

Solving Quadratically Constrained Geometrical Problems using Lagrangian Duality

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Abstract

In this paper we consider the problem of solving different pose and registration problems under rotational constraints. Traditionally, methods such as the iterative closest point algorithm have been used to solve these problems. They may however get stuck in local minima due to the non-convexity of the problem. In recent years methods for finding the global optimum, based on Branch and Bound convex under-estimators, have been developed. These methods are provably optimal, however since they are based on global optimization methods they are in general more time consuming than local methods.

In this paper we adopt a dual approach. Rather than trying to find the globally optimal solution we investigate how good solutions it is possible to obtain using Lagrange duality. Our approach allows us to formulate a single convex semidefinite program that approximates the original problem well. Furthermore, we show that it is possible to obtain bounds on the global optimum from this program using standard duality theory.¹

1 Geometric Registration

A frequently occurring and by now a classical problem in computer vision, robotic manipulation and photogrammetry is the registration problem. That is, finding the transformation between two coordinate systems, see [5] and the references therein. The problem appears in several contexts: relating two stereo reconstructions, solving the hand-eye calibration problem and finding the absolute pose of an object given 3D measurements.

There are a number of solutions proposed and perhaps the most well-known is by Horn *et al* [6]. They

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derive a closed-form solution for the Euclidean (or similarity) transformation that minimizes the sum of squares error between the transformed points and the measured points.

The more general problem of finding the registration between two 3-D shapes was considered in [2], where the *iterative closest point* (ICP) algorithm was proposed to solve the problem. The algorithm is able to cope with different geometric primitives, like point sets, line segments and different kinds of surface representations. However, the algorithm requires a good initial transformation in order to converge to the globally optimal solution, otherwise only a local optimum is obtained.

In [8], the method of Horn *et al* [6] was generalized by incorporating point, line and plane features in a common framework. Given *point-to-point*, *point-to-line*, or *point-to-plane* correspondences, it was demonstrated how the transformation (Euclidean or similarity) relating the two coordinate systems can be computed based on a geometrically meaningful cost-function. The algorithm was based on relaxing the non-convex problem by convex under-estimators and then using branch and bound to focus in on the global solution [1].

In this paper we adopt a dual approach. Rather than trying to find the globally optimal solution we investigate how good solutions it is possible to obtain using Lagrange duality. Our approach allows us to formulate a single convex semidefinite program that approximates the original problem well. We show on both synthetic and real data that the approximation is very close to the global optimum, in particular if the noise level is low. Furthermore, we show that it is possible to obtain bounds on the global optimum from this program using standard duality theory.

Semidefinite programming has previously been used for solving large scale binary quadratic optimization problems [7]. These problems are often NP-complete, however due to a result in [4], semidefinite programming have been shown to produce good approximations while still being a polynomial algorithm. On the other hand for very large problems, polynomial execu-

tion time might still be too much. In the vision community it has mostly been used for large scale combinatorial problems resulting in long execution times. However, since the matrices that arise in the applications that we consider in this paper are quite small (typically less than 100×100), it provides an efficient way to produce lower estimates and approximative solutions.

2 Lagrangian Duality

In its most general form, the problems we are interested in can be described as geometric registration problems. The objective is to relate measurements in one coordinate frame to an object model in another. In this section we consider the case where the measurements are 3D-points and the model consists of planes in 3D-space. The goal is to find a Euclidian transformation $(R, t) \in SO(3) \times \mathbb{R}^3$ that places the measurement point x_i as close to the corresponding model plane π_i as possible. It was shown in [8], that given a number of planes π_i in one coordinate system and points x_i in another, $i = 1, \dots, m$ this problem can be formulated as the optimization problem

$$\min_{\substack{R \in SO(3) \\ t \in \mathbb{R}^3}} \sum_{i=1}^m d(Rx_i + t, \pi_i), \quad (1)$$

where $d(x, \pi)$ denotes the distance between point x and the plane π . If we let y_i be an arbitrary point on π_i and n_i be the unit normal, then we can rewrite the problem as

$$\min_{\substack{R \in SO(3) \\ t \in \mathbb{R}^3}} \sum_{i=1}^m \|n_i^T (Rx_i + t - y_i)\|_2^2. \quad (2)$$

If we disregard the constraint $R \in SO(3)$ this is a linear least squares problem in the unknowns R and t . Note that since set of feasible matrices R is larger if this constraint is disregarded it is easy to see that the least squares solution provides a lower bound on the minimum value of the original problem. However since $R \in SO(3)$ we also have $R^T R = I$ giving the following quadratically constrained problem

$$\begin{aligned} \min \quad & \sum_{i=1}^m \|n_i^T (Rx_i + t - y_i)\|_2^2 \\ \text{s.t.} \quad & R^T R - I = 0 \end{aligned} \quad (3)$$

Since $R^T R - I$ is symmetric, equation (4) consists of 6 quadratic constraints. It is well known that problems with no more than 2 quadratic constraints can be solved in polynomial time. Hence we are forced to search for approximative solutions. Next we rewrite the problem in vector form. We let $v^T = [r_{11} \ r_{12} \ \dots \ r_{33} \ t_1 \ t_2 \ t_3]$. The objective function is quadratic and (3) can therefore be written

$$\min v^T A v + 2b^T v, \quad (5)$$

where A and b are determined from the model and measurement data. Note that $A \succeq 0$. By introducing the 12×3 -block matrices $E_1 = [I \ 0 \ 0 \ 0]$, $E_2 = [0 \ I \ 0 \ 0]$ and $E_3 = [0 \ 0 \ I \ 0]$ we may rewrite the problem as

$$\min \quad v^T A v + 2b^T v \quad (6)$$

$$\text{s.t.} \quad v^T E_i^T E_j v = \delta_{ij} \quad (7)$$

where $\delta_{ij} = 1$ if $i = j$ and 0 otherwise. As we mentioned earlier the problem is easily solved if the constraints are dropped. This gives us a lower bound on the objective value. We may modify the objective function by adding $\mu_{ij}(v^T E_i^T E_j v - \delta_{ij})$ without changing the original problem. For each different μ_{ij} we may obtain a different relaxation. Lagrangian duality can be viewed as finding the relaxation that gives the largest objective value.

The Lagrangian of (6) becomes

$$\begin{aligned} L(v, \mu) = v^T A v + 2b^T v + \sum_{i,j} \mu_{ij} (v^T E_i^T E_j v - \delta_{ij}) = \\ v^T \left(A + \sum_i \mu_{ij} E_{ij} \right) v + 2b^T v - \sum_{i,j} \mu_{ij} \delta_{ij}, \end{aligned} \quad (8)$$

where $E_{ij} = \frac{1}{2}(E_i^T E_j + E_j^T E_i)$. Thus finding the largest relaxation can be written as

$$\max_{\mu} \min_v L(v, \mu) \quad (9)$$

For a fixed μ the inner minimization is only finite if the matrix $A + \sum_{i,j} \lambda_{ij} E_{ij}$ is positive semidefinite. In this case the minimizer can be computed analytically by taking the gradient with respect to v , which yields

$$v = -(A + \sum_{i,j} \lambda_{ij} E_{ij})^{-1} b \quad (10)$$

Inserting into (8) we see that (9) can be written

$$\begin{aligned} \max_{\mu} \quad & -b^T \left(A + \sum_{i,j} \mu_{ij} E_{ij} \right)^{-1} b + \sum_{i,j} \delta_{ij} \mu_{ij} \\ \text{s.t.} \quad & A + \sum_{i,j} \mu_{ij} E_{ij} \succeq 0 \end{aligned} \quad (11)$$

The above problem is the dual program when the primal variables have been eliminated. It can be shown that the objective function is concave and the constraint is convex, hence this program can be solved efficiently. However to be able to use a standard solver, such as SeDuMi [9] we need to write it as a standard linear semidefinite program. This can be done using the Schur-complement (see [3]). First we add the artificial variable γ and rewrite the program as

$$\max_{\gamma, \mu} \quad \gamma \quad (12)$$

$$\text{s.t.} \quad A + \sum_{i,j} \mu_{ij} E_{ij} \succeq 0 \quad (13)$$

$$-b^T (A + \sum_{i,j} \mu_{ij} E_{ij})^{-1} b + \sum_{i,j} \delta_{ij} \mu_{ij} - \gamma \geq 0 \quad (14)$$

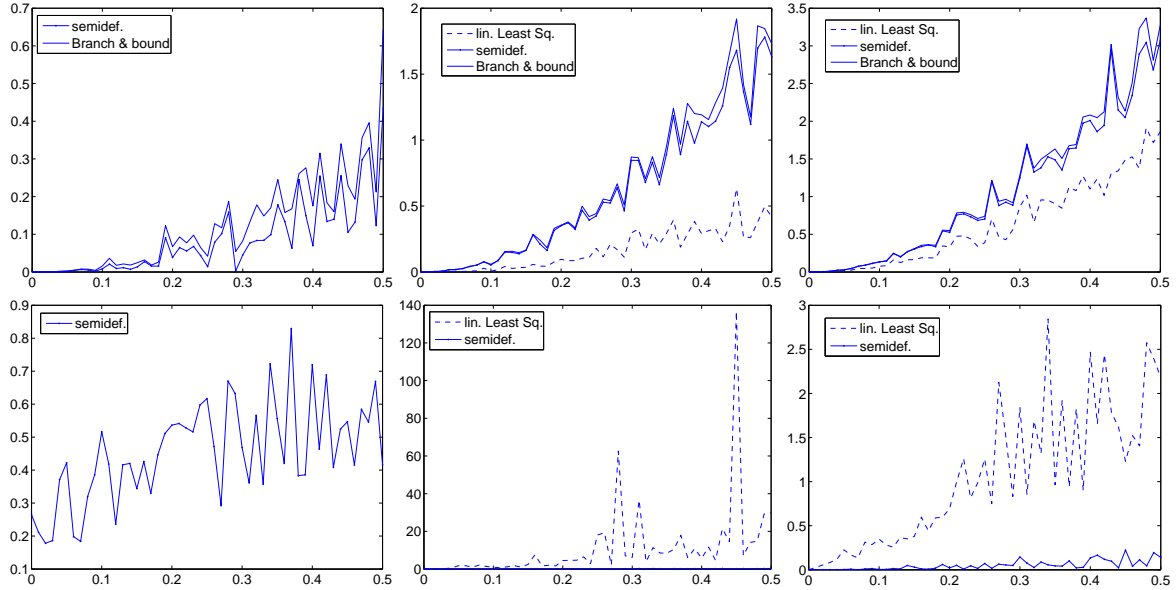


Figure 1: Top: The objective value of the obtained solution as a function of the noise level for three different problem sizes, $m = 7, 14$ and 20 . For comparison we also plot the global minimum, given by the Branch & Bound method, and the Linear Least Squares solution. Bottom: The value of $\|R^T R - I\|_2^2$ for the obtained solution as a function of the noise level.

Let

$$A(\mu, \gamma) = \begin{bmatrix} \sum_{i,j} \delta_{ij} \mu_{ij} - \gamma & b^T \\ b & A + \sum_{i,j} \mu_{ij} E_{ij} \end{bmatrix} \quad (15)$$

Using the Schur complement [3], the dual problem can finally be written as a linear semidefinite program

$$\max_{\gamma, \mu} \quad \gamma \quad (16)$$

$$\text{s.t.} \quad A(\mu, \gamma) \succeq 0. \quad (17)$$

This program is usually referred to the dual semidefinite program. Using the same technique as above it is possible to derive the primal semidefinite program

$$\min_{V \succeq 0} \quad \text{tr} \left(\begin{bmatrix} A & b \\ b^T & 0 \end{bmatrix} V \right) \quad (18)$$

$$\text{s.t.} \quad \text{tr}(E_{ij} V) = \delta_{ij} \quad (19)$$

$$\text{tr} \left(\begin{bmatrix} 0 & 0 \\ 0 & 1 \end{bmatrix} V \right) = 1. \quad (20)$$

Note that if V^* solves the primal problem and V^* can be written $V^* = v^*(v^*)^T$ then v^* is the global optimum of problem (6).

When using a primal-dual interior point as SeDuMi one obtains both a solution to the primal and the dual programs. To generate good approximate solutions we will take the eigenvector corresponding to the largest eigenvalue of the primal solution V as our approximate solution. If V is close to a rank one matrix this should be a good approximation.

3 Experiments

In this section we present a few experiments that shows that even though the program we have derived is a relaxation it provides surprisingly good solutions, indicating that the lagrangian relaxation is strong.

In our first experiment we use synthetic data to test the performance for various noise levels. The data was created by randomly placing m planes such that they intersect the sphere centered around zero with radius 10. Then we randomly selected one point from each plane and added noise of standard deviation σ . The first row of figure 1 shows the objective value of the obtained solution for different noise levels and different problem sizes. The relaxation can be seen to approximate the global optimum well, particularly for low noise level. Also it is a considerable improvement compared to the linear least squares solution. Note that for the case of 7 point-to-plane correspondences there is not enough data to compute a least squares solution. Also note that as the number of measurements increases the least squares solution will approach the correct solution. This is to be expected since the measurements are corrupted with normally distributed noise. The second row of figure 1 shows the value of $\|R^T R - I\|_2^2$ for the approximate solutions. The semidefinite solution appears to be quite close to being a rotation matrix while the linear least squares is far off, although as expected its performance improves as the number of measurements increases. Note that the y-scale is not

the same in all pictures.

For our next experiment we used real data. Figure 2 shows the setup for this experiment. We used a MicroScribe-3DLX 3d scanner to measure the 3D-coordinates of some points on two different objects. The 3D-scanner consists of a pointing arm with five degrees of freedom which measures 3D-coordinates. In total we measured 49 points on the toy model visible in figure 2. By request of the designer we will refer to the toy model as the space station. We created a computer

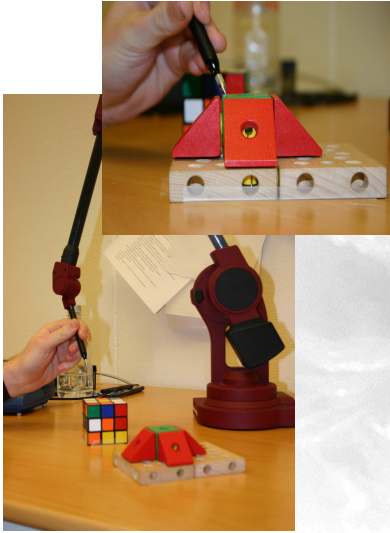


Figure 2: The experimental setup for the spacestation experiment.

model consisting of planes, lines and points (see figure 3). Among the measured points 27 were known to belong to a certain plane in the model, 12 to a line and 10 to point. Note that we only considered point-to-plane correspondences in the previous section, however it is easy to extend to other correspondences as well. Figure 3 shows the model and the resulting registration. The points marked with black crosses are points measured on the planes, the points marked with green rings are measured on lines and the points marked with red stars are measured on corners. In table 1 we show the total error for the different correspondences. For comparison we have also included the errors obtained with the other methods. In this case the solution of the primal semidefinite program has rank one and therefore finds the optimal solution (same as Branch & Bound). Note that Horn's method only measures the point-to-point errors resulting in a higher total error.

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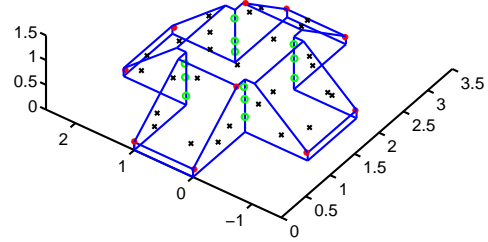


Figure 3: The model of the spacestation and the optimal registration. In this case both the branch and bound method and the semidefinite method gives the same solution.

Residuals:	B & B	Horn	Least Sq.	Semidef
point-point	0.0083	0.0063	0.0221	0.0083
point-line	0.0018	0.0036	0.0015	0.0018
point-plane	0.0046	0.0098	0.0046	0.0046
Total	0.0147	0.0197	0.0282	0.0147

Table 1: Resulting reconstruction errors for the space station problem when using the different methods. In this case the semidefinite program gives the same solution as the branch and bound method.

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