barooah and Hespanha 2007; Russel et al. 2011; Tron et al. 2015; Molavi and Jadbabaie 2011; Aragues et al. 2012), namely the problem of localizing a set of nodes in space from pairwise differences.  $\Sigma = SO(d)$  corresponds to rotation synchronization (also known as rotation averaging) (Sharp et al. 2002; Martinec and Pajdla 2007; Singer 2011; Thunberg et al. 2011; Crandall et al. 2011; Fredriksson and Olsson 2012; Boumal et al. 2013; Hartley et al. 2013; Chatterjee and Govindu 2013; Wang and Singer 2013; Arrigoni et al. 2014a; Tron et al. 2016) and  $\Sigma = SE(d)$  results in rigid-motion synchronization (also known as motion averaging or pose graph optimization) (Fusiello et al. 2002; Govindu 2004; Torsello et al. 2011; Tron and Danilidis 2014; Bernard et al. 2015; Arrigoni et al. 2016c, b; Rosen et al. 2015, 2019), which find application in structure from motion, registration of 3D point sets and simultaneous localization and mapping (SLAM).  $\Sigma = SL(d)$  produces homography synchronization (Schroeder et al. 2011), which is an essential step in the context of image stitching or mosaicking. Finally,  $\Sigma = S_d$  and  $\Sigma = I_d$  give rise to *permutation synchroniza*tion (Pachauri et al. 2013; Shen et al. 2016; Yu et al. 2016) and partial permutation synchronization (Arrigoni et al. 2017), respectively, which are related to multi-view matching, as it will be clarified in Sect. 8.2.

While it is clear that synchronization over  $\mathbb{R}^d$  can be reduced to a linear system of equations, the other instances of synchronization involve the minimization of a non convex cost function, which make the problem difficult to solve. However, if  $\Sigma$  admits a matrix representation, i.e. it can be embedded in  $\mathbb{R}^{d \times d}$ , then synchronization reduces to an eigenvalue decomposition, resulting in an efficient and closed-form solution. Specifically, the unknown states are derived from the top eigenvectors of a matrix constructed from the pairwise measures. An equivalent null-space formulation can also be derived. This procedure was introduced in Singer (2011) for  $\Sigma = SO(2)$ , extended in Arie-Nachimson et al. (2012), Singer and Shkolnisky (2011) to  $\Sigma = SO(3)$ , and further generalized in Arrigoni et al. (2016c), Bernard et al. (2015) to  $\Sigma = SE(3)$ . The same formulation appeared in Schroeder et al. (2011) and Pachauri et al. (2013) for  $\Sigma = SL(d)$  and  $\Sigma = S_d$  respectively. The latter was extended to  $\Sigma = \mathcal{I}_d$  in Arrigoni et al. (2017).

Other techniques exploiting a matrix representation of the group can be found in the literature, which express the synchronization problem in terms of well studied mathematical tools, such as semidefinite programming (Singer 2011; Wang and Singer 2013; Rosen et al. 2015) or matrix completion (Arrigoni et al. 2018), resulting in iterative solutions. These approaches, however, are not applicable to all groups. For example, the semidefinite programming formulation derives from specific properties of O(d) and SE(d), namely the fact that the matrix constructed from the pairwise measures is positive semidefinite when considering the Orthogonal Group (Singer 2011; Wang and Singer 2013), and the property that the convex hull of the Special Euclidean Group admits a semidefinite representation (Rosen et al. 2015). The matrix completion formulation (Arrigoni et al. 2018), instead, is based on the property that the matrix containing the pairwise measures is low-rank and it is incoherent (see Candès and Tao 2010). Unfortunately in the  $\Sigma = S_d$  case the matrix is sparse, being composed of permutation matrices, and hence it does not satisfy the incoherence assumption.

Other approaches include iterating local solutions on the measurement graph or explicit minimization of a cost function. The former comprises (Sharp et al. 2002), where the error is distributed over a set of cycles, and (Torsello et al. 2011; Aftab et al. 2015), where each state is updated in turn in a distributed fashion. The latter includes Quasi-Newton iterations (Fusiello et al. 2002), the Levenberg-Marquardt algorithm (Crandall et al. 2011), Riemannian trust-region optimization (Boumal et al. 2013), Riemannian gradient descent (Tron and Danilidis 2014), integer quadratic programming (Carlone and Censi 2014), and the Gauss-Seidel method (Yu et al. 2016). These techniques, however, heavily depend on the chosen group and its parametrization. For instance, unit quaternions are used in Castellani et al. (2002) and dual quaternions in Torsello et al. (2011), which represent rotations and rigid-motions in 3-space, respectively. See also the references in Carlone and Censi (2014).

#### 1.1 Scope and Outline

In this paper we provide a comprehensive survey on closedform solutions to group synchronization, namely linear least squares (Barooah and Hespanha 2007; Russel et al. 2011) and spectral decomposition (Singer 2011; Schroeder et al. 2011; Singer and Shkolnisky 2011; Arie-Nachimson et al. 2012; Pachauri et al. 2013; Bernard et al. 2015; Arrigoni et al. 2016c, 2017). The former solves synchronization over  $\mathbb{R}^d$  in an optimal way. The latter, although suboptimal, is theoretically appealing since it can be applied to any group admitting a matrix representation (e.g. homographies, rigid motions, permutations, ...), as opposed to other techniques which are based on ad-hoc minimizations of specific cost functions. These solutions are extremely fast and they easily cope with weights on individual relative measures, allowing a robust extension via Iteratively Reweighted Least Squares (IRLS) (Holland and Welsch 1977).

We also set forth a theoretical unified framework where several synchronization problems are seen as instances of a more abstract principle, gathering several works that were developed in different communities (including Computer Vision, Photogrammetry, Robotics and Graph Theory), and we show how this framework can be profitably used in several applications.

The paper is organized as follows. Section 2 formally defines the synchronization problem, which is grounded on the notion of group-labelled graph. Section 3 is devoted to synchronization over  $(\mathbb{R}^d, +)$ , which is expressed as a linear system of equations. Section 4 addresses synchronization over  $(GL(d), \cdot)$ , which can be cast to a spectral decomposition or a null-space problem. Then, several subgroups of GL(d) are analysed in Sect. 5, namely SL(d), O(d), SE(d-1), GA(d-1) and  $S_d$ , in which cases the solution needs to be projected onto the group, as closure is not guaranteed. Section 6 shows that the spectral solution can be extended to synchronization over  $\mathcal{I}_d$ , which is an inverse monoid and a subsemigroup of  $S_d$ . Some considerations about the optimization problem associated with the spectral solution are reported in Sect. 7. Finally, Sect. 8 describes how some Computer Vision problems can be addressed in terms of synchronization and Sect. 9 briefly discusses pros and cons of spectral synchronization. The results presented in this paper require some basic notions from graph theory, which are covered in "Appendix B", and the definitions of the Kronecker, Hadamard and Khatri-Rao products, which are given in "Appendix A".

# 2 Theoretical Framework

Let us start by introducing the notion of *group-labelled graph* (Edelman and Saks 1979). Let  $(\Sigma, *)$  be a group with unit

element  $1_{\Sigma}$ , and let  $\mathcal{G} = (\mathcal{V}, \mathcal{E})$  be a simple directed graph with vertex set  $\mathcal{V} = \{1, 2, ..., n\}$  and edge set  $\mathcal{E}$ , with  $m = |\mathcal{E}|$ . A  $\Sigma$ -*labelled* graph is a directed graph with a labelling of its edge set by elements of  $\Sigma$ , that is a tuple  $\Gamma = (\mathcal{V}, \mathcal{E}, z)$  where

$$z: \mathcal{E} \to \Sigma \tag{1}$$

is such that if  $(i, j) \in \mathcal{E}$  then  $(j, i) \in \mathcal{E}$  and

$$z(j,i) = z(i,j)^{-1}.$$
 (2)

Thus, we may also view  $\mathcal{G}$  as an undirected graph.

**Definition 1** Let  $\Gamma = (\mathcal{V}, \mathcal{E}, z)$  be a  $\Sigma$ -labelled graph. We say that a circuit  $\{(i_1, i_2), (i_2, i_3), \dots, (i_{\ell}, i_1)\}$  is a *null cycle*<sup>1</sup> if and only if the composition of the edge labels along the circuit returns the identity, namely

$$z(i_1, i_2) * z(i_2, i_3) * \dots * z(i_\ell, i_1) = 1_{\Sigma}.$$
 (3)

The "null" term clearly refers to an additive notation for the group (which is used in Kawarabayashia and Wollan 2006; Chudnovsky et al. 2008; Guillemot 2011), without implying that  $\Sigma$  needs to be Abelian. Note that if the group is not commutative, then it may happen that cyclic shifts of the same circuit yield different elements of the group. Nevertheless the notion of null cycle is well defined, as either all of the cyclic shifts are equal to  $1_{\Sigma}$  or none of them, as observed also in Guillemot (2011). As noted in Barooah and Hespanha (2008), the concept of null cycle resembles Kirchoff's voltage law, stating that the electrical potential differences around any cycle sum to zero.

**Definition 2** Let  $\Gamma = (\mathcal{V}, \mathcal{E}, z)$  be a  $\Sigma$ -labelled graph. Let  $x : \mathcal{V} \to \Sigma$  be a vertex labelling. We say that *x* is a *consistent labelling* if and only if

$$z(e) = x(i) * x(j)^{-1} \quad \forall \ e = (i, j) \in \mathcal{E}.$$
 (4)

A  $\Sigma$ -labelled graph admitting a consistent labelling is also called *balanced* (Joglekar et al. 2012). Equation (4) means that each edge label is the ratio of the corresponding vertex labels, as shown in Fig. 3. Such condition is referred to as *consistency constraint* and it is equivalent to

$$z(e) * x(j) = x(i) \quad \forall e = (i, j) \in \mathcal{E}.$$
(5)

It is understood that a consistent labelling is defined up to a global (right) product with any group element, in the sense that if  $x : \mathcal{V} \to \Sigma$  is consistent then also  $y : \mathcal{V} \to \Sigma$ , y(i) = x(i) \* s, is consistent, for any (fixed)  $s \in \Sigma$ .



Fig. 3 Consistent labelling

The notion of consistent labelling is strictly related to that of null cycle, as stated by the following result.

**Proposition 1** (Guillemot 2011) Let  $\Gamma = (\mathcal{V}, \mathcal{E}, z)$  be a  $\Sigma$ labelled graph. There exists a polynomial algorithm which either finds a non-null cycle in  $\Gamma$  or finds a consistent labelling of  $\Gamma$ .

The following procedure draws an outline of the proof. First, compute a spanning tree (see "Appendix B") and use Eq. (5) to label nodes, starting from the root labelled with the identity  $1_{\Sigma}$ : this is a consistent labelling by construction. Then add one by one the edges not belonging to the spanning tree, thereby creating a circuit. If the cycle is null then the edge can be added and leave the labelling consistent, otherwise a non-null cycle has been found.

**Corollary 1** (Guillemot 2011) The  $\Sigma$ -labelled graph  $\Gamma = (\mathcal{V}, \mathcal{E}, z)$  has a consistent labelling if and only if it does not contain a non-null cycle.

One direction is an immediate consequence of Proposition 1. The opposite direction follows from the observation that, if Eq. (4) holds, then all the terms in the left side in Eq. (3) simplify, yielding the identity.

## 2.1 Group Feedback Edge Set

The problem of finding non-null cycles in a group-labelled graph is studied in Graph Theory community under the name of "group feedback edge set" problem (Guillemot 2011). Specifically, the goal is to break non-null cycles by deleting *k* edges, where  $k \in \mathbb{N}$  is assumed to be known.

**Definition 3** Let  $\Gamma = (\mathcal{V}, \mathcal{E}, z)$  be a  $\Sigma$ -labelled graph. The *Group Feedback Edge Set* (GFES) problem is defined as follows: on input  $(\Gamma, k)$  for some  $k \in \mathbb{N}$ , decide whether there

<sup>&</sup>lt;sup>1</sup> A circuit is also a cycle; definitions are given in "Appendix B".

exists a subset of the edges  $S \subseteq \mathcal{E}$  with  $|S| \leq k$  such that the labelled graph of the remaining edges  $\Gamma' = (\mathcal{V}, \mathcal{E} \setminus S, z)$ does not contain a non-null cycle.

With some abuse of notation, in Definition 3 we denote with  $(\mathcal{V}, \mathcal{E} \setminus \mathcal{S}, z)$  the  $\Sigma$ -labelled graph with edges in  $\mathcal{S}$ removed from  $\mathcal{E}$ , even though formally z has in its domain edges that do not exist in  $\mathcal{E} \setminus \mathcal{S}$ .

The set S satisfying Definition 3 (if it exists) is called the *feedback edge set* of  $\Gamma$ . The interpretation is that S identifies edges with *outlying labels* that prohibit a consistent labelling to be found. Note that in the presence of noise we have to relax Eq. (3) and consider the following

$$\delta(z(i_1, i_2) * z(i_2, i_3) * \dots * z(i_{\ell}, i_1), 1_{\Sigma}) / \sqrt{\ell} \le \tau$$
(6)

where  $\tau \geq 0$  is a given threshold and it is assumed that  $\Sigma$  admits a metric function  $\delta : \Sigma \times \Sigma \rightarrow \mathbb{R}^+$ . The normalization factor  $\sqrt{\ell}$  takes into account error propagation when considering long cycles (Enqvist et al. 2011).

Outlying labels can be detected through the methods in Wahlström (2014), Cygan et al. (2012), which come from Graph theory community. Alternatively, Computer Vision solutions can be used, which include outlier rejection heuristics (Enqvist et al. 2011; Arrigoni et al. 2014b), Random Sample Consensus (RANSAC) (Govindu 2006; Olsson and Enqvist 2011), and Bayesian inference (Zach et al. 2010; Moulon et al. 2013; Bourmaud et al. 2014). However, these strategies are computationally demanding and do not scale well with the size of the graph.

#### 2.2 Group Synchronization

Let us assume that  $\Sigma$  is equipped with a metric function  $\delta : \Sigma \times \Sigma \to \mathbb{R}^+$  and let  $\rho : \mathbb{R}^+ \to \mathbb{R}^+$  be a non-negative non-decreasing function with a unique minimum at 0 and  $\rho(0) = 0$ . Some instances are the quadratic loss function  $\rho(y) = y^2$  or robust loss functions used in M-estimators (Holland and Welsch 1977).

**Definition 4** Let  $\Gamma = (\mathcal{V}, \mathcal{E}, z)$  be a  $\Sigma$ -labelled graph. Let  $\tilde{x} : \mathcal{V} \to \Sigma$  be a vertex labelling. We define the *consistency error* of  $\tilde{x}$  as the quantity

$$\epsilon(\widetilde{x}) = \sum_{(i,j)\in\mathcal{E}} \rho\Big(\delta\big(\widetilde{z}(i,j), z(i,j)\big)\Big)$$
(7)

where  $\tilde{z}$  is the edge labelling induced by  $\tilde{x}$ , namely  $\tilde{z}(i, j) = \tilde{x}(i) * \tilde{x}(j)^{-1}$ .

A vertex labelling is consistent if and only if it has zero consistency error. In practical applications a labelling with zero error hardly exists, since the edge labels are corrupted by noise, thus the goal is to address the following problem.



Fig. 4 The synchronization problem

**Definition 5** Given a  $\Sigma$ -labelled graph  $\Gamma = (\mathcal{V}, \mathcal{E}, z)$ , the *group synchronization problem* consists in finding a vertex labelling with minimum consistency error.

In other words, one wants to recover the unknown group elements (vertex labels) given a redundant set of noisy measurements of their ratios (edge labels), as shown in Fig. 4.

A related approach (e.g. Sharp et al. 2002) consists in minimizing the cost function (7) with respect to the edge labelling  $\tilde{z}$  while imposing the constraint that all the cycles are null (a.k.a. *cycle consistency*). For the  $\Sigma = \mathbb{R}^d$  case this is equivalent to synchronization (Proposition 2), whereas analogous results are not known in general.

The synchronization problem requires the graph to be connected, but *error compensation* happens only within cycles. The minimum number of relative measures is n - 1, which makes  $\mathcal{G}$  a tree. In this case every vertex can be labeled by simply propagating Eq. (5) along the tree, starting from the root labeled with the identity element. In this case, however, there is no remedy to error propagation: the error affecting an edge label propagates down to the leaves of the tree without compensation, as shown (e.g.) by experiments in Govindu (2004). In the synchronization problem, instead, the goal is to exploit *redundant* relative measures in a global fashion to improve the final estimate.

If the measures are also corrupted by outliers, one needs to solve a GFES problem beforehand, using a relaxed notion of null cycle, i.e., Eq. (6). Alternatively, a robust loss function can be used in (7) without detecting outliers explicitly (as done in Sect. 9.5).

#### 2.3 Synchronization over an Inverse Monoid

As observed in Arrigoni et al. (2017), the notion of synchronization can be extended to the case where  $\Sigma$  is an *inverse monoid*. One example is  $\Sigma = \mathcal{I}_d$ , which is a subsemigroup of  $S_d$ , resulting in *partial permutation synchronization* (Arrigoni et al. 2017).

**Definition 6** An *inverse semigroup*  $(\Sigma, *)$  is a semigroup in which for all  $s \in \Sigma$  there exists an element  $t \in \Sigma$  such that s = s \* t \* s and t = t \* s \* t. In this case, we write  $t = s^{-1}$  and call t the *inverse* of s. If  $\Sigma$  has an identity element  $1_{\Sigma}$  (i.e. it is a *monoid*), then it is called an *inverse monoid*.

**Remark 1** Inverses in an inverse semigroup have many of the same properties as inverses in a group, for instance,  $(a * b)^{-1} = b^{-1} * a^{-1}$  for all  $a, b \in \Sigma$ .

If  $\Sigma$  is an inverse monoid, then Eqs. (4) and (7) still make sense, with the provision that  $x(j)^{-1}$  now denotes the inverse of x(j) in the semigroup. Note that  $x(j)^{-1} * x(j)$  and  $x(j) * x(j)^{-1}$  are not necessarily equal to the identity, thus Eq. (4) is not equivalent to (5). The solution to the synchronization problem over an inverse monoid is defined up to a global (right) product with any element  $y \in \Sigma$  such that  $y * y^{-1} = 1_{\Sigma} = y^{-1} * y$ .

As concerns the notion of null-cycle, Eq. (3) still makes sense in the case of an inverse monoid, since it only involves products between elements of the set and the existence of a unit element  $1_{\Sigma}$ . However, Proposition 1 (and hence one direction of Corollary 1) does not straightforwardly extend to this case since the proof exploits Eq. (5), which is no longer equivalent to the consistency constraint. Regarding the other direction of Corollary 1, note that, if the consistency constraint holds, then the left side in Eq. (3) rewrites

$$z(i_{1}, i_{2}) * z(i_{2}, i_{3}) * \dots * z(i_{\ell}, i_{1})$$
  
=  $x(i_{1}) * \underbrace{x(i_{2})^{-1} * x(i_{2})}_{\neq 1_{\Sigma}} * x(i_{3})^{-1} * \dots * x(i_{\ell}) * x(i_{1})^{-1}$   
(8)

which, in general, does not coincide with the identity. Developing a complete theory for synchronization over an inverse monoid is outside the scope of this paper.

# **3** Synchronization over $\mathbb{R}^d$

In this section we consider the synchronization of real vectors with addition, namely  $(\Sigma, *) = (\mathbb{R}^d, +)$ , which is also known as *translation synchronization*, and we derive a direct solution following (Barooah and Hespanha 2007; Russel et al. 2011).

A vertex labelling  $\mathbf{x} : \mathcal{V} \to \mathbb{R}^d$  is consistent with a given edge labelling  $\mathbf{z} : \mathcal{E} \to \mathbb{R}^d$  if and only if<sup>2</sup>

$$\mathbf{x}_i - \mathbf{x}_j = \mathbf{z}_{ij} \ \forall \ (i, j) \in \mathcal{E}.$$
(9)

Note that we can view each component in Eq. (9) as a synchronization over  $(\mathbb{R}, +)$ .

If we denote the *incidence vector* of the edge (i, j) with

$$\boldsymbol{b}_{ij} = \begin{bmatrix} 0, \dots, 1, \dots, -1, \dots, 0 \end{bmatrix}^{\mathsf{T}}$$
(10)

then Eq. (9) can be written as

$$[\mathbf{x}_1, \dots, \mathbf{x}_n] \mathbf{b}_{ij} = \mathbf{z}_{ij} \quad \forall (i, j) \in \mathcal{E}$$
(11)

or, equivalently, in matrix form

$$XB = Z \tag{12}$$

where *B* is the  $n \times m$  incidence matrix of the *directed* graph  $\mathcal{G}$ , which has the vectors  $\mathbf{b}_{ij}$  as columns, *X* is the  $d \times n$  matrix obtained by juxtaposing all the vertex labels, namely  $X = [\mathbf{x}_1 \dots \mathbf{x}_n]$ , and *Z* is the  $d \times m$  matrix obtained by juxtaposing all the edge labels (ordered as in *B*), namely  $Z = [\mathbf{z}_{12} \dots \mathbf{z}_{ij} \dots]$ . See "Appendix B.1" for the definition of incidence matrix and related properties.

Applying the vectorization operator  $vec(\cdot)$  to both sides in (12) and using formula (87) we get

$$(B^{\mathsf{T}} \otimes I_d) \operatorname{vec}(X) = \operatorname{vec}(Z) \tag{13}$$

where  $I_d$  denotes the  $d \times d$  identity matrix and  $\otimes$  denotes the Kronecker product, defined in "Appendix A", which has the effect of "inflating" the incidence matrix in order to cope with the vector representation of the group elements.

Under the assumption that the graph is connected we have rank(B) = n - 1 and hence, using (89), we have rank( $B^{T} \otimes I_{d}$ ) = dn - d. The rank deficiency corresponds to the translation ambiguity. Since the solution to the synchronization problem is defined up to a global group element, we are allowed w.l.o.g. to arbitrarily set  $\mathbf{x}_{k} = \mathbf{0} = 1_{\Sigma}$  for a chosen  $k \in \mathcal{V}$ . Removing  $\mathbf{x}_{k}$  from the unknowns and the corresponding row in *B* leaves a full-rank matrix.

With a suitable choice of  $\delta(\cdot, \cdot)$  and  $\rho$  in Eq. (7), the consistency error of the synchronization problem writes

$$\epsilon(X) = \|(B^{\mathsf{T}} \otimes I_d) \operatorname{vec}(X) - \operatorname{vec}(Z)\|^2$$
(14)

where  $\|\cdot\|$  denotes the Euclidean norm. Thus the least-squares solution of Eq. (13) solves the synchronization problem. This

<sup>&</sup>lt;sup>2</sup> For simplicity of notation, hereafter we will use subscripts instead of parenthesis to denote indices of a node/edge labelling.

method is adopted in Barooah and Hespanha (2007), Russel et al. (2011) to localize a group of agents in a sensor network and in Arrigoni et al. (2016a) to recover camera positions in a structure-from-motion pipeline.

**Remark 2** If  $\mathbf{c} \in \{-1, 0, 1\}^m$  denotes the indicator vector of a circuit in  $\mathcal{G}$ , the cycle is null if and only if

$$Z\mathbf{c} = \mathbf{0}.\tag{15}$$

If the equations coming from all the circuits in a *cycle basis* are stacked, then the cycle consistency writes:

$$ZC^{\mathsf{T}} = 0 \tag{16}$$

where  $C \in \{-1, 0, 1\}^{(m-n+1)\times m}$  denotes the cycle matrix associated to the basis. See "Appendix B" or Carlone et al. (2011) for the definitions of cycle basis and cycle matrix. Using the vectorization operator and formula (87), cycle consistency can also be expressed as

$$(C \otimes I_d) \operatorname{vec} (Z) = \mathbf{0}.$$
(17)

It can be shown that the edge labels produced by the synchronization process are the closest to the input edge labels among those that yield null-cycles.

**Proposition 2** (Russel et al. 2011) If  $\tilde{X}$  is the least-squares solution to Eq. (13), then the induced edge labelling  $\tilde{Z} = \tilde{X}B$  solves the following constrained minimization problem

$$\min_{\widetilde{Z}} \|Z - \widetilde{Z}\|^2 \quad s.t. \ \widetilde{Z}C^{\mathsf{T}} = 0 \tag{18}$$

where C denotes the cycle matrix associated to a cycle basis of  $\mathcal{G}$ .

This result was also known in topography, in the context of the least-squares adjustment of leveling networks (Bjerhammar 1973).

# 4 Synchronization over GL(d)

In this section we consider the synchronization problem over the General Linear Group GL(d), which is the set of all  $d \times d$  invertible matrices, where the group operation \* reduces to matrix multiplication and  $1_{\Sigma} = I_d$ . A vertex labelling  $X : \mathcal{V} \to \mathbb{R}^{d \times d}$  is consistent with a given edge labelling  $Z : \mathcal{E} \to \mathbb{R}^{d \times d}$  if and only if  $Z_{ij} = X_i \cdot X_j^{-1}$ .

All the vertex/edge labels can be collected in two matrices  $X \in \mathbb{R}^{dn \times d}$  and  $Z \in \mathbb{R}^{dn \times dn}$  respectively, which are "matrices of matrices" defined as follows

$$X = \begin{bmatrix} X_1 \\ X_2 \\ \dots \\ X_n \end{bmatrix}, \quad Z = \begin{bmatrix} I_d & Z_{12} \dots & Z_{1n} \\ Z_{n1} & I_d & \dots & Z_{2n} \\ \dots & \dots & \dots \\ Z_{n1} & Z_{n2} \dots & I_d \end{bmatrix}.$$
 (19)

For a complete graph, the consistency constraint can be expressed in matrix form as

$$Z = X X^{-\flat} \tag{20}$$

where  $XX^{-b}$  contains the edge labels induced by X and  $X^{-b} \in \mathbb{R}^{d \times dn}$  denotes the block-matrix containing the inverse of each  $d \times d$  block of X, i.e.  $X^{-b} = [X_1^{-1} X_2^{-1} \dots X_n^{-1}]$ . Note that Eq. (20) implies that rank(Z) = d.

Remark 3 By computation it can be verified that

$$X^{-\flat}X = nI_d \tag{21}$$

and hence Z/n is idempotent.

If the graph is not complete then *Z* is not fully specified. In this case missing edges are represented as zero entries, i.e.  $Z_A := Z \circ (A \otimes \mathbf{1}_{d \times d})$  represents the matrix of the available measures, where  $\circ$  is the Hadamard product and *A* is the adjacency matrix of the graph  $\mathcal{G}$ , which gets "inflated" by the Kronecker product with  $\mathbf{1}_{d \times d}$  to match the block structure of the measures. Being a matrix of 0/1, the effect of its entrywise product with *Z* is to zero the unspecified entries and leave the others unchanged. See "Appendices B and A" for the definitions of Hadamard product and adjacency matrix, respectively. Hence the consistency constraint writes

$$Z_A = (XX^{-\flat}) \circ (A \otimes \mathbf{1}_{d \times d}).$$
(22)

With a suitable choice of  $\delta(\cdot, \cdot)$  and  $\rho$  in Eq. (7) the consistency error of the synchronization problem writes

$$\epsilon(X) = \|Z_A - (XX^{-\flat}) \circ (A \otimes \mathbf{1}_{d \times d})\|_F^2$$
(23)

where  $\|\cdot\|_F$  denotes the Frobenius norm. The Hadamard product with  $A \otimes \mathbf{1}_{d \times d}$  mirrors the summation over the edges in  $\mathcal{E}$  in the definition of the consistency error. The minimization of  $\epsilon$  is a non-linear least squares problem, for which closed-form solutions do not seem to exist. However, two direct solutions to a related version of the problem exist (Singer 2011; Singer and Shkolnisky 2011; Arie-Nachimson et al. 2012), which can be derived by considering the exact (noiseless) case.

#### 4.1 Spectral Solution

Let us consider the noiseless case, i.e.  $\epsilon = 0$ , and let us start assuming that the graph is complete. Using the consistency constraint and (21) we obtain

$$ZX = nX \tag{24}$$

which means that—in the absence of noise—the columns of X are d (independent) eigenvectors of Z corresponding to the eigenvalue n. Note that, since Z has rank d, all the other eigenvalues are zero, thus n is also the *largest* eigenvalue of Z.

We now consider the case of missing edges in which the graph is not complete and the adjacency matrix comes into play.

**Proposition 3** (Singer 2011; Singer and Shkolnisky 2011; Arie-Nachimson et al. 2012) *The columns of X are d (independent) eigenvectors of*  $(D \otimes I_d)^{-1}Z_A$  associated to the eigenvalue 1.

**Proof** In the case of missing data, it can be seen that Eq. (24) generalizes to

$$Z_A X = (D \otimes I_d) X \tag{25}$$

where D is the degree matrix of the graph (see "Appendix B.1"). Indeed, the i-th block-row in the above equation is

$$\sum_{j \text{ s.t. } (i,j)\in\mathcal{E}} Z_{ij}X_j = [D]_{ii}X_i$$
(26)

which is satisfied since  $Z_{ij} = X_i X_j^{-1}$ .

Note that the incomplete data matrix  $Z_A$  has full rank in general, thus 1 is not the unique nonzero eigenvalue of  $(D \otimes I_d)^{-1}Z_A$ , in contrast to the case of Eq. (24). However, it can be shown that 1 is the *largest* eigenvalue of such a matrix.

**Proposition 4** (Arrigoni et al. 2016c) The matrix  $(D \otimes I_d)^{-1}Z_A$  has real eigenvalues. The largest eigenvalue is 1 with multiplicity d.

**Proof** By computation it can be verified that

$$Z_A = \text{blkdiag}(X)(A \otimes I_d) \text{blkdiag}(X)^{-1}$$
(27)

where blkdiag(*X*) produces a  $dn \times dn$  block-diagonal matrix with  $d \times d$  blocks  $X_1, \ldots, X_n$  along the diagonal. Note that the diagonal matrix  $(D \otimes I_d)^{-1}$  commutes with blkdiag(*X*), since each  $d \times d$  block along its diagonal is a multiple of the identity matrix. Thus

$$(D \otimes I_d)^{-1} Z_A$$
  
= blkdiag(X)(D \otimes I\_d)^{-1} (A \otimes I\_d) blkdiag(X)^{-1}  
= blkdiag(X)((D^{-1}A) \otimes I\_d) blkdiag(X)^{-1} (28)

where the last equality follows from properties (85) and (86). Hence  $(D \otimes I_d)^{-1}Z_A$  is similar to the matrix  $(D^{-1}A) \otimes I_d$ , i.e., they have the same eigenvalues. The matrix  $D^{-1}A$  is the transition matrix of the graph  $\mathcal{G}$  (see "Appendix B.1"), which—as a consequence of the Perron–Frobenius theorem—has real eigenvalues and 1 is the largest eigenvalue (with multiplicity 1), if the graph is connected. Since the eigenvalues of the Kronecker product of two matrices are the product of the eigenvalue of  $(D^{-1}A) \otimes I_d$  is 1 with multiplicity d.

The proof of Proposition 4 has pointed out that—provided that Z is decomposable as  $Z = XX^{-b}$ —the matrix  $(D \otimes I_d)^{-1}Z_A$  has a particular structure that yields real eigenvalues, although it is not symmetric. In particular, the eigenvalues do not depend on the measured data, but they depend only on the structure of the graph  $\mathcal{G}$  (through the matrices D and A).

When noise is present, i.e.  $\epsilon \neq 0$ , the eigenvectors of  $(D \otimes I_d)^{-1}Z_A$  corresponding to the *d* largest eigenvalues are an estimate of the vertex labelling *X*. The presence of noise, however, cripples the structure of  $Z_A$ , i.e.  $Z_A \neq (XX^{-\flat}) \circ (A \otimes \mathbf{1}_{d \times d})$ , thus the eigenvalues and the eigenvectors may be complex. As a consequence, after computing the leading eigenvectors, the imaginary part is zeroed. This approach is adopted in Arrigoni et al. (2016a) in the *d* = 1 case to synchronize distances between camera pairs in a structure-from-motion pipeline. To the best of our knowledge, no general results are known linking this spectral solution to the synchronization cost function. Results in the special case where  $\Sigma = O(d)$  are reported in Sect. 7.1.

Ambiguity Since the eigenvalue 1 is repeated, the corresponding eigenvectors span a linear subspace, and hence any basis for such a space is a solution. However, a change of the basis in the eigenspace corresponds to right-multiply the eigenvectors by an invertible  $d \times d$  matrix, i.e., the solution to synchronization is defined up the action of an element of GL(d), as expected.

Projection Let U be the matrix containing the d leading eigenvectors of  $(D \otimes I_d)^{-1} Z_A$ , where the imaginary part (if any) is zeroed. The blocks of U are in general non-singular, hence they belong to GL(d). This means that the solution is *intrinsic*, and no projection is required.

**Remark 4** The top eigenvector can be computed by the power iteration method, which, considering (e.g.) the case of a complete graph, starts with a random vector  $\mathbf{x}_0 \in \mathbb{R}^{dn}$  and iterates the relation  $\mathbf{x}_{k+1} = Z\mathbf{x}_k / ||Z\mathbf{x}_k||$ , thus it requires to compute  $Z^k$ , for  $k = 1, 2, ..., k_{\text{max}}$ . It is observed in Singer (2011) that multiplying the matrix *Z* by itself integrates the consistency relation of triplets, while high order iterations exploit consistency relations of longer cycles. Indeed

$$Z_{ij}^2 = \sum_{k=1}^n Z_{ik} Z_{kj}$$
(29)

$$Z_{ij}^{3} = \sum_{k=1}^{n} \sum_{h=1}^{n} Z_{ik} Z_{kh} Z_{hj} \dots$$
(30)

Thus the top eigenvector integrates the consistency relation of all cycles.

## 4.2 Null-Space Solution

We now show that synchronization over  $(GL(d), \cdot)$  can also be expressed as a null-space problem. If  $\epsilon = 0$  Eq. (24) is equivalent to

$$(nI_{dn} - Z)X = 0 \tag{31}$$

which means that, if the graph is complete, the vertex labelling X coincides with the d-dimensional null-space of  $nI_{dn} - Z$ . In the case of missing edges, let us rewrite (25) as

$$(D \otimes I_d - Z_A)X = 0 \tag{32}$$

thus X belongs to the null-space of  $D \otimes I_d - Z_A$ .

Let us observe that the matrix  $D \otimes I_d$  coincides with  $(D \otimes \mathbf{1}_{d \times d}) \circ Z$ , since *Z* has identity blocks along its diagonal and  $D \otimes \mathbf{1}_{d \times d}$  is block-diagonal. Using the distributive property of the involved products, we obtain an equivalent expression for  $D \otimes I_d - Z_A$ 

$$D \otimes I_d - Z_A = ((D - A) \otimes \mathbf{1}_{d \times d}) \circ Z = (L \otimes \mathbf{1}_{d \times d}) \circ Z$$
(33)

where L = D - A is the Laplacian matrix of  $\mathcal{G}$  (see "Appendix B.1"), which gets "inflated" to a  $d \times d$ -block structure by the Kronecker product with  $\mathbf{1}_{d \times d}$ , to match the block structure of Z.

Note that in practice one cannot measure the matrix  $(L \otimes \mathbf{1}_{d \times d}) \circ Z$ , since the full *Z* is not available. In fact, only the product  $Z \circ (A \otimes \mathbf{1}_{d \times d})$  is available. Therefore, the left side in (33) will be used in real scenarios. However, the right side emphasizes the presence of the Laplacian matrix, which is useful to prove that the null-space of  $D \otimes I_d - Z_A$  is d-dimensional, as happens in the case of a complete graph.

**Proposition 5** (Arrigoni et al. 2016c) *The matrix*  $D \otimes I_d - Z_A$  *has a d-dimensional null-space.* 

**Proof** By computation it can be verified that

$$D \otimes I_d - Z_A = (L \otimes \mathbf{1}_{d \times d}) \circ Z$$
  
= blkdiag(X)(L \otimes I\_d) blkdiag(X)^{-1} (34)

which means that  $D \otimes I_d - Z_A$  and  $L \otimes I_d$  are similar, thus they have the same rank. The rank of the Laplacian matrix is n - 1, under the assumption that the graph is connected (see "Appendix B.1"). Since the rank of the Kronecker product of two matrices is the product of the rank of the matrices, we obtain

$$\operatorname{rank}(D \otimes I_d - Z_A) = \operatorname{rank}(L)\operatorname{rank}(I_d) = dn - d \quad (35)$$

thus we have the thesis.

When noise is present, an estimate of X is given by the right singular vectors of  $D \otimes I_d - Z_A$  corresponding to the d least singular values, which solve the following problem

$$\min_{X^{\mathsf{T}}X=nI_d} \| (D \otimes I_d - Z_A)X \|_F^2.$$
(36)

A formal relationship between this cost function and the consistency error of the synchronization problem has still to be found.

The considerations about ambiguity/projection made for the spectral method apply also to the null-space solution, modulo the fact that singular vectors are real even in the absence of noise, so no rounding is needed.

#### 4.3 Additive Solution

We observe that synchronization over the General Linear Group can be cast to a translation synchronization, exploiting the fact that GL(d) has the structure of a Lie group (Varadarajan 1984), where the associated Lie algebra consists of all  $d \times d$  real matrices with the commutator operator serving as the Lie bracket, namely [Y, W] = YW - WY. Informally, a Lie group can be locally viewed as topologically equivalent to a vector space, and the local neighbourhood of any group element can be adequately described by its tangent space, whose elements form a Lie algebra. The Lie algebra and the Lie group are related by the exponential mapping, and the inverse mapping from the Lie group to the Lie algebra is given by matrix logarithm.

By taking the logarithm, the consistency constraint of the synchronization problem over GL(d), that is  $Z_{ij} = X_i X_j^{-1}$ , can be transformed into the consistency constraint of an additive group, namely

$$\log(Z_{ij}) = \log(X_i) - \log(X_j) \tag{37}$$

assuming that the each of the above matrices admits a unique real logarithm. Specifically, by vectorizing each side in (37), a relation of the form (9) is obtained, which defines a translation synchronization problem. Thus the solution can be found by solving a linear system in the least-squares sense, as done in Sect. 3, or via IRLS (to gain robustness to outliers), as explained in Sect. 9.5. In other words, the synchronization problem is addressed in the Lie algebra rather than in the group.

However, as observed in Govindu (2004), the Euclidean distance in the Lie algebra does not coincide with the Riemannian distance in the group, but it constitutes a first-order approximation, as stated by the Baker–Campbell–Hausdorff formula (Varadarajan 1984). For this reason, in Govindu (2004, 2006), Chatterjee and Govindu (2013) the solution is found by iterating between solving the linear system in the Lie algebra and remapping onto the group.

As a final note, it is straightforward to see that the approach of Sect. 4 also applies to synchronization over  $(\mathbb{R}\setminus\{0\}, \cdot)$ , for it coincides with  $(GL(1), \cdot)$ .

# 5 Synchronization over Subgroups of GL(d)

The analysis carried out in Sect. 4 can be extended to the case where  $\Sigma$  is a subgroup of GL(d), i.e., it can be embedded in  $\mathbb{R}^{d \times d}$ , where the group operation \* reduces to matrix multiplication and  $1_{\Sigma} = I_d$ . In this case Propositions 4 and 5 still hold, and the synchronization problem can be addressed either via the spectral solution, which computes the top *d* eigenvectors of  $(D \otimes I_d)^{-1}Z_A$  (which may be complex in the presence of noise), or via the null-space solution, which computes the least *d* right singular vectors of  $D \otimes I_d - Z_A$ . Alternatively, the (approximate) null-space of  $I_{dn} - (D \otimes I_d)^{-1}Z_A$  can be computed, as done in Schroeder et al. (2011).

Let U be the  $dn \times d$  matrix containing either the output of the spectral method or the null-space solution. Recall that U is not uniquely determined (even in the absence of noise) but it is defined up to (right) multiplication by an element of GL(d), since any basis for the null-space of  $D \otimes I_d - Z_A$  (or, equivalently, any basis for the eigenspace of  $(D \otimes I_d)^{-1}Z_A$  associated with eigenvalue 1) is a solution. Thus a procedure that reduces the ambiguity up to an element of  $\Sigma$ is required, since the solution to the synchronization problem is inherently defined up to an element of the group.

Note that closure is not always guaranteed. In other words, the spectral and the null space methods produce an *extrinsic* estimate of the vertex labelling X which needs to be eventually projected onto  $\Sigma$ . This approach is clearly suboptimal with respect to working *in* the group, and represents the price to be paid for simplicity and computational efficiency.

In the following sections we will analyze synchronization is some subgroups of GL(d) (see Table 1 and Fig. 1). They basically differ from each other by the ambiguity fixing and the projection stages.

**Remark 5** Synchronization over a subgroup of GL(d) can also be addressed via the approach detailed in Sect. 4.3,

as done in Govindu (2004, 2006), Chatterjee and Govindu (2013). Particularly interesting are the  $\Sigma = SO(3)$  and  $\Sigma = SE(3)$  cases, where the associated Lie algebras are described by 3 and 6 parameters, respectively, and the exponential and logarithm maps admit closed form expressions (Moakher 2002; Cardoso and Leite 1999).

#### 5.1 Synchronization over SL(d)

The Special Linear Group SL(d) is the set of  $d \times d$  matrices with unit determinant. Synchronization over SL(3) is studied in Schroeder et al. (2011) within the context of multiple-view homography estimation, that is why the problem is referred to as *homography synchronization*.

Ambiguity U is the solution up to multiplication by element of GL(d), which can be reduced to SL(d) after permutation of two columns of U s.t.  $det(U_1) > 0$  and division by  $\sqrt[d]{det(U_1)}$ , where  $U_1$  denotes the first  $d \times d$  block in U.

Projection In order to obtain elements of SL(d) from U, each  $d \times d$  block in U, denoted by  $U_i$ , must be scaled to unit determinant, which can be done by dividing  $U_i$  by  $\sqrt[d]{\det(U_i)}$ . In the case of the spectral solution, before performing such projection, the imaginary part of the eigenvectors is zeroed.

#### 5.2 Synchronization over O(d)

The Orthogonal Group O(d) is the set of orthogonal transformations in *d*-space, which admits a matrix representation through  $d \times d$  orthogonal matrices. An important subgroup of O(d) is the Special Orthogonal Group SO(d), that is the set of orthogonal matrices with determinant 1, which represent rotations in *d*-space. Synchronization over SO(d) is also known as *rotation (angular) synchronization* (Singer 2011) or *multiple rotation averaging* (Hartley et al. 2013; Wilson et al. 2016). A comprehensive survey on existing solutions can be found in Carlone et al. (2015), Tron et al. (2016).

From the theoretical perspective, synchronization over SO(3) is analyzed in depth in Hartley et al. (2013), Wilson et al. (2016). In Hartley et al. (2013) the consistency error (7) is studied under the choice  $\rho(y) = y^p$  (with  $p \ge 1$ ) and several distance measures are considered, including quaternion, angular (geodesic) and chordal distances, where each metric is related to a particular parametrization of the rotation space. In Wilson et al. (2016) it is shown that smaller and well-connected graphs are easier than larger and noisy ones, based on a local convexity analysis. Further theoretical analysis is reported in Boumal et al. (2014) where Cramèr–Rao bounds for synchronization over SO(d) are derived, namely lower bounds on the variance of unbiased estimators, assuming a certain noise model.

The spectral method was introduced in Singer (2011) for  $\Sigma = SO(2)$  and extended in Singer and Shkolnisky (2011),

Arie-Nachimson et al. (2012) to  $\Sigma = SO(3)$ . A related approach is adopted in Martinec and Pajdla (2007), where the  $\Sigma = SO(3)$  case is considered and the least eigenvectors of  $D \otimes I_3 - Z_A$  are computed (instead of the least right singular vectors).

**Remark 6** Note that both Z and  $Z_A$  are symmetric even in the presence of noise, since, by assumption, a group-labelled graph satisfies  $Z_{ji} = Z_{ij}^{-1}$  for all  $(i, j) \in \mathcal{E}$ , which becomes  $Z_{ji} = Z_{ij}^{\mathsf{T}}$  in the  $\Sigma = O(d)$  case. The matrix  $(D \otimes I_d)^{-1} Z_A$ is not symmetric, but it is similar to the *symmetric* matrix  $(D \otimes I_d)^{-1/2} Z_A (D \otimes I_d)^{-1/2}$ , thus its eigenvalues are real and it admits an orthonormal basis of real eigenvectors.

Ambiguity The solution U is defined up to an element of O(d), since a change of the orthonormal basis in the eigenspace (or null-space) corresponds to (right) multiplication by an orthogonal  $d \times d$  matrix. This reduces to SO(d)after permutation of columns of U such that  $det(U_1)$  is positive.

Projection Each  $d \times d$  block of U is not guaranteed to belong to O(d) and has to be projected onto the group. Such projection can be obtained by solving an orthogonal Procrustes problem, e.g. via Singular Value Decomposition (SVD) (Keller 1975). Specifically, if  $Q \in \mathbb{R}^{d \times d}$  is a given matrix, then the nearest orthogonal matrix (in the Frobenius norm sense) is given by

$$R = WV^{\mathsf{T}} \in O(d) \tag{38}$$

where  $Q = WSV^{\mathsf{T}}$  denotes the singular value decomposition of Q. If the synchronization problem over  $\Sigma = SO(d)$  is considered, the projection step is slightly different. Specifically, the nearest rotation matrix (in the Frobenius norm sense) is given by

$$R = W \operatorname{diag}([1, \dots, 1, \operatorname{det}(WV^{\mathsf{T}})])V^{\mathsf{T}} \in SO(d).$$
(39)

Note that the final rounding step which is required for GL(d), i.e., zeroing the imaginary part of the eigenvectors, is not necessary here due to Remark 6.

## 5.3 Synchronization over GA(d)

The General Affine Group GA(d) is the set of invertible affine transformations in *d*-space, which admits a matrix representation through  $(d + 1) \times (d + 1)$  matrices

$$GA(d) = \left\{ \begin{bmatrix} M & \mathbf{t} \\ \mathbf{0}^{\mathsf{T}} & 1 \end{bmatrix}, \ s.t. \ M \in GL(d), \ \mathbf{t} \in \mathbb{R}^d \right\}.$$
(40)

Synchronization over GA(d) is referred to as *affine synchronization* (Bernard et al. 2015; Santellani et al. 2018).

Ambiguity The solution U is unique up to the action of an element of GL(d + 1), which can be reduced to GA(d) by computing a linear combination of the columns of U such that the output matrix has the vector  $[\mathbf{0}_{1\times d}1]$  in rows multiple of d + 1. More precisely, let  $F \in \mathbb{R}^{n \times (d+1)n}$  be the 0/1-matrix such that  $FU \in \mathbb{R}^{n \times (d+1)}$  consists of the rows of U with indices multiple of d + 1. The coefficients  $\mathbf{a}, \mathbf{b} \in \mathbb{R}^{d+1}$  of the linear combination are solution of

$$FU\mathbf{a} = \mathbf{0}_{n \times 1}, \quad FU\mathbf{b} = \mathbf{1}_{n \times 1} \tag{41}$$

where the first equation has a *d*-dimensional solution space. If  $\mathbf{a}_1, \ldots, \mathbf{a}_d$  denotes a basis for the null-space of FU, then U is transformed into  $U[\mathbf{a}_1, \ldots, \mathbf{a}_d, \mathbf{b}]$ . In the presence of noise, Eq. (41) is solved in the least squares sense.

*Projection* In order to project the solution onto GA(d), the rows of U multiple of d + 1 are forced to  $[\mathbf{0}_{1 \times d} \mathbf{1}]$ , and the imaginary part of the eigenvectors is zeroed (in the case of the spectral solution).

#### 5.4 Synchronization over SE(d)

The Special Euclidean Group SE(d) is the set of direct isometries (or rigid motions) in *d*-space, which admits a matrix representation through  $(d + 1) \times (d + 1)$  matrices

$$SE(d) = \left\{ \begin{bmatrix} R & \mathbf{t} \\ \mathbf{0}^{\mathsf{T}} & 1 \end{bmatrix}, \ s.t. \ R \in SO(d), \mathbf{t} \in \mathbb{R}^d \right\}.$$
(42)

Synchronization over SE(d) is also known as *rigid-motion* synchronization (Arrigoni et al. 2016c; Bernard et al. 2015) or motion averaging (Govindu 2004, 2017) or pose-graph optimization (Carlone et al. 2015). The spectral and null-space solutions can be regarded as the extension of the rotation synchronization approach introduced in Singer (2011), Singer and Shkolnisky (2011), Arie-Nachimson et al. (2012), and were developed independently by Bernard et al. (2015) and Arrigoni et al. (2016c).

Ambiguity The solution U is unique up to the action of an element of GL(d+1). In order to reduce the ambiguity up to an element of SE(d) we must design a suitable transformation that maps, e.g., the first block of U onto an element of SE(d). In order to do that let us apply a linear combination of the columns of U such that the (d + 1)th row becomes  $[\mathbf{0}_{1\times d}1]$ , as in the  $\Sigma = GA(d)$  case, and subsequently transform U such that  $U_1$ , the first  $d \times d$  block of U, becomes an element of SO(d). Let  $U_1 = RP$  be the polar decomposition of  $U_1$ , with  $R \in O(d)$  and P symmetric positive definite. Since  $U_1$ is invertible, then  $R = U_1P^{-1}$ , and  $P^{-1}$  is the sought transformation. In the presence of noise, a least square solution that brings every row multiple of d + 1 as close as possible to  $[\mathbf{0}_{1\times d}1]$  is sought, as in the affine case. **Projection** In order to project the solution onto SE(d)—as in Belta and Kumar (2002)—the rows of U multiple of d + 1 are forced to  $[\mathbf{0}_{1\times d}1]$  and the  $d \times d$  rotation blocks are projected onto SO(d). In the case of the spectral solution, before performing such projection, the imaginary part of the eigenvectors is zeroed.

*Two-Step Synchronization* Since the Special Euclidean Group is the the semi-direct product of SO(d) and  $\mathbb{R}^d$ , synchronization over SE(d) can be alternatively addressed by breaking the problem into rotation and translation and solving the two sub-problems separately, as done (e.g.) in Arrigoni et al. (2016a). Let  $X_i \in SE(d)$  denote the (unknown) label of node *i* 

$$X_i = \begin{bmatrix} R_i \ \mathbf{t}_i \\ \mathbf{0}^{\mathsf{T}} \ 1 \end{bmatrix}$$
(43)

where  $R_i \in SO(d)$  and  $\mathbf{t}_i \in \mathbb{R}^d$  denote the rotation and translation components of the rigid motion, respectively. Similarly, each edge label  $Z_{ij} \in SE(d)$  can be expressed as

$$Z_{ij} = \begin{bmatrix} R_{ij} \ \mathbf{t}_{ij} \\ \mathbf{0}^{\mathsf{T}} \ \mathbf{1} \end{bmatrix}$$
(44)

with  $R_{ij} \in SO(d)$  and  $\mathbf{t}_{ij} \in \mathbb{R}^d$ . Using this notation, the consistency constraint for synchronization over SE(d), namely  $Z_{ij} = X_i X_j^{-1}$ , can be equivalently rewritten as

$$R_{ij} = R_i R_j^{\mathsf{T}} \tag{45}$$

$$\mathbf{t}_{ij} = -R_i R_j^{\mathsf{T}} \mathbf{t}_j + \mathbf{t}_i.$$
(46)

Note that Eq. (45) defines a rotation synchronization problem, thus the rotation components of the unknown vertex labels can be recovered as explained in Sect. 5.2. Equation (46) can be equivalently written as

$$-R_i^{\mathsf{T}}\mathbf{t}_{ij} = R_j^{\mathsf{T}}\mathbf{t}_j - R_i^{\mathsf{T}}\mathbf{t}_i = \mathbf{x}_i - \mathbf{x}_j$$
(47)

using the substitution  $\mathbf{x}_i = -R_i^{\mathsf{T}}\mathbf{t}_i$ . Thus—assuming that rotations have been computed beforehand—recovering the translation components of the vertex labels can be reduced to a translation synchronization problem (as defined in Sect. 3), where the edge labels are given by  $\mathbf{z}_{ij} = -R_i^{\mathsf{T}}\mathbf{t}_{ij}$ .

## 5.5 Synchronization over $S_d$

The Symmetric Group  $S_d$  is the set of bijections between d objects, which admits a matrix representation through  $d \times d$  permutation matrices. A permutation matrix is such that exactly one entry in each row and column is equal to 1 and all other entries are 0. Synchronization over  $S_d$  is also known

as *permutation synchronization*, which finds application in multi-view matching (see Sect. 8.2). The spectral solution was introduced in Pachauri et al. (2013) for a complete graph (based on Singer (2011)) and subsequently extended in Shen et al. (2016) to the case of missing data.

Ambiguity The solution U is defined up to an element of O(d), as a consequence of Remark 6. However, the solution to permutation synchronization is inherently defined up to an element of  $S_d$ . Let Q be the unknown orthogonal transformation such that X = UQ, where X is the solution, a matrix whose blocks are in  $S_d$ . Let  $U_1$  be the first d rows of U, then from  $X_1 = U_1Q$  we have  $Q = U_1^TX_1$ , i.e.,  $Q = U_1^T$  up to a permutation.  $U_1$  is indeed orthogonal because  $U_1U_1^T = X_1X_1^T = I_d$ .

**Projection** Even after fixing the ambiguity, the blocks of U will not be permutation matrices (in general). We need to project them onto  $S_d$  by solving the so-called permutation Procrustes problem, e.g., with the Kuhn–Munkres algorithm (Kuhn 1955). Note that the spectral solution returns real eigenvectors, as in the  $\Sigma = O(d)$  case.

# 6 Synchronization over $\mathcal{I}_d$

Let us consider now the Symmetric Inverse Semigroup  $\mathcal{I}_d$ , that is the set of bijections between (different) subsets of d objects, which admits a matrix representation through  $d \times d$  partial permutation matrices. A *partial* permutation matrix—which represents matches between different objects (see Fig. 6)—has at most one nonzero entry in each row and column, and these nonzero entries are all 1. The synchronization problem over  $\mathcal{I}_d$  is also known as *partial permutation synchronization* (Arrigoni et al. 2017).

The set  $\mathcal{I}_d$  is an inverse monoid with respect to matrix multiplication (and a subsemigroup of  $\mathcal{S}_d$ ) where the inverse is given by matrix transposition (see Definition 6). Let  $X_i \in$  $\mathcal{I}_d$  denote the (unknown) label of vertex *i* and let  $Z_{ij} \in \mathcal{I}_d$ denote the (known) label of edge  $(i, j) \in \mathcal{E}$ , which are linked by the consistency constraint

$$Z_{ij} = X_i X_j^{\mathsf{T}}.\tag{48}$$

If  $[X_i]_{(h,k)} = 1$  for some index *h* we say that "node *i* sees object *k*". Note that  $X_i^T X_i$  is not equal, in general, to the identity, unless  $X_i \in S_d$ . Indeed,  $[X_i^T X_i]_{(k,k)} = 1$  if node *i* sees object *k* and  $[X_i^T X_i]_{(k,k)} = 0$  otherwise. However, it can be checked that  $X_i^T X_i \leq I_d$ .

**Remark 7** If  $\{(i_1, i_2), (i_2, i_3), \dots, (i_{\ell}, i_1)\}$  denotes a circuit in  $\mathcal{G}$ , then—assuming that the consistency constraint (48) holds—the composition of edge labels along its edges becomes

$$Z_{i_1i_2}Z_{i_2i_3} \cdot \cdots \cdot Z_{i_\ell i_1}$$

$$= X_{i_1} \underbrace{X_{i_2}^{\mathsf{T}} X_{i_2}}_{\leq I_d} X_{i_3}^{\mathsf{T}} \cdot \cdots \cdot X_{i_\ell} X_{i_1}^{\mathsf{T}} \leq I_d.$$
(49)

In contrast to the case of total permutations, where synchronization implies that compositions of edge labels over circuits must be equal to the identity (see Corollary 1), in the case of partial permutations we obtain that compositions of edge labels over circuits must be a subset of the identity, as observed also in Bernard et al. (2019). In other words, with reference to the multi-view matching application (see Sect. 8.2), due to potential missing matches (i.e. zero rows/columns in  $Z_{ij}$ ) along a cyclic path, some of the original matches may vanish, and only those matches that are seen in all the images survive.

## 6.1 Spectral Solution

Despite the fact that the group structure is missing,<sup>3</sup> it can be shown that a spectral solution can be derived under the assumption that the graph is complete. Let X and Z be two block-matrices containing the vertex labels and edge labels respectively—defined as in (19)—so that the consistency constraint becomes  $Z = XX^T$  with Z of rank d. Note that here the diagonal of Z is not filled with identity matrices, in general. When all the objects seen by node i are different from those seen by node j we have  $X_i X_j^T = 0$ , resulting in a zero block in Z.

**Proposition 6** (Arrigoni et al. 2017) *The columns of X are d* (*orthogonal*) *eigenvectors of Z and the corresponding eigenvalues are contained in the diagonal of the following d*  $\times$  *d matrix* 

$$V := X^{\mathsf{T}} X = \sum_{i=1}^{n} X_{i}^{\mathsf{T}} X_{i}.$$
 (50)

**Proof** Using (50) and the consistency constraint, we obtain

$$ZX = XV \tag{51}$$

which is a spectral decomposition, i.e. the columns of X are d eigenvectors of Z and the corresponding eigenvalues are on the diagonal of V. Recall that Z admits an orthonormal basis of real eigenvectors since it is symmetric.

Although  $\mathcal{I}_d$  is not a group, an eigenvalue decomposition problem has been obtained, where the non-zero eigenvalues are contained in the diagonal of V. Specifically, the k-th eigenvalue counts how many nodes see object k, thus all the eigenvalues are integer numbers lower than or equal to n. This implies that, when the number of objects is larger than the number of nodes (i.e., d > n)—which is likely to happen in the multi-view matching application—the eigenvalues are repeated. In the case of total permutations (i.e.  $\Sigma = S_d$ ) all the nodes see all the objects, thus  $V = nI_d$  and all the eigenvalues are equal, hence Eq. (51) reduces to (24). In the presence of noise, the eigenvectors of Z corresponding to the d largest eigenvalues are computed.

This result holds for a complete graph. When it is not, Proposition 6 can not be straightforwardly extended as we did for  $\Sigma = GL(d)$  case.

**Remark 8** Equation (51) could also be expressed as a null-space problem, but in that case the matrix V, which is unknown, have to be estimated somehow.

Ambiguity Note that the reverse of Proposition 6 is not true in general, i.e., the matrix U is not necessarily equal to the vertex labelling X. Indeed, U is not uniquely determined if the eigenvalues of Z are repeated. So we have to face the problem of how to select, among the infinitely many Us, the one that resembles X, a matrix composed of partial permutations. Note that the reasoning reported in Sect. 5.5 does not apply here, since the first block  $U_1$  is not orthogonal in general, due to the presence of zero rows. A key observation is reported in the following proposition, suggesting that such a problem can be solved via clustering techniques.

**Proposition 7** (Arrigoni et al. 2017) Let U be the  $nd \times d$ matrix composed by the d leading eigenvectors of Z; then U has d + 1 different rows (in the absence of noise). One of these is the zero row.

**Proof** Let  $\lambda_1, \lambda_2, \ldots, \lambda_\ell$  denote all the *distinct* eigenvalues of Z (with  $\ell \leq d$ ), and let  $m_1, m_2, \ldots, m_\ell$  be their multiplicities such that  $\sum_{k=1}^{\ell} m_k = d$ . Let  $U_{\lambda_k}$  denote the  $m_k$ columns of U corresponding to the eigenvalue  $\lambda_k$ , and let  $X_{\lambda_k}$ be the corresponding columns of X. Up to a permutation of the columns, we have

$$U = [U_{\lambda_1} \ U_{\lambda_2} \ \dots \ U_{\lambda_\ell}], \quad X = [X_{\lambda_1} \ X_{\lambda_2} \ \dots \ X_{\lambda_\ell}].$$
(52)

Since  $U_{\lambda_k}$  and  $X_{\lambda_k}$  are (orthogonal) eigenvectors corresponding to the same eigenvalue, there exists an orthogonal matrix  $Q_k \in \mathbb{R}^{m_k \times m_k}$  representing a change of basis in the eigenspace of  $\lambda_k$ , such that  $U_{\lambda_k} = X_{\lambda_k} Q_k$ . In matrix form this rewrites

$$U = X \underbrace{\text{blkdiag}(Q_1, Q_2, \dots, Q_\ell)}_{Q}.$$
(53)

Note that the rows of *X* are the rows of  $I_d$  plus the zero row. Since *Q* is invertible (hence injective), U = XQ has only d + 1 different rows as well.

<sup>&</sup>lt;sup>3</sup> Note that  $0 \in \mathcal{I}_d$ , thus we can not distinguish between the case of missing measures and the case of missing correspondences between nodes, as it was the case for the subgroups of GL(d).

Specifically, an estimate of the vertex labelling can be obtained by clustering the rows of U into d + 1 clusters (e.g. with k-means), then assigning the centroid which is closest to zero to the zero row, and arbitrarily assigning each of the other d centroids to a row of  $I_d$ . This arbitrary assignment corresponds to the fact that the solution to partial permutation synchronization is defined up to an element of  $S_d$ .

Projection Even after fixing the ambiguity, valid permutation matrices may not be obtained. Indeed, since there are no constraints in the clustering phase, it may happen that different rows of a  $d \times d$  block in U are assigned to the same cluster, resulting in more than one entry per column equal to 1. For this reason, for each  $d \times d$  block in U, the partial permutation matrix that best maps such block into the set of centroids has to be computed (e.g. via the Kuhn–Munkres algorithm Kuhn 1955), and such permutation is output as the sought solution.

# 7 Spectral and Other Relaxations

In this section we concentrate mainly on the  $\Sigma = O(d)$  case. We frame the spectral solution as an instance of a constraint relaxation pattern by setting it side by side with rank relaxation (Arrigoni et al. 2014a, 2016b; Zhou et al. 2015) and semidefinite programming (Singer 2011; Chen et al. 2014; Rosen et al. 2015), which are not closed-form, though. Moreover we highlight the link between the spectral solution and the consistency error of the synchronization problem (which is available only for  $\Sigma = O(d)$ ).

Note that this is a special case, since the inverse equals matrix transposition, thus the consistency constraint rewrites  $Z_{ij} = X_i X_i^{\mathsf{T}}$  which, if the graph is complete, is equivalent to

$$Z = XX^{\mathsf{T}}.$$
(54)

Such a decomposition implies that, if  $\epsilon = 0$ , the matrix Z is *symmetric and positive semidefinite*, besides being low-rank. In the case of missing edges, the consistency constraint translates into

$$Z_A = (XX^{\mathsf{T}}) \circ (A \otimes \mathbf{1}_{d \times d}) \tag{55}$$

and the consistency error of the synchronization problem becomes

$$\epsilon(X) = \sum_{(i,j)\in\mathcal{E}} \|Z_{ij} - X_i X_j^{\mathsf{T}}\|_F^2$$
  
=  $\|Z_A - (XX^{\mathsf{T}}) \circ (A \otimes \mathbf{1}_{d \times d})\|_F^2.$  (56)

In particular, since the Frobenius norm of a matrix can be defined in terms of its trace, Eq. (56) can be expressed as

$$\epsilon(X)$$

$$= \sum_{i,j=1}^{n} \operatorname{tr}(Z_{ij}^{\mathsf{T}} Z_{ij}) + \operatorname{tr}(X_j X_i^{\mathsf{T}} X_i X_j^{\mathsf{T}}) - 2\operatorname{tr}(X_i^{\mathsf{T}} Z_{ij} X_j)$$
$$= \sum_{i,j=1}^{n} \operatorname{tr}(Z_{ij}^{\mathsf{T}} Z_{ij}) + d - 2\operatorname{tr}(X_i^{\mathsf{T}} Z_{ij} X_j)$$
(57)

where a complete graph is considered and the last equality holds since  $X_i^{\mathsf{T}} X_i = I_d = X_i^{\mathsf{T}} X_j$ . Therefore

$$\min_{X \in O(d)^n} \epsilon(X) \longleftrightarrow \max_{X_1, \dots, X_n \in O(d)} \sum_{i, j=1}^n \operatorname{tr}(X_i^{\mathsf{T}} Z_{ij} X_j)$$
$$= \max_{X \in O(d)^n} \operatorname{tr}(X^{\mathsf{T}} Z X).$$
(58)

Solving the synchronization problem over O(d) is difficult since the feasible set is non-convex, and the cost function may have multiple local minima in different regions of attraction, as shown in Hartley et al. (2013). Several relaxations will be considered.

#### 7.1 Spectral Relaxation

Let us start with the case of a complete graph and let us consider the following minimization problem

$$\min_{X^{\mathsf{T}}X=nI_d} \|Z - XX^{\mathsf{T}}\|_F^2$$
(59)

where the columns of X are constrained to be orthogonal rather than imposing that each  $d \times d$  block in X belongs to O(d). This is called the *spectral relaxation*, and, reasoning as in Eq. (58), it can be shown to be equivalent to the following generalized Rayleigh problem

$$\max_{X^{\mathsf{T}}X=nI_d} \operatorname{tr}(X^{\mathsf{T}}ZX).$$
(60)

**Proposition 8** Equation (60), with Z symmetric, admits a closed-form solution given by the d leading eigenvectors of Z.

**Proof** Let  $\mathcal{F}$  be the unconstrained cost function corresponding to problem (60), namely

$$\mathcal{F}(X) = \operatorname{tr}(X^{\mathsf{T}}ZX) + \operatorname{tr}(\Lambda(X^{\mathsf{T}}X - nI_d))$$
(61)

where  $\Lambda \in \mathbb{R}^{d \times d}$  is a symmetric matrix of unknown Lagrange multipliers. Setting to zero the partial derivatives of  $\mathcal{F}$  with respect to X we obtain

$$\frac{\partial \mathcal{F}}{\partial X} = 2ZX + 2X\Lambda = 0 \Rightarrow ZX = -X\Lambda.$$
(62)

Let  $\mathbf{u}_i$  denote *d* eigenvectors of *Z* (normalized so that  $||\mathbf{u}_i|| = \sqrt{n}$ ) for i = 1, ..., d and let  $\lambda_i$  be the corresponding eigenvalues. Then  $X = [\mathbf{u}_1 \ \mathbf{u}_2 \ ... \ \mathbf{u}_d]$  satisfies both (62) and the constraint  $X^T X = nI_d$ , with  $\Lambda = -\text{diag}([\lambda_1, \lambda_2, ..., \lambda_d])$ . In other words, any set of *d* eigenvectors is a stationary point for the objective function  $\mathcal{F}$ . The corresponding stationary value is given by  $n(\lambda_1 + \lambda_2 + ... \lambda_d)$ , hence the maximum is attained if  $\mathbf{u}_i$  are the eigenvectors of *Z* corresponding to the *d* largest eigenvalues.

If the graph is not complete, we consider a different definition for the consistency error

$$\epsilon(X) = \sum_{(i,j)\in\mathcal{E}} [D]_{ii}^{-1} \|Z_{ij} - X_i X_j^{\mathsf{T}}\|_F^2$$
(63)

in which the term associated to the edge (i, j) is weighted with the inverse of the degree of node *i*. Reasoning as before we get

$$\min_{X \in O(d)^n} \epsilon(X)$$

$$\iff \max_{X_1, \dots, X_n \in O(d)} \sum_{(i,j) \in \mathcal{E}} [D]_{ii}^{-1} \operatorname{tr}(X_i^{\mathsf{T}} Z_{ij} X_j)$$

$$= \max_{X \in O(d)^n} \operatorname{tr}(X^{\mathsf{T}}(D \otimes I_d)^{-1} Z_A X)$$
(64)

which, if the spectral relaxation is adopted, becomes

$$\max_{X^{\mathsf{T}}X=nI_d} \operatorname{tr}(X^{\mathsf{T}}(D \otimes I_d)^{-1} Z_A X)$$
(65)

whose solution is given by the *d* leading eigenvectors of  $(D \otimes I_d)^{-1} Z_A$ .

We have shown, following (Arie-Nachimson et al. 2012), that the spectral solution minimizes the consistency error of the synchronization problem under relaxed constraints. Note that this result is due to the special structure of O(d), it also holds for subgroups of O(d), and it is not available in general for synchronization over GL(d).

Concerning the Symmetric Inverse Semigroup, since Z is symmetric, Proposition 8 holds. However, the computation above linking Problem (60) to the synchronization cost function is no longer valid since the equation  $X_i^T X_i = I_d = X_i^T X_j$  does not hold for  $\Sigma = \mathcal{I}_d$ .

# 7.2 Rank Relaxation

Let  $\Omega = A \otimes \mathbf{1}_{d \times d}$  denote the *pattern* (also known as the *sampling set*) of  $Z_A$ , that is the index set of available entries. Using this notation, the synchronization problem can be expressed as

$$\min_{X \in \Sigma^n} \epsilon(X) = \min_{X \in \Sigma^n} \| (Z - XX^{-\flat}) \circ \Omega \|_F^2 \iff (66)$$

$$\min_{\widetilde{Z}} \| (Z - \widetilde{Z}) \circ \Omega \|_F^2 \quad s.t. \ \widetilde{Z} = X X^{-\flat}, \ X \in \Sigma^n$$
(67)

where the problem of finding a consistent vertex labelling X is reduced to that of finding an edge labelling  $\widetilde{Z}$  induced by X.

If the *rank relaxation* is adopted (Arrigoni et al. 2014b), i.e. the matrix  $\widetilde{Z}$  is enforced to have rank (at most) *d* (while the remaining properties on  $\widetilde{Z}$  are not enforced), then Problem (67) becomes

$$\min_{\widetilde{Z}} \| (Z - \widetilde{Z}) \circ \Omega \|_F^2 \quad s.t. \ \operatorname{rank}(\widetilde{Z}) \le d \tag{68}$$

which is a *matrix completion* problem (Candès and Tao 2010), that is the problem of recovering a low-rank matrix starting from an incomplete subset of its entries (possibly corrupted by noise), which can be solved via (e.g.) the OPTSPACE algorithm (Keshavan et al. 2010).

In order to handle outliers, a *robust matrix completion* framework can be considered instead of (68), namely

$$\min_{\widetilde{Z},S} \| (Z - \widetilde{Z} - S) \circ \Omega \|_F^2$$
  
s.t. rank( $\widetilde{Z}$ )  $\leq d$ , S is sparse in  $\Omega$  (69)

where the additional variable *S* represents outliers, which are sparse over the measurement graph (by assumption). Available algorithms to solve problem (69) include R- GODEC (Arrigoni et al. 2014a), GRASTA (He et al. 2012) and L1-ALM (Zheng et al. 2012). This approach was introduced in Arrigoni et al. (2014a) for  $\Sigma = SO(d)$  and extended in Arrigoni et al. (2016b) to  $\Sigma = SE(d)$ .

In the  $\Sigma = S_d$  case, however, the optimization variable Z is sparse, being composed of binary matrices, and hence it does not satisfy the incoherence assumption (see Candès and Tao 2010) that make "generic" matrix completion algorithms work in practice. To overcome this drawback, the authors of Zhou et al. (2015) consider the following problem instead of (68)

$$\min_{\widetilde{Z}} \| (Z - \widetilde{Z}) \circ \Omega \|_{F}^{2} + \alpha \| \widetilde{Z} \|_{1}$$
s.t. rank $(\widetilde{Z}) \le d, \ 0 \le \widetilde{Z} \le 1$ 
(70)

where the regularization term  $\alpha \| \widetilde{Z} \|_1$  is included to promote a sparse solution, and the optimization variable is enforced to lie in the interval [0, 1] (while the binary constraints are not enforced). The resulting cost function is minimized via the Alternating Direction Method of Multipliers (ADMM) (Boyd et al. 2011).

## 7.3 Semidefinite Relaxation

Let us consider the  $\Sigma = O(d)$  case. Reasoning as in Eq. (67), we can express the synchronization problem as

$$\min_{\widetilde{Z}} \| (Z - \widetilde{Z}) \circ \Omega \|_F^2 \quad s.t. \ \widetilde{Z} = X X^{\mathsf{T}}, \ X \in \Sigma^n.$$
(71)

If the *semidefinite relaxation* is employed (Singer 2011; Singer and Shkolnisky 2011; Arie-Nachimson et al. 2012), i.e. the optimization variable  $\tilde{Z}$  is constrained to be symmetric positive semidefinite and covered by identity blocks along its diagonal (while the remaining properties on  $\tilde{Z}$  are not enforced), then Problem (71) reduces to a semidefinite program

$$\min_{\widetilde{Z}} \| (Z - \widetilde{Z}) \circ \Omega \|_F^2 \quad s.t. \ \widetilde{Z} \succeq 0, \ \widetilde{Z}_{ii} = I_d$$
(72)

which can be solved (e.g.) though interior point methods (Wright 1997). In Wang and Singer (2013) the  $\ell_1$ -norm is used in (72) in place of the  $\ell_2$ -norm, exploiting the fact that the former is more robust to outliers than the latter.

In order to deal with the  $\Sigma = S_d$  case, the authors of Chen et al. (2014) introduce in (72) the non-negative constraint  $\widetilde{Z} \ge 0$  and a sparsifying regularization term  $\alpha \|\widetilde{Z}\|_1$ , similarly to (70), resulting in the following problem

$$\min_{\widetilde{Z}} \| (Z - \widetilde{Z}) \circ \Omega \|_{F}^{2} + \alpha \| \widetilde{Z} \|_{1}$$
  
s.t.  $\widetilde{Z} \succeq 0, \ \widetilde{Z}_{ii} = I_{d}, \ \widetilde{Z} \ge 0$  (73)

which is solved via the ADMM algorithm.

Concerning the Special Euclidean Group, in Rosen et al. (2015) a cost function tightly related to (67) is considered and the feasible set SE(d) is relaxed to its convex hull, which admits a semidefinite representation (Saunderson et al. 2015). A convex relaxation is also employed in Rosen et al. (2019), Carlone et al. (2016) where the authors, using the theory of Lagrangian duality, develop an algorithm for certifying the global optimality of a candidate solution to rigid-motion synchronization.

It should be noted that although semidefinite relaxation solutions are not closed-form, they are nonetheless convex problems, which are relatively easy to compute.

# 8 Synchronization in Computer Vision

Among the several applications that have been mentioned in the introduction (image mosaicking, multi-view matching, clock synchronization, etc.) we concentrate here on the few ones where the formulation of the problem in terms of synchronization might require some clarification.



**Fig. 5** In the center, two nodes with partial visibility match their three common objects. At the extrema the ground truth ordering of the objects. Each node sees some of the objects (white circles are missing objects) and puts them in a different order, i.e., it gives them different numeric labels

## 8.1 Homography Synchronization

When elements of SL(d) are identified with homographies of the (d-1)-dimensional projective space, synchronization over SL(3) can be easily seen as a convenient way to align multiple images into a mosaic, starting from pairwise homographies. This works because any real non-singular  $3 \times 3$ matrix can be scaled to a *real* unit determinant matrix, as first mentioned in Schroeder et al. (2011). The same trick does not apply when *d* is even because of complex roots.

## 8.2 Permutation Synchronization

Consider a set of d objects, which is attached to each node in the measurement graph in a random order, i.e., each node has its own local labelling of the objects with integers in the range  $\{1, \ldots, d\}$ , represented as a permutation. It is assumed that pairs of nodes can match these objects, establishing which objects are the same in the two nodes, despite the different naming, and the goal is to infer a global labelling of the objects, such that the same object receives the same label in all the nodes (see Fig. 5).

A more concrete problem statement can be given in terms of *multi-view matching* (Chen et al. 2014; Zhou et al. 2015; Maset et al. 2017), where nodes are *images* and objects are *features*. A set of matches between pairs of images is computed in isolation, and the goal is to jointly update them so as to maximize their consistency. In general, not all the features are visible (or matchable) in all the images, so each matching is modelled as a partial permutation (Fig. 6).

Specifically, the permutation matrix P representing the matching between node B and node A is constructed as follows:

$$[P]_{h,k} = \begin{cases} 1 \text{ if obj. } k \in B \text{ is matched with obj. } h \in A; \\ 0 \text{ otherwise.} \end{cases}$$



Fig. 6 Partial permutation matrix representing the matches between two nodes

If row  $[P]_{h,.}$  contains only zeros, then object *h* in node A does not have a matching object in node B. If column  $[P]_{.,k}$  is made of zeros, then object *k* in node B does not have a matching object in node A.

Let  $Z_{ij}$  denote the permutation matrix representing the matching between node j and node i, and let  $X_i$  (resp.  $X_j$ ) denote the unknown permutation that reveals the true identity of the objects in node i (resp. j). It can be easily verified that  $Z_{ij} = X_i X_j^{\mathsf{T}}$ , thus the problem of finding the global labelling of the objects can be modelled as a synchronization problem over the Symmetric Inverse Semigroup  $\mathcal{I}_d$  (if permutations are partial), or over the Symmetric Group  $\mathcal{S}_d$  (if permutations are total).

## 8.3 Rigid-Motion Synchronization

Elements of SE(d) represent the angular attitude and position of a *d*-dimensional reference frame. These two properties are collectively referred to as *motion* in Computer Vision,

orientation in Photogrammetry, or pose in Robotic (although pose is also used in Vision). Synchronization over SE(d) is tantamount to recovering the location and attitude of a set of reference frames organized in a network, where the links of this network are relative transformations of one frame with respect to (some of) the others. This is also called *network* orientation (Fraser 20058), pose graph optimization (Carlone et al. 2015), or sensor network localization (Cucuringu et al. 2012). If we restrict the attention to the angular attitude (leaving out the position) then we get a rotation synchronization. Similarly, if position only is considered, it results in translation synchronization.

Such local frames can be local coordinates where 3D points are represented, in which case we are dealing with *multiple point-set registration* (Pulli 1999), or camera reference frames, in which case we are in the context of (global) *structure-from-motion* (Ozyesil et al. 2017).

In the first case, the goal is to find the rigid transformation that brings multiple (n > 2) 3D point sets into alignment. The problem can be solved in *point space* or in *frame space*. In the former case all the transformations are simultaneously optimized with respect to a cost function that depends on the distance between corresponding points (Pennec 1996; Benjemaa and Schmitt 1998; Krishnan et al. 2007; Toldo et al. 2010; Fantoni et al. 2012). In the latter case the optimization criterion is related to the internal coherence of the network of transformations applied to the local coordinate frames (Sharp et al. 2002; Fusiello et al. 2002; Torsello et al. 2011; Govindu and Pooja 2014; Arrigoni et al. 2016b). This is exactly an instance of rigid-motion synchronization, as shown in Fig. 7, where the input edge labels are typically computed via the Iterative Closest Point algorithm (Besl and McKay 1992).

In the structure from motion application the goal is to recover both scene structure (3D coordinates of scene points)



Fig. 7 The goal of the multiple point-set registration problem is to find the rigid transformations that bring multiple 3D point sets into alignment, where each rigid transformation is represented by a direct isometry



**Fig. 8** The epipolar graph: each node represents one camera and each edge is labelled with the essential matrix of the corresponding camera pair, which encodes the relative rotation and translation (up to scale)

and camera motion starting from a set of images. In this case G is known as the *epipolar graph* (Moulon et al. 2013; Arrigoni et al. 2015) or the *viewing graph* (Levi and Werman 2003) (see Fig. 8).

Structure from motion is a well studied problem that can be addressed in several ways [see the recent survey by Ozyesil et al. (2017)]. One stream of research focuses on methods that compute the motion *before* recovering the structure, and in this paper we are specifically interested in frame-space methods, that do not make use of points in order to compute the global motion. These methods are usually faster than sequential and hierarchical methods (e.g. Snavely et al. 2006; Schonberger and Frahm 2016; Toldo et al. 2015), while ensuring a fair distribution of the errors among the cameras, being global. On the other hand, accuracy is usually worse than that achieved by point-space methods, such as bundle adjustment.

We are interested in these "structure-free" methods because they reduce to a rigid-motion synchronization, modulo the fact that the relative translations are only known as directions, as the magnitude is unknown.

#### 8.3.1 Localization

We consider here the two-step formulation of synchronization over SE(3), described in Sect. 5.4. Since the magnitude of translations are unknown, we are required to estimate such magnitudes either directly or indirectly (i.e., by computing camera locations from the directions only). We will only hint at these solutions here, since they fall outside the domain of synchronization, although being related to it. As a matter of fact, the starting point is the translation synchronization equation

$$(B^{\dagger} \otimes I_3) \operatorname{vec}(X) = \operatorname{vec}(Z) \tag{74}$$

and from here two paths can be followed:

- First recover the *magnitude* of translations (Arrigoni et al. 2015; Tron et al. 2015) and then solve a *translation synchronization*;
- Solve the problem straight from the direction information: *bearing-only localization* (Brand et al. 2004; Zhao and Zelazo 2016).

Direct computation of magnitudes If the relative measures  $\mathbf{z}_{ij}$  are expanded into magnitude  $\alpha_{ij} = \|\mathbf{z}_{ij}\|$  and direction  $\mathbf{u}_{ij} = \mathbf{z}_{ij}/\|\mathbf{z}_{ij}\|$ , then the matrix *Z* rewrites

$$Z = U \operatorname{diag}(\boldsymbol{\alpha}) \tag{75}$$

where U denotes the  $3 \times m$  matrix obtained by juxtaposing all the edge directions, namely  $U = [\mathbf{u}_{12} \dots \mathbf{u}_{ij} \dots]$ , and  $\boldsymbol{\alpha} \in \mathbb{R}^m$  denotes the vector containing all the *unknown* magnitudes, namely  $\boldsymbol{\alpha} = [\alpha_{12} \dots \alpha_{ij} \dots]^{\mathsf{T}}$ . Thus, using Property (92), we get

$$\operatorname{vec}(Z) = (I_m \odot U)\boldsymbol{\alpha} \tag{76}$$

where  $\odot$  denotes the Khatri–Rao product (see "Appendix A"). Hence, Eq. (74) rewrites

$$(B^{\mathsf{T}} \otimes I_3) \operatorname{vec}(X) = (I_m \odot U) \boldsymbol{\alpha}.$$
(77)

This equation is also called the *edge-based bearing constraint* in Tron et al. (2015). Let us consider the cycle matrix  $C \in \{-1, 0, 1\}^{(m-n+1)\times m}$  associated with a cycle basis of  $\mathcal{G} = (\mathcal{V}, \mathcal{E})$  and let us multiply left and right sides in (77) by  $(C \otimes I_3)$ 

$$(C \otimes I_3)(B^{\mathsf{T}} \otimes I_3) \operatorname{vec}(X) = (C \otimes I_3)(I_m \odot U)\boldsymbol{\alpha}.$$
(78)

Using Properties (86), (91) and (108) we get

$$(CB^{\mathsf{T}} \otimes I_3) \operatorname{vec}(X) = \mathbf{0} = (C \odot U)\boldsymbol{\alpha}.$$
<sup>(79)</sup>

Note that Eq. (79) is nothing else but the cycle consistency (Eq. (17)) rewritten in terms of directions and magnitudes.

**Proposition 9** (Arrigoni et al. 2015; Tron et al. 2015) *The* unknown translation magnitudes can be uniquely (up to a global scale) recovered if and only if  $\operatorname{rank}(C \odot U) = m - 1$ . In this case the solution is given by the 1-dimensional nullspace of  $C \odot U$ .

*Bearing-only localization* Let us multiply (74) by the blockdiagonal matrix

$$S = \text{blkdiag}\left(\{[\mathbf{u}_{ij}]_{\times}\}_{(i,j)\in\mathcal{E}}\right)$$
(80)

	Localization		Synchronization
Direction:	Distance:	Translation:	Rotation
$\frac{\mathbf{x}_i - \mathbf{x}_j}{\ \mathbf{x}_i - \mathbf{x}_i\ }$	$\ \mathbf{x}_i - \mathbf{x}_j\ $	$\mathbf{x}_i - \mathbf{x}_j$	

Fig. 9 The relationship between synchronization and localization

where  $[\mathbf{u}_{ij}]_{\times}$  denotes the 3 × 3 skew-symmetric matrix corresponding to the cross-product with  $\mathbf{u}_{ii}$ , yielding

$$S(B^{\top} \otimes I_3) \operatorname{vec}(X) = S \operatorname{vec}(Z) = \mathbf{0}.$$
(81)

This step has the effect of substituting Z, which is unknown, with S (derived from U) which is known instead. Expanding this equation for a single edge (i, j) of the graph yields the more custom expression:

$$\mathbf{u}_{ij} \times (\mathbf{x}_i - \mathbf{x}_j) = \mathbf{0}. \tag{82}$$

This equation is also called the *node-based bearing constraint* in Tron et al. (2015). Its solution yields the node locations *X*, hence implicitly recovering the magnitudes. It is observed in Kennedy et al. (2012) that least squares solution of (81), which is used (e.g.) in Govindu (2001), is equivalent to the method presented in Brand et al. (2004). A bearingonly formulation is also adopted in Goldstein et al. (2016), Jiang et al. (2013), Moulon et al. (2013), Ozyesil and Singer (2015), Tron and Vidal (2014), Wilson and Snavely (2014) within the context of structure from motion.

This last section led us to glimpse the problem of *localization*. The goal of localization is to compute the position of *n* nodes in *d*-space given measures on the edges. As in a translation synchronization problem the nodes correspond to positions, i.e., elements of  $\mathbb{R}^d$ , but the available measures are *not* differences of states, i.e. translations, for they can be directions (bearings) or distances. Figure 9 hopefully helps to clarify this relationship.

The problem of localization from directions and the conditions under which an epipolar graph is *localizable* are discussed with continuity of notation by the same authors of this paper in Arrigoni and Fusiello (2019).

# 9 Discussion

In this section we report some considerations about the pros and cons of the closed-form solutions reviewed in this paper.

#### 9.1 Simplicity

These methods are particularly attractive for their simplicity. As an example we report the Matlab code of rotation and translation synchronization in 3-space (Listings 1 and 2).

```
Listing 1 Rotation synchronization
```

```
1
   function R = rotation_synch(Z,A)
2
     n = size(A,1);
3
     iD = diag(1./sum(A,2));
4
     [Q,~]=eigs( kron(iD, eye(3))*Z, 3);
5
6
     % Remove ambiguity
7
       = Q/(Q(1:3, 1:3));
     Ο
8
9
     % Projection onto SO(3)
10
     R = cell(1, n);
11
     for i=1:n
        [U,~,V] = svd(Q(3*i-2:3*i,:));
12
13
       R{i} = U*diag([1,1,det(U*V')])*V';
14
     end
15
   end
```

Listina 2	Translation	synchronization

1	func	tion T = translation_synch(U,B)
2		
3		B(1,:) = []; % remove node 1
4		F= <b>kron</b> (B′, <b>eye</b> (3));
5		$X = F \setminus U$ (:);
6		X=[0;0;0;X]; % add node 1
7		X= <b>reshape</b> (X,3,[]);
8		
9		<pre>T = num2cell(reshape(X,3,[]),</pre>
10		[1, <b>size</b> (X,2)]);
11	end	

In particular, rotation synchronization can be taken as a prototype of synchronization in any subgroup of GL(d): the only step that changes is the projection. In fact, the removal of the ambiguity with division by the first block works in all cases, since the identity belongs to any subgroup.

This does not hold for  $\mathcal{I}_d$ , which is not a subgroup of GL(d). In this case, removal of the ambiguity reduces to a clustering problem, as explained in Sect. 6. See also Tron et al. (2017) for the relation between clustering and partial permutation synchronization. Note that synchronization over  $\mathcal{I}_d$  requires to know the value of d in advance, which is not available in the multi-view matching application, since it corresponds to the total number of features (or tracks) present in the images. To overcome this drawback, a practical solution to multi-view matching is developed in Maset et al. (2017), where the authors compute refined edge labels (instead of vertex labels), in order to avoid the ambiguity issue.

#### 9.2 Spectral Solution Versus Null-Space Solution

In the presence of noise the null-space and spectral solutions do not coincide, in general. In the case of a complete graph, for instance, they coincide if and only if the unique non zero eigenvalue of Z is exactly n, which is unlikely to happen in practice.

An empirical comparison between the two approaches for the  $\Sigma = SE(3)$  case—is reported in Arrigoni et al. (2016c) where it is shown that the spectral solution achieves the same accuracy as the null-space method but it is faster. In particular, it turns out that the final rounding step, i.e. zeroing the imaginary part of the eigenvectors, do not compromise the accuracy achieved by the spectral method.

Note that in both cases the matrices inherit the same sparsity pattern as the adjacency matrix A, thus sparse solvers can be exploited, e.g., eigs for the spectral method and svds for the null-space solution, as done in Listing 1. As a matter of fact, svds (F) calls eigs ([0 F; F' 0]), as reported in the function documentation, and consequently it runs more slowly, for the dimension of the matrix is double. From the computational complexity point of view, every iteration of the Lanczos method (implemented by eigs) is linear in n(Golub and Van Loan 1996), if the matrix is sparse, but the number of iterations cannot be bounded by a constant.

# 9.3 Efficiency

The considered instances of synchronization translate into efficient closed form solutions, such as spectral decomposition or linear least squares. As a matter of fact, in the experiments reported in the literature (Arie-Nachimson et al. 2012; Bernard et al. 2015; Arrigoni et al. 2016c) they are consistently the fastest method. For instance, the rigid-motion synchronization pipeline in Arrigoni et al. (2016a) takes 7 s for the Madrid Metropolis dataset (Wilson and Snavely 2014), whose epipolar graph contains about 300 nodes and 65% of missing edges (with respect to the complete graph).

Note also that the spectral/null-space solutions aim at recovering the vertex labelling X, thus they require to store a  $dn \times d$  matrix in addition to the  $dn \times dn$  matrix  $Z_A$  containing the input measures, which is *sparse* if the graph is not complete. Semidefinite and rank relaxations, instead, aim at computing a refined edge label for each pair of nodes, represented by the optimization variable  $\tilde{Z}$  (see Sects. 7.2 and 7.3), thus they require an additional amount of memory to store a *full*  $dn \times dn$  matrix.

# 9.4 Accuracy

The least-squares solution to translation synchronization is statistically optimal, whereas the spectral (or, equivalently, null-space) solution to the other instances of synchronization provides an extrinsic estimate, whose quality is—in general—inferior to those provided by intrinsic methods. Note also that the spectral technique is a weaker relaxation than semidefinite programming (namely it enforces less constraints on the optimization variable), as explained in Sect. 7, and it does not guarantee exact recovery (in contrast to Wang and Singer 2013; Rosen et al. 2019; Carlone et al. 2016).

In Arrigoni et al. (2016c) the spectral method is compared to Rosen et al. (2015), which minimizes a geometric error tightly related to the synchronization cost function with a convex relaxation. The latter returns a more accurate solution, but it requires a significant amount of time (for instance, it takes around ten minutes on a synthetic scenario with n =100 and 30% of missing edges, whereas the spectral approach is able to compute a solution in less than 2 s with up to n = 1000).

Experiments on multiple point-set registration conducted in Arrigoni et al. (2018)-graphs with up to 95% of missing edges-demonstrate that the spectral solution and the semidefinite relaxation proposed in Rosen et al. (2019) are comparable in accuracy but the latter requires additional time. However, the same comparison<sup>4</sup> on the SLAM dataset used in Rosen et al. (2019)-with more than 99% of missing edgesreveals that (Rosen et al. 2019) achieves a running time comparable with the spectral method on these very sparse graphs. The authors Rosen and Carlone (2017) suggest that some computational enhancements can be incorporated in semidefinite relaxations in order to speed up the computation further. Finally, the same study performed in Arrigoni et al. (2018) indicates also that methods based on matrix completion (Arrigoni et al. 2014a), instead, fail when the percentage of missing data is higher than 90%.

Although the accuracy obtained by the spectral method is not optimal, it is however high, as demonstrated by experiments performed in Arrigoni et al. (2016c) in a variety of scenarios (for instance, it gets an average rotation error of the order of 1 degree on the popular Stanford 3D Scanning Repository<sup>5</sup>). As a consequence, it can be seen as a good and fast initialization for a subsequent local refinement (e.g. bundle adjustment in structure from motion).

### 9.5 Robustness

The least-squares solution for translation synchronization of Sect. 3 can be made resistant to outliers (i.e. rogue edge labels) by replacing  $\rho(y) = y^2$  in Eq. (14) with another function  $\rho(y)$  with sub-quadratic growth, and solving the resulting minimization problem, e.g., with Iteratively Reweighted Least Squares (IRLS) (Holland and Welsch 1977). This technique iteratively solves weighted least squares problems

<sup>&</sup>lt;sup>4</sup> Unpublished experiments made by the authors.

<sup>&</sup>lt;sup>5</sup> http://graphics.stanford.edu/data/3Dscanrep/.

where the weights are computed at each iteration as a function of the residuals of the current solution.

As for synchronization over GL(d), it is easy to see that the analysis of Sect. 4 can be extended to handle weighted measurements, which translates in relaxing the entries of the adjacency matrix A to assume real values in [0, 1], where 0 still indicates a missing measurement and the other values reflect the reliability of the edge labels. This allows to apply an IRLS-like scheme: first, an estimate for the vertex labelling with given edge weights is obtained via either the spectral approach or the null-space solution; then, the weights are updated based on the current estimate of the vertex labels, and these steps are iterated until convergence or a maximum number of iterations is reached.

Experiments in Arrigoni et al. (2016c) report good performances in the  $\Sigma = SE(d)$  case (resilience to about 40% of wrong edge labels) and it is shown in Arrigoni et al. (2018) that the spectral relaxation is comparable to methods performing robust matrix completion in terms of robustness to outliers.

### 9.6 Limitations and Open Questions

In summary, the spectral and null-space methods provide fast, sub-optimal (but fairly accurate) solutions in the presence of noise/outliers.

Further efficiency is needed in certain applications involving very large-scale problems, such as multi-view matching. For instance, the method in Maset et al. (2017), which provides a practical approach to partial permutation synchronization based on a spectral decomposition, took one and half hour to compute a solution in the Dino dataset (Seitz et al. 2006), where d = 493 and n = 363 (resulting in a matrix of size  $\approx 180000$ ). A possible way to speed-up the computation is to split the synchronization problem into smaller subproblems in a distributed fashion, similarly to (Cucuringu et al. 2012; Bhowmick et al. 2014).

Thanks to the group-labelling interpretation, the formalism of synchronization permits to address a variety of Computer Vision applications using the same Linear Algebra approach without relying on features or points, since the problem is formulated in frame space, or, more abstractly, in a group/semigroup. An unexplored application involves projective frames, which are represented by  $4 \times 4$  invertible matrices defined up to scale. As mentioned in Sect. 8.1, the idea of scaling each matrix by the *d*-th root of its determinant, which was used in Bartoli and Sturm (2003), does not apply here because complex roots may appear, hence further research is required in order to formulate the problem in terms of synchronization.

From the theoretical perspective, interesting open questions include the link between the spectral solution and the consistency error when the data matrix is not symmetric, and the relation between the notion of null-cycle and the synchronization problem over an inverse monoid.

# **10 Conclusion**

In this paper we gathered several disparate works within the common framework of synchronization and showed how this framework can be profitably used in several Computer Vision applications. Besides exhibiting a nice and clean formulation, synchronization can also benefit from efficient and closedform solutions such as spectral decomposition or linear least squares. We hope that this survey will serve as a starting point for more research in this field.

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# A Kronecker, Hadamard and Khatri–Rao Products

This appendix is devoted to the Kronecker, Hadamard and Khatri–Rao products (Van Loan 2000; Minka 2000; Liu and Trenkler 2008), which are widely used in this paper.

Let *A* and *B* be two real matrices of dimension  $m \times r$  and  $n \times s$  respectively. The *Kronecker product* of *A* and *B* (Van Loan 2000), denoted by  $A \otimes B$ , is defined as

$$A \otimes B = \begin{bmatrix} [A]_{1,1}B & [A]_{1,2}B \dots & [A]_{1,r}B \\ [A]_{2,1}B & [A]_{2,2}B \dots & [A]_{2,r}B \\ \dots & \dots & \dots \\ [A]_{m,1}B & [A]_{m,2}B \dots & [A]_{m,r}B \end{bmatrix}$$
(83)

where each  $[A]_{i,j} B$  is a block of dimension  $n \times s$ , thus  $A \otimes B$  has dimension  $mn \times rs$ . The Kronecker product is associative, distributive (with respect to the sum of matrices), but not commutative, and it satisfies the following properties

$$(A \otimes B)^{\mathsf{T}} = A^{\mathsf{T}} \otimes B^{\mathsf{T}} \tag{84}$$

$$(A \otimes B)^{-1} = A^{-1} \otimes B^{-1}$$
(85)

$$(A \otimes B)(C \otimes D) = (AC) \otimes (BD)$$
(86)

$$\operatorname{vec}(AXB) = (B^{\mathsf{T}} \otimes A)\operatorname{vec}(X)$$
 (87)

where  $vec(\cdot)$  denotes the vectorization operator which transforms a matrix into a vector by stacking the columns of the matrix one underneath the other.

Let  $A = U_A \Sigma_A V_A^{\mathsf{T}}$  and  $B = U_B \Sigma_B V_B^{\mathsf{T}}$  be the singular value decompositions of A and B, respectively, then

$$A \otimes B = (U_A \otimes U_B)(\Sigma_A \otimes \Sigma_B)(V_A \otimes V_B)^{\mathsf{T}}$$
(88)

which implies

$$\operatorname{rank}(A \otimes B) = \operatorname{rank}(A) \operatorname{rank}(B).$$
(89)

Thus the Kronecker product of two matrices is invertible if and only if both the factors are invertible.

Consider now two real matrices A and B of dimension  $m \times r$  and  $n \times r$  respectively, and denote the columns of A by  $\mathbf{a}_1, \ldots, \mathbf{a}_r$  and those of *B* by  $\mathbf{b}_1, \ldots, \mathbf{b}_r$ . The *Khatri–Rao* product of A and B (Khatri and Rao 1968; Liu and Trenkler 2008), denoted by  $A \odot B$ , is defined as

$$A \odot B = \begin{bmatrix} \mathbf{a}_1 \otimes \mathbf{b}_1 \ \mathbf{a}_2 \otimes \mathbf{b}_2 \cdots \mathbf{a}_r \otimes \mathbf{b}_r \end{bmatrix}$$
(90)

where each  $\mathbf{a}_i \otimes \mathbf{b}_i$  is a vector of dimension *mn*, thus  $A \odot B$ has dimension  $mn \times r$ . The Khatri–Rao product is associative, distributive, but not commutative, and it satisfies the following equalities

$$(A \otimes B)(C \odot D) = (AC) \odot (BD) \tag{91}$$

$$\operatorname{vec}(A \operatorname{diag}(\mathbf{x})B) = (B^{\top} \odot A)\mathbf{x}$$
(92)

where diag(**x**) transforms the vector  $\mathbf{x} = [x_1 \dots x_r]^T$  into a diagonal matrix with elements  $x_1, \ldots, x_r$  along the diagonal.

To the best of our knowledge, equalities expressing the rank of  $A \odot B$  in terms of the rank of the factors are not present in the literature, in contrast to the case of the Kronecker product. Some inequalities are reported in Khatri and Rao (1968), where it is shown, for instance, that  $rank(A \odot B) \ge rank(A)$ , if all the columns of B corresponding to independent columns of A are non-null.

Let A and B be two real matrices of dimension  $m \times r$ . The Hadamard product (or entry-wise product) of A and B (Minka 2000), denoted by  $A \circ B$ , has dimension  $m \times r$  as well, and it is simply the product of the corresponding elements

$$A \circ B = \begin{bmatrix} [A]_{1,1}[B]_{1,1} & [A]_{1,2}[B]_{1,2} & \dots & [A]_{1,r}[B]_{1,r} \\ [A]_{2,1}[B]_{2,1} & [A]_{2,2}[B]_{2,2} & \dots & [A]_{2,r}[B]_{2,r} \\ \dots & & \dots \\ [A]_{m,1}[B]_{m,1} & [A]_{m,2}[B]_{m,2} & \dots & [A]_{m,r}[B]_{m,r} \end{bmatrix}.$$
(93)

The Hadamard product is associative, distributive, commutative, and it satisfies the following properties

$$(A \odot B) \circ (C \odot D) = (A \circ C) \odot (B \circ D)$$
(94)

$$(A \odot B)^{\mathsf{T}} (A \odot B) = (AA^{\mathsf{T}}) \circ (BB^{\mathsf{T}})$$
(95)

$$\operatorname{vec}(A \circ B) = \operatorname{diag}(\operatorname{vec}(A)) \operatorname{vec}(B)$$
(96)  
$$\operatorname{diag}(\mathbf{x}) \land \operatorname{diag}(\mathbf{x}) = A \circ (\mathbf{x}\mathbf{y}^{\mathsf{T}})$$
(97)

$$\operatorname{rank}(A \circ B) < \operatorname{rank}(A) \operatorname{rank}(B).$$
(97)

$$\operatorname{rank}(A \circ B) \le \operatorname{rank}(A) \operatorname{rank}(B).$$
 (98)

## **B** Results from Graph Theory

In this section we review some useful concepts from graph theory. A complete treatment of this subject can be found in Bollobas (1998), Kavitha et al. (2009).

A graph is a pair  $\mathcal{G} = (\mathcal{V}, \mathcal{E})$  where  $\mathcal{V}$  is a finite set and  $\mathcal{E}$  is a family of pairs of elements of  $\mathcal{V}$ . We use *n* and *m* to denote the number of vertices and edges respectively, namely n = $|\mathcal{V}|$  and  $m = |\mathcal{E}|$ . A weighted graph is a graph together with a weight function  $\omega : \mathcal{E} \to \mathbb{R}^+$ . If the graph is unweighted, we set  $\omega : \mathcal{E} \to 1$  and call w the uniform weight function. An edge occurring more than once is referred to as a *multiple* edge, and a graph without multiple edges is called simple. An edge of the form (v, v) is called a *loop*. In an undirected graph, the *degree* of a vertex v is the number of times that v occurs as an endpoint of an edge. In a directed graph, the outdegree and indegree of a vertex v are the number of times that v occurs as the tail and head of an edge, respectively.

A subgraph  $\mathcal{G}' = (\mathcal{V}', \mathcal{E}')$  of  $\mathcal{G}$  is a graph with  $\mathcal{V}' \subseteq \mathcal{V}$ and  $\mathcal{E}' \subseteq \mathcal{E}$ . If  $\mathcal{E}'$  is a subset of  $\mathcal{E}$ , then  $\mathcal{G} \setminus \mathcal{E}'$  denotes the graph obtained by removing all the edges in  $\mathcal{E}'$  from  $\mathcal{G}$ . If  $\mathcal{V}'$  is a subset of  $\mathcal{V}$ , then  $\mathcal{G} \setminus \mathcal{V}'$  denotes the graph obtained by removing all the vertices in  $\mathcal{V}'$  and their incident edges from  $\mathcal{G}$ . A *path* from v to w is a subgraph  $\mathcal{G}' = (\mathcal{V}', \mathcal{E}')$  with  $\mathcal{V}' = \{v_0 = v, v_1, \dots, v_k = w\}$ and  $\mathcal{E}' = \{(v_0, v_1), (v_1, v_2), \dots, (v_{k-1}, v_k)\}$ . An undirected graph is called *connected* if there exists a path from each vertex to any other, and a directed graph is called connected if the underlying undirected graph is connected. Any maximal connected subgraph  $\mathcal{H}$  is called a *connected component*. A graph is a *tree* if it is connected and it has n - 1 edges. The disjoint union of trees is called a forest. The number of edges in a forest is n - cc, where cc denotes the number of connected components in  $\mathcal{G}$ . A subgraph  $\mathcal{G}'$  of a connected graph  $\mathcal{G}$  is called a *spanning tree* if it has the same vertices of  $\mathcal{G}$  and it is a tree. If G is not connected, any union of spanning trees for each connected component is called a spanning forest.

A *cycle* in a graph  $\mathcal{G}$  is a vector  $\mathbf{c} \in \mathbb{Q}^m$  such that for any vertex  $v \in \mathcal{V}$  it holds

$$\sum_{e \in \delta_+(v)} [\mathbf{c}]_e = \sum_{e \in \delta_-(v)} [\mathbf{c}]_e$$
(99)

where  $\delta_+(v)$  and  $\delta_-(v)$  denote the edges leaving and entering v, respectively, and  $[\mathbf{c}]_e$  denotes the component of **c** indexed by edge *e*. A cycle is *simple* if  $[\mathbf{c}]_e \in \{-1, 0, 1\}$  for all  $e \in \mathcal{E}$ , and a simple cycle is a *circuit* if its support (i.e. the set of edges with  $[\mathbf{c}]_e \neq 0$ ) is connected and for any vertex  $v \in \mathcal{V}$ there are at most two edges in the support incident to v. The set of cycles forms a vector space over  $\mathbb{Q}$ , which is called the *cycle space* of  $\mathcal{G}$ , and a *cycle basis* is a set of circuits forming a basis of such a space. It can be shown (Bollobas 1998; Kavitha et al. 2009) that if  $\mathcal{G}$  is connected the dimension of the cycle space is given by the *cyclomatic number* 

$$\nu = m - n + 1. \tag{100}$$

## **B.1 Matrices Associated with Graphs**

The *adjacency matrix* A of a graph G is the  $n \times n$  matrix whose elements indicate whether pairs of vertices are adjacent or not, namely

$$[A]_{i,j} = \begin{cases} 1 & \text{if } (i,j) \in \mathcal{E} \\ 0 & \text{otherwise.} \end{cases}$$
(101)

If  $\mathcal{G}$  does not contain loops, then A has zero diagonal. Note that the adjacency matrix is symmetric if the graph is undirected.

The *incidence matrix B* of a directed graph G is the  $n \times m$  matrix defined by

$$[B]_{k,e} = \begin{cases} -1 & \text{if } k \text{ is the head of edge } e, \\ 1 & \text{if } k \text{ is the tail of edge } e, \\ 0 & \text{otherwise.} \end{cases}$$
(102)

The rows of *B* correspond to vertices and the columns correspond to edges. Note that each column has exactly two non zero entries, which correspond to the endpoints of the edge associated to that column. The *incidence matrix B* of an undirected graph  $\mathcal{G}$  is defined considering a particular orientation of the edges. It is shown in Bollobas (1998) that, if  $\mathcal{G}$  is connected, then

$$\operatorname{rank}(B) = n - 1. \tag{103}$$

The *degree matrix* D of an undirected graph G is the  $n \times n$  diagonal matrix such that  $[D]_{i,i}$  contains the degree of node i. Equivalently, it can be defined as

$$D = \operatorname{diag}(A\mathbf{1}_{n \times 1}) \tag{104}$$

where  $\mathbf{1}_{n \times 1}$  denotes a  $n \times 1$  matrix filled by ones, thus  $A\mathbf{1}_{n \times 1}$  is the sum of the rows of A. In the case of a directed graph, either the indegree or the outdegree can be used. The *transition matrix* P is defined as

$$P = D^{-1}A. (105)$$

The Laplacian matrix L of a graph  $\mathcal{G}$  is defined as

$$L = D - A. \tag{106}$$

It can be checked that, independently of the orientation of the edges, the following equation holds

$$L = BB^{\mathsf{T}} \tag{107}$$

which implies that *L* is symmetric and positive semidefinite, and, if the graph is connected,  $\operatorname{rank}(L) = \operatorname{rank}(B) = n - 1$ , hence *L* is singular. Note that the vector  $\mathbf{1}_{n \times 1}$  is in the nullspace of *L*.

The notion of adjacency matrix can be extended to the case of a weighted graph, which translates in letting the entries of A to assume values in [0, 1]. Specifically,  $[A]_{i,j}$  contains the weight of edge (i, j), and  $[A]_{i,j} = 0$  still indicates that  $(i, j) \notin \mathcal{E}$ . In this case Eqs. (104), (105) and (106) still make sense, which define the degree matrix, the transition matrix and the Laplacian matrix of a weighted graph, respectively.

The *cycle matrix C* corresponding to a cycle basis of a connected graph  $\mathcal{G}$  is the  $(m - n + 1) \times m$  matrix having the incidence vectors of the circuits in the basis in its rows. The following equation (Bollobas 1998) expresses the relation between the cycle matrix and the incidence matrix

$$CB^{\mathsf{T}} = 0. \tag{108}$$

Note that, if the graph is undirected, the matrices A and L are symmetric, thus their eigenvalues are real. The matrix P is not symmetric, but it is similar to the symmetric matrix N defined as

$$N = D^{-1/2} A D^{-1/2} (109)$$

since  $P = D^{-1/2}ND^{1/2}$ . The matrices N and P have the same eigenvalues, so P has real eigenvalues.

**Theorem 1** (Perron–Frobenius (Meyer 2000)) If an  $n \times n$ matrix has non-negative entries then it has a non-negative real eigenvalue  $\lambda$  which has maximum absolute value among all the eigenvalues. This eigenvalue has a non-negative real eigenvector. If, in addition, the matrix has no block-triangular decomposition, then  $\lambda$  has multiplicity 1 and the corresponding eigenvector is positive.

As explained in Lovász (2007), the Perron–Frobenius theorem implies that, if  $\mathcal{G}$  is connected, the largest eigenvalue of *A* has multiplicity 1. Likewise, the largest eigenvalue of the transition matrix is 1 and it ha multiplicity 1. It is easy to check that the eigenvector associated to such eigenvalue is  $\mathbf{1}_{n \times 1}$ .

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