

ITERATIVE MINIMIZATION SCHEMES FOR SOLVING THE SINGLE SOURCE LOCALIZATION PROBLEM*

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Abstract. We consider the problem of locating a single radiating source from several noisy measurements using a maximum likelihood (ML) criteria. The resulting optimization problem is nonconvex and nonsmooth, and thus finding its global solution is in principle a hard task. Exploiting the special structure of the objective function, we introduce and analyze two iterative schemes for solving this problem. The first algorithm is a very simple explicit fixed-point-based formula, and the second is based on solving at each iteration a nonlinear least squares problem, which can be solved globally and efficiently after transforming it into an equivalent quadratic minimization problem with a single quadratic constraint. We show that the nonsmoothness of the problem can be avoided by choosing a specific “good” starting point for both algorithms, and we prove the convergence of the two schemes to stationary points. We present empirical results that support the underlying theoretical analysis and suggest that, despite of its nonconvexity, the ML problem can effectively be solved globally using the devised schemes.

Key words. single source location problem, Weiszfeld algorithm, nonsmooth and nonconvex minimization, fixed-point methods, nonlinear least squares, generalized trust region, semidefinite relaxation

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1. Introduction.

1.1. The source localization problem. Consider the problem of locating a single radiating source from noisy range measurements collected using a network of passive sensors. More precisely, consider an array of m sensors, and let $\mathbf{a}_j \in \mathbb{R}^n$ denote the coordinates of the j th sensor.¹ Let $\mathbf{x} \in \mathbb{R}^n$ denote the unknown source’s coordinate vector, and let $d_j > 0$ be a noisy observation of the range between the source and the j th sensor:

$$(1.1) \quad d_j = \|\mathbf{x} - \mathbf{a}_j\| + \varepsilon_j, \quad j = 1, \dots, m,$$

where $\boldsymbol{\varepsilon} = (\varepsilon_1, \dots, \varepsilon_m)^T$ denotes the unknown noise vector. Such observations can be obtained, for example, from the time-of-arrival measurements in a constant-velocity propagation medium. The source localization problem is the following.

The source localization problem: Given the observed range measurements $d_j > 0$, find a “good” approximation of the source \mathbf{x} satisfying (1.1).

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¹In practical applications $n = 2$ or 3 .

The source localization problem has received significant attention in the signal processing literature and specifically in the field of mobile phones localization [12, 5, 13]. It is also worth mentioning that the interest in wireless localization problems has increased since the first ruling of the Federal Communications Commission (FCC) for the detection of emergency calls in the United States in 1996 [17]. Currently, a high percentage of Enhanced 911 (E911) calls originate from mobile phones. Due to the unknown location of the wireless E911 calls, these calls do not receive the same quality of emergency assistance that fixed network 911 calls enjoy. To deal with this problem, the FCC issued an order on July 12, 1996, requiring all wireless service providers to report accurate mobile station location information to the E911 operator.

In addition to emergency management, mobile position information is also useful in mobile advertising, asset tracking, fleet management, location-sensitive billing [12], interactive map consultation, and monitoring of the mentally impaired [5].

1.2. The maximum likelihood criteria. In this paper we adopt the maximum-likelihood (ML) approach for solving the source localization problem (1.1); see, e.g., [4]. When ϵ follows a Gaussian distribution with a covariance matrix proportional to the identity matrix, the source \mathbf{x} is the ML estimate that is the solution of the problem:

$$(1.2) \quad (\text{ML}): \quad \min_{\mathbf{x} \in \mathbb{R}^n} \left\{ f(\mathbf{x}) \equiv \sum_{j=1}^m (\|\mathbf{x} - \mathbf{a}_j\| - d_j)^2 \right\}.$$

Note that, in addition to the statistical interpretation, the latter problem is a least squares problem in the sense that it minimizes the squared sum of the errors.

An alternative approach for estimating the source location \mathbf{x} is by solving the following least squares (LS) problem in the squared domain:

$$(1.3) \quad (\text{LS}): \quad \min_{\mathbf{x} \in \mathbb{R}^n} \sum_{j=1}^m (\|\mathbf{x} - \mathbf{a}_j\|^2 - d_j^2)^2.$$

Despite of its nonconvexity, the LS problem can be solved globally and efficiently by transforming it into a problem of minimizing a quadratic function subject to a single quadratic constraint [1] (more details will be given in section 3.2). However, the LS approach has two major disadvantages compared to the ML approach: first, the LS formulation lacks the statistical interpretation of the ML problem. Second, as demonstrated by the numerical simulations in section 4, the LS estimate provides less accurate solutions than those provided by the the ML approach.

The ML problem, like the LS problem, is nonconvex. However, as opposed to the LS problem for which a global solution can be computed efficiently [1], the ML problem seems to be a difficult problem to solve efficiently. A possible reason for the increased difficulty of the ML problem is its nonsmoothness. One approach for approximating the solution of the ML problem is via semidefinite relaxation (SDR) [4, 1]. We also note that the source localization problem formulated as (ML) can be viewed as a special instance of sensor network localization problems in which several sources are present; see, for example, the recent work in [3]; for this class of problems, semidefinite programming-based algorithms have been developed.

In this paper we depart from the SDR techniques and seek other efficient approaches to solve the ML problem. This is achieved by exploiting the special structure of the objective function which allows us to devise fixed-point-based iterative schemes

for solving the nonsmooth and nonconvex ML problem (1.2). The first scheme admits a very simple explicit iteration formula given by

$$\mathbf{x}^{k+1} = \mathcal{M}_1(\mathbf{x}^k, \mathbf{a}) \quad (\text{where } \mathbf{a} \equiv (\mathbf{a}_1, \dots, \mathbf{a}_m)),$$

while the second iterative scheme is of the form

$$\mathbf{x}^{k+1} \in \underset{\mathbf{x}}{\operatorname{argmin}} \mathcal{M}_2(\mathbf{x}, \mathbf{x}^k, \mathbf{a})$$

and requires the solution of an additional subproblem which will be shown to be efficiently solved. The main goals of this paper are to introduce the building mechanism of these two schemes, to develop and analyze their convergence properties, and to demonstrate their computational viability for solving the ML problem (1.2), as well as their effectiveness when compared with the LS and SDR approaches.

1.3. Paper layout. In the next section, we present and analyze the first scheme, which is a simple fixed-point-based method. The second algorithm, which is based on solving a sequence of least squares problems of a similar structure to that of (1.3), is presented and analyzed in section 3. The construction of both methods is motivated by two different interpretations of the well-known Weiszfeld method for the Fermat–Weber location problem [16]. For both schemes, we show that the nonsmoothness of the problem can be avoided by choosing a specific “good” starting point. Empirical results presented in section 4 provide a comparison between the two devised algorithms, as well as a comparison to different approaches such as LS and SDR. In particular, the numerical results suggest that, despite its nonconvexity, the ML problem can, for all practical purposes, be globally solved using the devised methods.

1.4. Notation. Throughout the paper, the following notation is used: vectors are denoted by boldface lowercase letters, e.g., \mathbf{y} , and matrices by boldface uppercase letters, e.g., \mathbf{A} . The i th component of a vector \mathbf{y} is written as y_i . Given two matrices \mathbf{A} and \mathbf{B} , $\mathbf{A} \succ \mathbf{B}$ ($\mathbf{A} \succeq \mathbf{B}$) means that $\mathbf{A} - \mathbf{B}$ is positive definite (semidefinite). The directional derivative of a function $f : \mathbb{R}^n \rightarrow \mathbb{R}$ at $\bar{\mathbf{x}}$ in the direction \mathbf{v} is defined (if it exists) by

$$(1.4) \quad f'(\mathbf{x}; \mathbf{v}) \equiv \lim_{t \rightarrow 0^+} \frac{f(\bar{\mathbf{x}} + t\mathbf{v}) - f(\bar{\mathbf{x}})}{t}.$$

The α -level set of a function $f : \mathbb{R}^n \rightarrow \mathbb{R}$ is defined by $\operatorname{Lev}(f, \alpha) = \{\mathbf{x} \in \mathbb{R}^n : f(\mathbf{x}) \leq \alpha\}$. The collection of m sensors $\{\mathbf{a}_1, \dots, \mathbf{a}_m\}$ is denoted by \mathcal{A} .

2. A simple fixed-point algorithm. In this section we introduce a simple fixed-point algorithm that is designed to solve the ML problem (1.2). The algorithm is inspired by the celebrated Weiszfeld algorithm for the Fermat–Weber problem, which is briefly recalled in section 2.1. In section 2.2 we introduce and analyze the fixed-point scheme designed to solve the ML problem.

2.1. A small detour: Weiszfeld algorithm for the Fermat–Weber problem. As was already mentioned, the ML problem (1.2) is nonconvex and nonsmooth, and thus finding its exact solution is in principle a difficult task. We propose a fixed-point scheme motivated by the celebrated Weiszfeld algorithm [16, 7] for solving the Fermat–Weber location problem:

$$(2.1) \quad \min_{\mathbf{x}} \left\{ s(\mathbf{x}) \equiv \sum_{j=1}^m \omega_j \|\mathbf{x} - \mathbf{a}_j\| \right\},$$

where $\omega_j > 0$ and $\mathbf{a}_j \in \mathbb{R}^n$ for $j = 1, \dots, m$. Of course, the Fermat–Weber problem is much easier to analyze and solve than the ML problem (1.2) since it is a well-structured nonsmooth convex minimization problem. This problem has been extensively studied in the location theory literature; see, for instance, [11]. Our objective here is to mimic the Weiszfeld algorithm [16] to obtain an algorithm for solving the nonsmooth and nonconvex ML problem (1.2). The Weiszfeld method is a very simple fixed-point scheme that is designed to solve the Fermat–Weber problem. One way to derive it is to write the first order global optimality conditions for the convex problem (2.1)

$$\nabla s(\mathbf{x}) = \sum_{j=1}^m \omega_j \frac{\mathbf{x} - \mathbf{a}_j}{\|\mathbf{x} - \mathbf{a}_j\|} = 0 \quad \forall \mathbf{x} \notin \mathcal{A}$$

as

$$\mathbf{x} = \frac{\sum_{j=1}^m \omega_j \frac{\mathbf{a}_j}{\|\mathbf{x} - \mathbf{a}_j\|}}{\sum_{j=1}^m \frac{\omega_j}{\|\mathbf{x} - \mathbf{a}_j\|}},$$

which naturally calls for the iterative scheme

$$(2.2) \quad \mathbf{x}^{k+1} = \frac{\sum_{j=1}^m \omega_j \frac{\mathbf{a}_j}{\|\mathbf{x}^k - \mathbf{a}_j\|}}{\sum_{j=1}^m \frac{\omega_j}{\|\mathbf{x}^k - \mathbf{a}_j\|}}.$$

For the convergence analysis of the Weiszfeld algorithm (2.2) and modified versions of the algorithm, see, e.g., [10, 15], and references therein.

2.2. The simple fixed-point algorithm: Definition and analysis. Similarly to the Weiszfeld method, our starting point for constructing a fixed-point algorithm to solve the ML problem is by writing the optimality conditions. Assuming that $\mathbf{x} \notin \mathcal{A}$ we have that \mathbf{x} is a stationary point for problem (ML) if and only if

$$(2.3) \quad \nabla f(\mathbf{x}) = 2 \sum_{j=1}^m (\|\mathbf{x} - \mathbf{a}_j\| - d_j) \frac{\mathbf{x} - \mathbf{a}_j}{\|\mathbf{x} - \mathbf{a}_j\|} = \mathbf{0},$$

which can be written as

$$\mathbf{x} = \frac{1}{m} \left\{ \sum_{j=1}^m \mathbf{a}_j + \sum_{j=1}^m d_j \frac{\mathbf{x} - \mathbf{a}_j}{\|\mathbf{x} - \mathbf{a}_j\|} \right\}.$$

The latter relation calls for the following fixed-point algorithm, which we term the *standard fixed point (SFP) scheme*.

ALGORITHM SFP.

$$(2.4) \quad \mathbf{x}^{k+1} = \frac{1}{m} \left\{ \sum_{j=1}^m \mathbf{a}_j + \sum_{j=1}^m d_j \frac{\mathbf{x}^k - \mathbf{a}_j}{\|\mathbf{x}^k - \mathbf{a}_j\|} \right\}, \quad k \geq 0.$$

Like in the Weiszfeld algorithm, the SFP scheme is not well defined if $\mathbf{x}^k \in \mathcal{A}$ for some k . In what follows we will show that by carefully selecting the initial vector \mathbf{x}^0

we can *guarantee* that the iterates are not in the sensors set \mathcal{A} , therefore establishing that the method is well defined. At this juncture, it is interesting to notice that the approach we suggest here for dealing with the points of nonsmoothness that occur at $\mathbf{x}^k \in \mathcal{A}$ is quite different from the common approaches for handling the nonsmoothness. For example, in order to avoid the nondifferentiable points of the Fermat–Weber objective function, several modifications of the Weiszfeld method were proposed; see, e.g., [10, 15], and references therein. However, there do not seem to have been any attempts in the literature to choose good initial starting points to avoid the nonsmoothness difficulty. A constructive procedure for choosing a good starting point for the SFP method will be given at the end of this section.

Before proceeding with the analysis of the SFP method, we record the fact that, much like the Weiszfeld algorithm (see [7]), the SFP scheme is a gradient method with a fixed step size.

PROPOSITION 2.1. *Let $\{\mathbf{x}^k\}$ be the sequence generated by the SFP method (2.4), and suppose that $\mathbf{x}^k \notin \mathcal{A}$ for all $k \geq 0$. Then*

$$(2.5) \quad \mathbf{x}^{k+1} = \mathbf{x}^k - \frac{1}{2m} \nabla f(\mathbf{x}^k).$$

Proof. The proof follows by a straightforward calculation, using the gradient of f computed in (2.3). \square

A gradient method does not necessarily converge without additional assumptions (e.g., assuming that ∇f is Lipschitz continuous and/or using a line search [2]). Nevertheless, we show below that scheme (2.4) *does* converge.

By Proposition 2.1 the SFP method can be compactly written as

$$(2.6) \quad \mathbf{x}^{k+1} = T(\mathbf{x}^k),$$

where $T : \mathbb{R}^n \setminus \mathcal{A} \rightarrow \mathbb{R}^n$ is the operator defined by

$$(2.7) \quad T(\mathbf{x}) = \mathbf{x} - \frac{1}{2m} \nabla f(\mathbf{x}).$$

In the convergence analysis of the SFP method, we will also make use of the auxiliary function:

$$(2.8) \quad h(\mathbf{x}, \mathbf{y}) \equiv \sum_{j=1}^m \|\mathbf{x} - \mathbf{a}_j - d_j r_j(\mathbf{y})\|^2 \quad \forall \mathbf{x} \in \mathbb{R}^n, \mathbf{y} \in \mathbb{R}^n \setminus \mathcal{A},$$

where

$$r_j(\mathbf{y}) \equiv \frac{\mathbf{y} - \mathbf{a}_j}{\|\mathbf{y} - \mathbf{a}_j\|}, \quad j = 1, \dots, m.$$

Note that for every $\mathbf{y} \notin \mathcal{A}$, the following relations hold for every $j = 1, \dots, m$:

$$(2.9) \quad \|r_j(\mathbf{y})\| = 1,$$

$$(2.10) \quad (\mathbf{y} - \mathbf{a}_j)^T r_j(\mathbf{y}) = \|\mathbf{y} - \mathbf{a}_j\|.$$

In Lemma 2.1 below, we prove several key properties of the auxiliary function h defined in (2.8).

LEMMA 2.1.

- (a) $h(\mathbf{x}, \mathbf{x}) = f(\mathbf{x})$ for every $\mathbf{x} \notin \mathcal{A}$.
- (b) $h(\mathbf{x}, \mathbf{y}) \geq f(\mathbf{x})$ for every $\mathbf{x} \in \mathbb{R}^n, \mathbf{y} \in \mathbb{R}^n \setminus \mathcal{A}$.
- (c) If $\mathbf{y} \notin \mathcal{A}$, then

$$(2.11) \quad T(\mathbf{y}) = \operatorname{argmin}_{\mathbf{x} \in \mathbb{R}^n} h(\mathbf{x}, \mathbf{y}).$$

Proof. (a) For every $\mathbf{x} \notin \mathcal{A}$,

$$\begin{aligned} f(\mathbf{x}) &= \sum_{j=1}^m (\|\mathbf{x} - \mathbf{a}_j\| - d_j)^2 \\ &= \sum_{j=1}^m (\|\mathbf{x} - \mathbf{a}_j\|^2 - 2d_j\|\mathbf{x} - \mathbf{a}_j\| + d_j^2) \\ &\stackrel{(2.9), (2.10)}{=} \sum_{j=1}^m (\|\mathbf{x} - \mathbf{a}_j\|^2 - 2d_j(\mathbf{x} - \mathbf{a}_j)^T r_j(\mathbf{x}) + d_j^2 \|r_j(\mathbf{x})\|^2) = h(\mathbf{x}, \mathbf{x}), \end{aligned}$$

where the last equation follows from (2.8).

(b) Using the definition of f and h given in (1.2) and (2.8), respectively, and the fact (2.9), a short computation shows that for every $\mathbf{x} \in \mathbb{R}^n, \mathbf{y} \in \mathbb{R}^n \setminus \mathcal{A}$,

$$\begin{aligned} h(\mathbf{x}, \mathbf{y}) - f(\mathbf{x}) &= 2 \sum_{j=1}^m d_j (\|\mathbf{x} - \mathbf{a}_j\| - (\mathbf{x} - \mathbf{a}_j)^T r_j(\mathbf{y})) \\ &\geq 0, \end{aligned}$$

where the last inequality follows from the Cauchy–Schwarz inequality and using again (2.9).

(c) For any $\mathbf{y} \in \mathbb{R}^n \setminus \mathcal{A}$, the function $\mathbf{x} \mapsto h(\mathbf{x}, \mathbf{y})$ is strictly convex on \mathbb{R}^n and consequently admits a unique minimizer \mathbf{x}^* satisfying

$$\nabla_{\mathbf{x}} h(\mathbf{x}^*, \mathbf{y}) = \mathbf{0}.$$

Using the definition of h given in (2.8), the latter identity can be explicitly written as

$$\sum_{j=1}^m (\mathbf{x}^* - \mathbf{a}_j - d_j r_j(\mathbf{y})) = \mathbf{0},$$

which by simple algebraic manipulation can be shown to be equivalent to $\mathbf{x}^* = \mathbf{y} - \frac{1}{2m} \nabla f(\mathbf{y})$, establishing that $\mathbf{x}^* = T(\mathbf{y})$. \square

Using Lemma 2.1 we are now able to prove the monotonicity property of the operator T with respect to f .

LEMMA 2.2. *Let $\mathbf{y} \notin \mathcal{A}$. Then*

$$f(T(\mathbf{y})) \leq f(\mathbf{y}),$$

and equality holds if and only if $T(\mathbf{y}) = \mathbf{y}$.

Proof. By (2.11) and the strict convexity of the function $\mathbf{x} \mapsto h(\mathbf{x}, \mathbf{y})$, one has

$$h(T(\mathbf{y}), \mathbf{y}) < h(\mathbf{x}, \mathbf{y}) \text{ for every } \mathbf{x} \neq T(\mathbf{y}).$$

In particular, if $T(\mathbf{y}) \neq \mathbf{y}$, then

$$(2.12) \quad h(T(\mathbf{y}), \mathbf{y}) < h(\mathbf{y}, \mathbf{y}) = f(\mathbf{y}),$$

where the last equality follows from Lemma 2.1(a). By Lemma 2.1(b), $h(T(\mathbf{y}), \mathbf{y}) \geq f(T(\mathbf{y}))$, which, combined with (2.12), establishes the desired strict monotonicity. \square

Theorem 2.1 given below states the basic convergence results for the SFP method. In the proof, we exploit the boundedness of the level sets of the objective function f , which is recorded in the following lemma.

LEMMA 2.3. *The level sets of f are bounded.*

Proof. The proof follows immediately from the fact that $f(\mathbf{x}) \rightarrow \infty$ as $\|\mathbf{x}\| \rightarrow \infty$. \square

THEOREM 2.1 (convergence of the SFP method). *Let $\{\mathbf{x}^k\}$ be generated by (2.4) such that \mathbf{x}^0 satisfies*

$$(2.13) \quad f(\mathbf{x}^0) < \min_{j=1, \dots, m} f(\mathbf{a}_j).$$

Then

- (a) $\mathbf{x}^k \notin \mathcal{A}$ for every $k \geq 0$;
- (b) for every $k \geq 0$, $f(\mathbf{x}^{k+1}) \leq f(\mathbf{x}^k)$, and equality is satisfied if and only if $\mathbf{x}^{k+1} = \mathbf{x}^k$.
- (c) the sequence of function values $\{f(\mathbf{x}^k)\}$ converges;
- (d) the sequence $\{\mathbf{x}^k\}$ is bounded;
- (e) every convergent subsequence $\{\mathbf{x}^{k_l}\}$ satisfies $\mathbf{x}^{k_l+1} - \mathbf{x}^{k_l} \rightarrow \mathbf{0}$;
- (f) any limit point of $\{\mathbf{x}^k\}$ is a stationary point of f .

Proof. (a) and (b) The proof follows by induction on k using Lemma 2.2.

(c) The proof readily follows from the monotonicity and lower boundedness (by zero) of the sequence $\{f(\mathbf{x}^k)\}$.

(d) By (b), all of the iterates \mathbf{x}^k are in the level set $\text{Lev}(f, f(\mathbf{x}^0))$ which, by Lemma 2.3, establishes the boundedness of the sequence $\{\mathbf{x}^k\}$.

(e) and (f) Let $\{\mathbf{x}^{k_l}\}$ be a convergent subsequence of $\{\mathbf{x}^k\}$ with limit point \mathbf{x}^* . Since $f(\mathbf{x}^{k_l}) \leq f(\mathbf{x}^0) < \min_{j=1, \dots, m} f(\mathbf{a}_j)$, it follows by the continuity of f that $f(\mathbf{x}^*) \leq f(\mathbf{x}^0) < \min_{j=1, \dots, m} f(\mathbf{a}_j)$, proving that $\mathbf{x}^* \notin \mathcal{A}$. By (2.6)

$$(2.14) \quad \mathbf{x}^{k_l+1} = T(\mathbf{x}^{k_l}).$$

Therefore, since the subsequence $\{\mathbf{x}^{k_l}\}$ and its limit point \mathbf{x}^* are not in \mathcal{A} , by the continuity of ∇f on $\mathbb{R}^n \setminus \mathcal{A}$, we conclude that the subsequence $\{\mathbf{x}^{k_l+1}\}$ converges to a vector $\bar{\mathbf{x}}$ satisfying

$$(2.15) \quad \bar{\mathbf{x}} = T(\mathbf{x}^*).$$

To prove (e), we need to show that $\bar{\mathbf{x}} = \mathbf{x}^*$. Since both \mathbf{x}^* and $\bar{\mathbf{x}}$ are limit points of $\{\mathbf{x}^k\}$ and since the sequence of function values converges (by (c)), then the continuity of f over \mathbb{R}^n implies that $f(\mathbf{x}^*) = f(\bar{\mathbf{x}})$. Invoking Lemma 2.2 for $\mathbf{y} = \mathbf{x}^*$, we conclude that $\bar{\mathbf{x}} = \mathbf{x}^*$, proving claim (e). Part (f) follows from the observation that the equality $\mathbf{x}^* = T(\mathbf{x}^*)$ is equivalent (by the definition of T) to $\nabla f(\mathbf{x}^*) = \mathbf{0}$. \square

Remark 2.1. It is easy to find a vector \mathbf{x}^0 satisfying condition (2.13). For example, Procedure INIT, that will be described at the end of this section, produces a point satisfying (2.13).

Combining claims (c) and (f) of Theorem 2.1, we immediately obtain convergence of the sequence of function values.

COROLLARY 2.1. *Let $\{\mathbf{x}^k\}$ be the sequence generated by the SFP algorithm satisfying (2.13). Then $f(\mathbf{x}^k) \rightarrow f^*$, where f^* is the function value at a stationary point of f .*

We were able to prove the convergence of the function values of the sequence. The situation is more complicated for the sequence itself, where we were able only to show that all limit points are stationary points. We can prove convergence of the sequence itself if we assume that all stationary points of the objective function are isolated.² The proof of this claim strongly relies on the following lemma from [8].

LEMMA 2.4 (see [8, Lemma 4.10]). *Let \mathbf{x}^* be an isolated limit point of a sequence $\{\mathbf{x}^k\}$ in \mathbb{R}^n . If $\{\mathbf{x}^k\}$ does not converge, then there is a subsequence $\{\mathbf{x}^{k_i}\}$ which converges to \mathbf{x}^* and an $\epsilon > 0$ such that $\|\mathbf{x}^{k_i+1} - \mathbf{x}^{k_i}\| \geq \epsilon$.*

We can now use the above lemma to prove a convergence result under the assumption that all stationary points of f are isolated.

THEOREM 2.2 (convergence of the sequence). *Let $\{\mathbf{x}^k\}$ be generated by (2.4) such that \mathbf{x}^0 satisfies (2.13). Suppose further that all stationary points of f are isolated. Then the sequence $\{\mathbf{x}^k\}$ converges to a stationary point.*

Proof. Let \mathbf{x}^* be a limit point of $\{\mathbf{x}^k\}$ (its existence follows from the boundedness of the sequence proved in Theorem 2.1(d)). By our assumption \mathbf{x}^* is an isolated point. Suppose in contradiction that the sequence does not converge. Then by Lemma 2.4 there exists a subsequence $\{\mathbf{x}^{k_i}\}$ that converges to \mathbf{x}^* satisfying $\|\mathbf{x}^{k_i+1} - \mathbf{x}^{k_i}\| \geq \epsilon$. However, this is in contradiction to (e) of Theorem 2.1. We thus conclude that $\{\mathbf{x}^k\}$ converges to a stationary point. \square

The analysis of the SFP method relies on the validity of condition (2.13) on the starting point \mathbf{x}^0 . We will now show that, thanks to the special structure of the objective function (ML), we can compute such a point through a simple procedure. This is achieved by establishing the following result.

LEMMA 2.5. *Let $\mathcal{A} \equiv \{\mathbf{a}_1, \dots, \mathbf{a}_m\}$ be the given set of m sensors, and let*

$$g_j(\mathbf{x}) = \sum_{i=1, i \neq j}^m (\|\mathbf{x} - \mathbf{a}_i\| - d_i)^2, \quad j = 1, \dots, m.$$

Then for every $j = 1, \dots, m$ the following apply:

- (i) *If $\nabla g_j(\mathbf{a}_j) \neq \mathbf{0}$, then $f'(\mathbf{a}_j; -\nabla g_j(\mathbf{a}_j)) < 0$. Otherwise, if $\nabla g_j(\mathbf{a}_j) = \mathbf{0}$, then $f'(\mathbf{a}_j; \mathbf{v}) < 0$ for every $\mathbf{v} \neq \mathbf{0}$. In particular, there exists a descent direction from every sensor point.*
- (ii) *Every $\bar{\mathbf{x}} \in \mathcal{A}$ is not a local optimum for the ML problem (1.2).*

Proof. (i) For convenience, for every $j = 1, \dots, m$ we denote

$$(2.16) \quad f_j(\mathbf{x}) = (\|\mathbf{x} - \mathbf{a}_j\| - d_j)^2$$

so that the objective function of problem (ML) can be written as

$$(2.17) \quad f(\mathbf{x}) = f_j(\mathbf{x}) + g_j(\mathbf{x})$$

for every $\mathbf{x} \in \mathbb{R}^n$ and $j = 1, \dots, m$. Note that f is not differentiable for every $\mathbf{x} \in \mathcal{A}$. Nonetheless, the directional derivative of f at \mathbf{x} in the direction $\mathbf{v} \in \mathbb{R}^n$ always exists

²We say that \mathbf{x}^* is an isolated stationary point of f , if there are no other stationary points in some neighborhood of \mathbf{x}^* .

and is given by

$$(2.18) \quad f'(\bar{\mathbf{x}}; \mathbf{v}) = \begin{cases} \nabla f(\bar{\mathbf{x}})^T \mathbf{v}, & \bar{\mathbf{x}} \notin \mathcal{A}, \\ \nabla g_j(\mathbf{a}_j)^T \mathbf{v} - 2d_j \|\mathbf{v}\|, & \bar{\mathbf{x}} = \mathbf{a}_j. \end{cases}$$

Indeed, the above formula for $\bar{\mathbf{x}} \notin \mathcal{A}$ is obvious. In the other case, suppose then that $\bar{\mathbf{x}} = \mathbf{a}_j$ for some $j \in \{1, \dots, m\}$. Noting that g_j is differentiable at \mathbf{a}_j , we have $g'_j(\mathbf{a}_j; \mathbf{v}) = \nabla g_j(\mathbf{a}_j)^T \mathbf{v}$, and using definition (2.16) for f_j , we get $f'_j(\mathbf{a}_j; \mathbf{v}) = -2d_j \|\mathbf{v}\|$, and hence with (2.17), we obtain the desired formula (2.18) for $f'(\mathbf{a}_j; \mathbf{v})$. Finally, if $\nabla g_j(\mathbf{a}_j) \neq \mathbf{0}$, then using (2.18) we have

$$f'(\mathbf{a}_j; -\nabla g_j(\mathbf{a}_j)) = -\|\nabla g_j(\mathbf{a}_j)\|^2 - 2d_j \|\nabla g_j(\mathbf{a}_j)\| < 0.$$

Otherwise, if $\nabla g_j(\mathbf{a}_j) = \mathbf{0}$, then for every $\mathbf{v} \neq \mathbf{0}$ we have

$$f'(\mathbf{a}_j; \mathbf{v}) = -2d_j \|\mathbf{v}\| < 0.$$

(ii) By part (i) there exists a descent direction from every sensor point $\bar{\mathbf{x}} \in \mathcal{A}$. Therefore, none of the sensor points can be a local optimum for problem (ML). \square

Using the descent directions provided by Lemma 2.5, we can compute a point $\bar{\mathbf{x}}$ satisfying

$$f(\bar{\mathbf{x}}) < \min_{j=1, \dots, m} f(\mathbf{a}_j)$$

by the following procedure.

PROCEDURE INIT.

1. $t = 1$.
2. **Set** k to be an index for which $f(\mathbf{a}_k) = \min_{j=1, \dots, m} f(\mathbf{a}_j)$.
3. **Set**

$$(2.19) \quad \mathbf{v}_0 = \begin{cases} -\nabla g_k(\mathbf{a}_k), & \nabla g_k(\mathbf{a}_k) \neq \mathbf{0}, \\ \mathbf{e}, & \nabla g_k(\mathbf{a}_k) = \mathbf{0}, \end{cases}$$

where \mathbf{e} is the vector of all ones.³

4. **While** $f(\mathbf{a}_k + t\mathbf{v}_0) \geq f(\mathbf{a}_k)$, **set** $t = t/2$. **End**
5. The output of the algorithm is $\mathbf{a}_k + t\mathbf{v}_0$.

The validity of this procedure stems from the fact that, by Lemma 2.5, the direction \mathbf{v}_0 defined in (2.19) is always a descent direction.

One of the advantages of the SFP scheme is its simplicity. However, the SFP method, being a gradient method, does have the tendency to converge to local minima. In the next section we will present a second and more involved algorithm to solve the ML problem. As we shall see in the numerical examples presented in section 4, the empirical performance of this second iterative scheme is significantly better than that of the SFP, both with respect to the number of required iterations and with respect to the probability of getting stuck in a local/nonglobal point.

3. A sequential weighted least squares algorithm. In this section we study a different method for solving the ML problem (1.2), which we call the sequential weighted least squares (SWLS) algorithm. The SWLS algorithm is also motivated by the construction of the Weiszfeld method, but from a different viewpoint; see section 3.1. Each iteration of the method consists of solving a nonlinear least squares problem, whose solution is found by the approach discussed in section 3.2. The convergence analysis of the SWLS algorithm is given in section 3.3.

³We could have chosen any other nonzero vector.

3.1. The SWLS algorithm. To motivate the SWLS algorithm, let us first go back to the Weiszfeld scheme for solving the classical Fermat–Weber location problem, whereby we rewrite the iterative scheme (2.2) in the following equivalent, but different, way:

$$(3.1) \quad \mathbf{x}^{k+1} = \operatorname{argmin}_{\mathbf{x} \in \mathbb{R}^n} \left\{ \sum_{j=1}^m \omega_j \frac{\|\mathbf{x} - \mathbf{a}_j\|^2}{\|\mathbf{x}^k - \mathbf{a}_j\|} \right\}.$$

The strong convexity of the objective function in (3.1) (recall that $\omega_j > 0$ for all j) implies that \mathbf{x}^{k+1} is uniquely defined as a function of \mathbf{x}^k . Therefore, the Weiszfeld method (2.2) for solving problem (2.1) can also be written as

$$\mathbf{x}^{k+1} = \operatorname{argmin}_{\mathbf{x} \in \mathbb{R}^n} q(\mathbf{x}, \mathbf{x}^k),$$

where

$$q(\mathbf{x}, \mathbf{y}) \equiv \sum_{j=1}^m \omega_j \frac{\|\mathbf{x} - \mathbf{a}_j\|^2}{\|\mathbf{y} - \mathbf{a}_j\|} \text{ for every } \mathbf{x} \in \mathbb{R}^n, \mathbf{y} \in \mathbb{R}^n \setminus \mathcal{A}.$$

The auxiliary function q was essentially constructed from the objective function s of the Fermat–Weber location problem, by replacing the norm terms $\|\mathbf{x} - \mathbf{a}_j\|$ with $\frac{\|\mathbf{x} - \mathbf{a}_j\|^2}{\|\mathbf{y} - \mathbf{a}_j\|}$, i.e., with $s(\mathbf{x}) = q(\mathbf{x}, \mathbf{x})$. Mimicking this observation for the ML problem under study, we will use an auxiliary function in which each norm term $\|\mathbf{x} - \mathbf{a}_j\|$ in the objective function (1.2) is replaced with $\frac{\|\mathbf{x} - \mathbf{a}_j\|^2}{\|\mathbf{y} - \mathbf{a}_j\|}$, resulting in the following auxiliary function:

$$(3.2) \quad g(\mathbf{x}, \mathbf{y}) \equiv \sum_{i=1}^m \left(\frac{\|\mathbf{x} - \mathbf{a}_i\|^2}{\|\mathbf{y} - \mathbf{a}_i\|} - d_i \right)^2, \quad \mathbf{x} \in \mathbb{R}^n, \mathbf{y} \in \mathbb{R}^n \setminus \mathcal{A}.$$

The general step of the algorithm for solving problem (ML), the SWLS method, is now given by

$$\mathbf{x}^{k+1} \in \operatorname{argmin}_{\mathbf{x} \in \mathbb{R}^n} g(\mathbf{x}, \mathbf{x}^k)$$

or more explicitly by the following algorithm.

ALGORITHM SWLS.

$$(3.3) \quad \mathbf{x}^{k+1} \in \operatorname{argmin}_{\mathbf{x} \in \mathbb{R}^n} \sum_{j=1}^m \left(\frac{\|\mathbf{x} - \mathbf{a}_j\|^2}{\|\mathbf{x}^k - \mathbf{a}_j\|} - d_j \right)^2.$$

The name SWLS stems from the fact that at each iteration k we are required to solve the following weighted least squares (WLS) version of the LS problem (1.3):

$$(3.4) \quad (\text{WLS}): \quad \min_{\mathbf{x}} \sum_{j=1}^m \omega_j^k \left(\|\mathbf{x} - \mathbf{c}_j\|^2 - \beta_j^k \right)^2,$$

with

$$(3.5) \quad \mathbf{c}_j = \mathbf{a}_j, \quad \beta_j^k = d_j \|\mathbf{x}^k - \mathbf{a}_j\|, \quad \omega_j^k = \frac{1}{\|\mathbf{x}^k - \mathbf{a}_j\|^2}.$$

Note that the SWLS algorithm as presented above is not defined for iterations in which $\mathbf{x}^k \in \mathcal{A}$. In our random numerical experiments (cf. section 4) this situation never occurred; i.e., \mathbf{x}^k did not belong to \mathcal{A} for every k . However, from a theoretical point of view this issue must be resolved. Similarly to the methodology advocated in the convergence analysis of the SFP method, our approach for avoiding the sensor points \mathcal{A} is by choosing a “good enough” initial vector. In section 3.3, we introduce a simple condition on the initial vector \mathbf{x}^0 under which the algorithm is well defined and proven to converge.

3.2. Solving the WLS subproblem. We will now show how the WLS subproblem (3.4) can be solved globally and efficiently by transforming it into a problem of minimizing a quadratic function subject to a single quadratic constraint. This derivation is a straightforward extension of the solution technique devised in [1] and is briefly described here for completeness.

For a given fixed k (for simplicity we omit the index k below), we first transform (3.4) into a constrained minimization problem:

$$(3.6) \quad \min_{\mathbf{x} \in \mathbb{R}^n, \alpha \in \mathbb{R}} \left\{ \sum_{j=1}^m \omega_j (\alpha - 2\mathbf{c}_j^T \mathbf{x} + \|\mathbf{c}_j\|^2 - \beta_j)^2 : \|\mathbf{x}\|^2 = \alpha \right\},$$

which can also be written as (using the substitution $\mathbf{y} = (\mathbf{x}^T, \alpha)^T$)

$$(3.7) \quad \min_{\mathbf{y} \in \mathbb{R}^{n+1}} \left\{ \|\mathbf{A}\mathbf{y} - \mathbf{b}\|^2 : \mathbf{y}^T \mathbf{D}\mathbf{y} + 2\mathbf{f}^T \mathbf{y} = 0 \right\},$$

where

$$\mathbf{A} = \begin{pmatrix} -2\sqrt{\omega_1} \mathbf{c}_1^T & \sqrt{\omega_1} \\ \vdots & \vdots \\ -2\sqrt{\omega_m} \mathbf{c}_m^T & \sqrt{\omega_m} \end{pmatrix}, \quad \mathbf{b} = \begin{pmatrix} \sqrt{\omega_1} (\beta_1 - \|\mathbf{c}_1\|^2) \\ \vdots \\ \sqrt{\omega_m} (\beta_m - \|\mathbf{c}_m\|^2) \end{pmatrix}$$

and

$$\mathbf{D} = \begin{pmatrix} \mathbf{I}_n & \mathbf{0}_{n \times 1} \\ \mathbf{0}_{1 \times n} & 0 \end{pmatrix}, \quad \mathbf{f} = \begin{pmatrix} 0 \\ -0.5 \end{pmatrix}.$$

Note that (3.7) belongs to the class of problems consisting of minimizing a quadratic function subject to a single quadratic constraint. Problems of this type are called generalized trust region subproblems (GTRS). GTRS problems possess necessary and sufficient optimality conditions from which efficient solution methods can be derived; see, e.g., [6, 9].

The SWLS scheme is of course more involved than the simpler SFP scheme. However, as explained above, the additional computations required in SWLS to solve the subproblem can be done efficiently and are worthwhile, since the SWLS algorithm usually possesses a much larger region of convergence to the global minimum than the SFP scheme, which in turn implies that it has the tendency of avoiding local minima and a greater chance of hitting the global minimum. This will be demonstrated on the numerical examples given in section 4.

TABLE 1
 Number of runs (out of 10000) for which Assumption 2 is satisfied for $\mathbf{x}^0 = \mathbf{x}_{\text{LS}}$.

σ	1e-3	1e-2	1e-1	1e+0
N_σ	10000	10000	9927	6281

3.3. Convergence analysis of the SWLS method. In this section we provide an analysis of the SWLS method. We begin by presenting our underlying assumptions in section 3.3.1, and in section 3.3.2 we prove the convergence results of the method.

3.3.1. Underlying assumptions. The following assumption will be made throughout this section.

Assumption 1. The matrix

$$\mathbf{A} = \begin{pmatrix} 1 & \mathbf{a}_1^T \\ 1 & \mathbf{a}_2^T \\ \vdots & \vdots \\ 1 & \mathbf{a}_m^T \end{pmatrix}$$

is of full column rank.

This assumption is equivalent to saying that $\mathbf{a}_1, \dots, \mathbf{a}_m$ do not reside in a lower-dimensional affine space (i.e., a line if $n = 2$ and a plane if $n = 3$).

To guarantee the well definiteness of the SWLS algorithm (i.e., $\mathbf{x}^k \notin \mathcal{A}$ for all k), we will make the following assumption on the initial vector \mathbf{x}^0 .

Assumption 2. $\mathbf{x}^0 \in \mathcal{R}$, where

$$(3.8) \quad \mathcal{R} := \left\{ \mathbf{x} \in \mathbb{R}^n : f(\mathbf{x}) < \frac{\min_j \{d_j\}^2}{4} \right\}.$$

A similar assumption was made for the SFP method (see condition (2.13)). Note that for the true source location \mathbf{x}_{true} one has $f(\mathbf{x}_{\text{true}}) = \sum_{j=1}^m \varepsilon_j^2$. Therefore, \mathbf{x}_{true} satisfies Assumption 2 if the errors ε_j are smaller in some sense than the range measurements d_j . This is a very reasonable assumption since in real applications the errors ε_i are often an order of magnitude smaller than d_i . Now, if the initial point \mathbf{x}^0 is good enough in the sense that it is close to the true source location, then Assumption 2 will be satisfied. We have observed through numerical experiments that the solution to the LS problem (1.3) often satisfies Assumption 2 as the following example demonstrates.

Example 3.1. Consider the source localization problem with $m = 5$ and $n = 2$. We performed Monte Carlo runs, where in each run the sensor locations \mathbf{a}_j and the source location \mathbf{x} were randomly generated from a uniform distribution over the square $[-20, 20] \times [-20, 20]$. The observed distances d_j are given by (1.1) with ε_j being independently generated from a normal distribution with mean zero and standard deviation σ . In our experiments σ takes on four different values: $1, 10^{-1}, 10^{-2}$, and 10^{-3} . For each σ , N_σ denotes the number of runs for which the condition $f(\mathbf{x}_{\text{LS}}) < \frac{\min_j d_j^2}{4}$ holds, and the results are given in Table 1. Clearly, Assumption 2 fails only for high noise levels.

The following simple and important property will be used in our analysis.

LEMMA 3.1. *Let $\mathbf{x} \in \mathcal{R}$. Then*

$$(3.9) \quad \|\mathbf{x} - \mathbf{a}_j\| > d_j/2, \quad j = 1, \dots, m.$$

Proof. Suppose in contradiction that there exists j_0 for which $\|\mathbf{x} - \mathbf{a}_{j_0}\| \leq d_{j_0}/2$. Then

$$f(\mathbf{x}) = \sum_{j=1}^m (\|\mathbf{x} - \mathbf{a}_j\| - d_j)^2 \geq (\|\mathbf{x} - \mathbf{a}_{j_0}\| - d_{j_0})^2 \geq \frac{d_{j_0}^2}{4} \geq \frac{\min\{d_j\}^2}{4},$$

which contradicts $\mathbf{x} \in \mathcal{R}$. \square

A direct consequence of Lemma 3.1 is that any element in \mathcal{R} cannot be one of the sensors.

COROLLARY 3.1. *If $\mathbf{x} \in \mathcal{R}$, then $\mathbf{x} \notin \mathcal{A}$.*

3.3.2. Convergence analysis of the SWLS method. We begin with the following result which plays a key role in the forthcoming analysis.

LEMMA 3.2. *Let δ be a positive number, and let $t > \delta/2$. Then*

$$(3.10) \quad \left(\frac{s^2}{t} - \delta\right)^2 \geq 2(s - \delta)^2 - (t - \delta)^2$$

for every $s > \sqrt{\frac{\delta t}{2}}$, and equality is satisfied if and only if $s = t$.

Proof. Rearranging (3.10) one has to prove

$$A(s, t) \equiv \left(\frac{s^2}{t} - \delta\right)^2 - 2(s - \delta)^2 + (t - \delta)^2 \geq 0.$$

Some algebra shows that the expression $A(s, t)$ can be written as follows:

$$(3.11) \quad A(s, t) = \frac{1}{t}(s - t)^2 \left(\left(\frac{s}{\sqrt{t}} + \sqrt{t}\right)^2 - 2\delta \right).$$

Using the conditions $t > \delta/2$ and $s > \sqrt{\frac{\delta t}{2}}$, we obtain

$$(3.12) \quad \left(\frac{s}{\sqrt{t}} + \sqrt{t}\right)^2 - 2\delta > \left(\sqrt{\frac{\delta}{2}} + \sqrt{\frac{\delta}{2}}\right)^2 - 2\delta = 0.$$

Therefore, from (3.11) and (3.12) it readily follows that $A(s, t) \geq 0$ and that equality holds if and only if $s = t$. \square

Thanks to Lemma 3.2, we establish the next result which is essential in proving the monotonicity of the SWLS method.

LEMMA 3.3. *Let $\mathbf{y} \in \mathcal{R}$. Then the function $g(\mathbf{x}, \mathbf{y})$ given in (3.2) is well defined on $\mathbb{R}^n \times \mathcal{R}$, and with*

$$(3.13) \quad \mathbf{z} \in \underset{\mathbf{x} \in \mathbb{R}^n}{\operatorname{argmin}} g(\mathbf{x}, \mathbf{y}),$$

the following properties hold:

- (a) $f(\mathbf{z}) \leq f(\mathbf{y})$, and the equality is satisfied if and only if $\mathbf{z} = \mathbf{y}$;
- (b) $\mathbf{z} \in \mathcal{R}$.

Proof. By Corollary 3.1, any $\mathbf{y} \in \mathcal{R}$ implies $\mathbf{y} \notin \mathcal{A}$, and hence the function g given by (cf. (3.2))

$$g(\mathbf{x}, \mathbf{y}) = \sum_{i=1}^m \left(\frac{\|\mathbf{x} - \mathbf{a}_i\|^2}{\|\mathbf{y} - \mathbf{a}_i\|} - d_i \right)^2$$

is well defined on $\mathbb{R}^n \times \mathcal{R}$. Now, by (3.13) and $\mathbf{y} \in \mathcal{R}$ we have

$$(3.14) \quad g(\mathbf{z}, \mathbf{y}) \leq g(\mathbf{y}, \mathbf{y}) = f(\mathbf{y}) < \frac{\min\{d_j\}^2}{4}.$$

In particular,

$$\left(\frac{\|\mathbf{z} - \mathbf{a}_j\|^2}{\|\mathbf{y} - \mathbf{a}_j\|} - d_j \right)^2 < \frac{d_j^2}{4}, \quad j = 1, \dots, m,$$

from which it follows that

$$(3.15) \quad \frac{\|\mathbf{z} - \mathbf{a}_j\|^2}{\|\mathbf{y} - \mathbf{a}_j\|} \geq \frac{d_j}{2}, \quad j = 1, \dots, m.$$

Invoking Lemma 3.2, whose conditions are satisfied by (3.15) and Lemma 3.1, we obtain

$$\left(\frac{\|\mathbf{z} - \mathbf{a}_j\|^2}{\|\mathbf{y} - \mathbf{a}_j\|} - d_j \right)^2 \geq 2(\|\mathbf{z} - \mathbf{a}_j\| - d_j)^2 - (\|\mathbf{y} - \mathbf{a}_j\| - d_j)^2.$$

Summing over $j = 1, \dots, m$, we obtain

$$\sum_{j=1}^m \left(\frac{\|\mathbf{z} - \mathbf{a}_j\|^2}{\|\mathbf{y} - \mathbf{a}_j\|} - d_j \right)^2 \geq 2 \sum_{j=1}^m (\|\mathbf{z} - \mathbf{a}_j\| - d_j)^2 - \sum_{j=1}^m (\|\mathbf{y} - \mathbf{a}_j\| - d_j)^2.$$

Therefore, together with (3.14), we get

$$f(\mathbf{y}) \geq g(\mathbf{z}, \mathbf{y}) \geq 2f(\mathbf{z}) - f(\mathbf{y}),$$

showing that $f(\mathbf{z}) \leq f(\mathbf{y})$. Now, assume that $f(\mathbf{y}) = f(\mathbf{z})$. Then by Lemma 3.2 it follows that the following set of equalities is satisfied:

$$(3.16) \quad \|\mathbf{y} - \mathbf{a}_j\| = \|\mathbf{z} - \mathbf{a}_j\|, \quad j = 1, \dots, m,$$

which after squaring and rearranging reads as

$$(\|\mathbf{y}\|^2 - \|\mathbf{z}\|^2) - 2\mathbf{a}_j^T(\mathbf{y} - \mathbf{z}) = 0, \quad j = 1, \dots, m.$$

Therefore,

$$\begin{pmatrix} 1 & \mathbf{a}_1^T \\ 1 & \mathbf{a}_2^T \\ \vdots & \vdots \\ 1 & \mathbf{a}_m^T \end{pmatrix} \begin{pmatrix} \|\mathbf{y}\|^2 - \|\mathbf{z}\|^2 \\ -2(\mathbf{y} - \mathbf{z}) \end{pmatrix} = 0.$$

Thus, by Assumption 1, $\mathbf{z} = \mathbf{y}$, and the proof of (a) is completed. To prove (b), using (a) and (3.14), we get

$$f(\mathbf{z}) \leq f(\mathbf{y}) < \min_{j=1, \dots, m} \frac{d_j^2}{4},$$

proving that $\mathbf{z} \in \mathcal{R}$. \square

We are now ready to prove the main convergence results for the SWLS method.

THEOREM 3.1 (convergence of the SWLS method). *Let $\{\mathbf{x}^k\}$ be the sequence generated by the SWLS method. Suppose that Assumptions 1 and 2 hold true. Then*

- (a) $\mathbf{x}^k \in \mathcal{R}$ for $k \geq 0$;
- (b) for every $k \geq 0$, $f(\mathbf{x}^{k+1}) \leq f(\mathbf{x}^k)$ and equality holds if and only if $\mathbf{x}^{k+1} = \mathbf{x}^k$;
- (c) the sequence of function values $\{f(\mathbf{x}^k)\}$ converges;
- (d) the sequence $\{\mathbf{x}^k\}$ is bounded;
- (e) every convergent subsequence $\{\mathbf{x}^{k_l}\}$ satisfies $\mathbf{x}^{k_l+1} - \mathbf{x}^{k_l} \rightarrow \mathbf{0}$;
- (f) any limit point of $\{\mathbf{x}^k\}$ is a stationary point of f .

Proof. (a) and (b) The proof follows by induction on k using Lemma 3.3.

(c) The proof follows from the fact that $\{f(\mathbf{x}^k)\}$ is bounded below (by zero) and is a nonincreasing sequence.

(d) By (b), all of the iterates \mathbf{x}^k are in the level set $\text{Lev}(f, f(\mathbf{x}^0))$ which, by Lemma 2.3, establishes the boundedness of the sequence $\{\mathbf{x}^k\}$.

(e) Let $\{\mathbf{x}^{k_l}\}$ be a convergent subsequence, and denote its limit by \mathbf{x}^* . By claims (a) and (b), we have for every k that

$$f(\mathbf{x}^k) \leq f(\mathbf{x}^0) < \min_{j=1, \dots, m} \frac{d_j^2}{4},$$

which combined with the continuity of f implies $\mathbf{x}^* \in \mathcal{R}$ and hence $\mathbf{x}^* \notin \mathcal{A}$, by Corollary 3.1. Now, recall that

$$\mathbf{x}^{k_l+1} \in \operatorname{argmin}_{\mathbf{x}} g(\mathbf{x}, \mathbf{x}^{k_l}).$$

To prove the convergence of $\{\mathbf{x}^{k_l+1}\}$ to \mathbf{x}^* , we will show that every subsequence converges to \mathbf{x}^* . Let $\{\mathbf{x}^{k_{l_p}+1}\}$ be a convergent subsequence, and denote its limit by \mathbf{y}^* . Since

$$\mathbf{x}^{k_{l_p}+1} \in \operatorname{argmin}_{\mathbf{x} \in \mathbb{R}^n} g(\mathbf{x}, \mathbf{x}^{k_{l_p}}),$$

the following holds:

$$g(\mathbf{x}, \mathbf{x}^{k_{l_p}}) \geq g(\mathbf{x}^{k_{l_p}+1}, \mathbf{x}^{k_{l_p}}) \text{ for every } \mathbf{x} \in \mathbb{R}^n.$$

Taking the limits of both sides in the last inequality and using the continuity of the function f , we have

$$g(\mathbf{x}, \mathbf{x}^*) \geq g(\mathbf{y}^*, \mathbf{x}^*) \text{ for every } \mathbf{x} \in \mathbb{R}^n,$$

and hence

$$(3.17) \quad \mathbf{y}^* \in \operatorname{argmin}_{\mathbf{x} \in \mathbb{R}^n} g(\mathbf{x}, \mathbf{x}^*).$$

Since the sequence of function values converges, it follows that $f(\mathbf{x}^*) = f(\mathbf{y}^*)$. Invoking Lemma 3.3 with $\mathbf{y} = \mathbf{x}^*$ and $\mathbf{z} = \mathbf{y}^*$, we obtain $\mathbf{x}^* = \mathbf{y}^*$, establishing claim (e).

(f) To prove the claim, note that (3.17) and $\mathbf{x}^* = \mathbf{y}^*$ imply that

$$\mathbf{x}^* \in \operatorname{argmin}_{\mathbf{x} \in \mathbb{R}^n} g(\mathbf{x}, \mathbf{x}^*).$$

Thus, by the first order optimality conditions we obtain the following:

$$0 = \nabla_{\mathbf{x}} g(\mathbf{x}, \mathbf{x}^*)|_{\mathbf{x}=\mathbf{x}^*} = 4 \sum_{j=1}^m (\|\mathbf{x}^* - \mathbf{a}_j\| - d_j) \frac{\mathbf{x}^* - \mathbf{a}_j}{\|\mathbf{x}^* - \mathbf{a}_j\|} = 2 \nabla f(\mathbf{x}^*). \quad \square$$

As a direct consequence of Theorem 3.1, we obtain the following convergence in function values.

COROLLARY 3.2. *Let $\{\mathbf{x}^k\}$ be the sequence generated by the algorithm. Then $f(\mathbf{x}^k) \rightarrow f^*$, where f^* is the function value at some stationary point \mathbf{x}^* of f .*

As was shown for the SFP algorithm, global convergence of the sequence generated by the SWLS algorithm can also be established under the same condition, i.e., assuming that f admits isolated stationary points.

THEOREM 3.2 (convergence of the sequence). *Let $\{\mathbf{x}^k\}$ be generated by (3.3) such that Assumptions 1 and 2 hold. Suppose further that all stationary points of f are isolated. Then the sequence $\{\mathbf{x}^k\}$ converges to a stationary point.*

Proof. The proof is the same as the proof of Theorem 2.2. \square

4. Numerical examples. In this section we present numerical simulations illustrating the performance of the SFP and SWLS schemes, as well as numerical comparisons with the LS approach and with the SDR of the ML problem. The simulations were performed in MATLAB, and the semidefinite programs were solved by SeDuMi [14].

Before describing the numerical results, for the reader's convenience, we first recall the SDR proposed in [4], which will be used in our numerical experiments comparisons. The first stage is to rewrite problem (ML) given in (1.2) as

$$\begin{aligned} \min_{\mathbf{x}, \mathbf{g}} \quad & \sum_{j=1}^m (g_j - d_j)^2 \\ \text{s.t.} \quad & g_j^2 = \|\mathbf{x} - \mathbf{a}_j\|^2, \quad j = 1, \dots, m. \end{aligned}$$

Making the change of variables

$$\mathbf{G} = \begin{pmatrix} \mathbf{g} \\ 1 \end{pmatrix} (\mathbf{g}^T \quad 1), \quad \mathbf{X} = \begin{pmatrix} \mathbf{x} \\ 1 \end{pmatrix} (\mathbf{x}^T \quad 1),$$

problem (1.2) becomes

$$\begin{aligned} \min_{\mathbf{X}, \mathbf{G}} \quad & \sum_{j=1}^m (G_{jj} - 2d_j G_{m+1,j} + d_j^2) \\ \text{s.t.} \quad & G_{jj} = \text{Tr}(\mathbf{C}_j \mathbf{X}), \quad j = 1, \dots, m, \\ & \mathbf{G} \succeq \mathbf{0}, \quad \mathbf{X} \succeq \mathbf{0}, \\ & G_{m+1,m+1} = X_{n+1,n+1} = 1, \\ & \text{rank}(\mathbf{X}) = \text{rank}(\mathbf{G}) = 1, \end{aligned}$$

where

$$\mathbf{C}_j = \begin{pmatrix} \mathbf{I} & -\mathbf{a}_j \\ -\mathbf{a}_j^T & \|\mathbf{a}_j\|^2 \end{pmatrix}, \quad j = 1, \dots, m.$$

Dropping the rank constraints in the above problem, we obtain the desired SDR of problem (1.2). The SDR is not guaranteed to provide an accurate solution to the ML problem, but it can always be considered as an approximation of the ML problem.

In the first example, we show that the SWLS scheme usually possesses a larger region of convergence to the global minimum than the scheme SFP. This last property is further demonstrated in the second example, which compares the SFP and SWLS methods and also demonstrates the superiority of the SWLS scheme. The last example

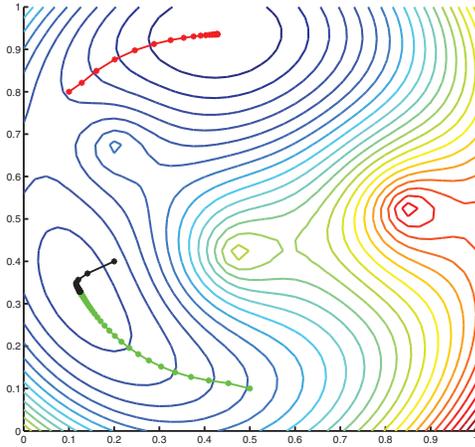


FIG. 1. The SFP method for three initial points.

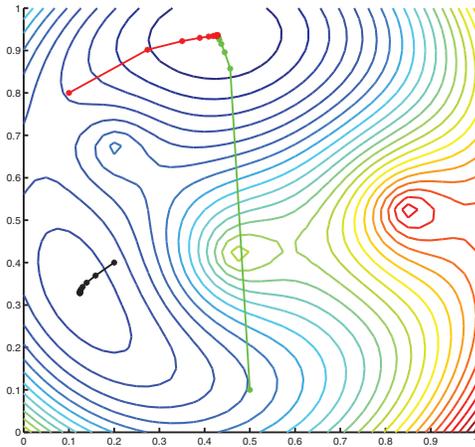


FIG. 2. The SWLS method for three initial points.

illustrates the attractiveness of the solution obtained by the SWLS method over the SDR and the LS approaches.

Example 4.1 (region of convergence of the SFP and SWLS methods). In this example we show typical behaviors of the SFP and SWLS methods. Consider an instance of the source localization problem in the plane ($n = 2$) with three sensors ($m = 3$) in the locations $(0.466, 0.418)$, $(0.846, 0.525)$ and $(0.202, 0.672)$. Figures 1 and 2 describe the results produced by the iterative schemes SFP and SWLS, respectively, for three initial trial points. The global minimum is $(0.4285, 0.9355)$, and there exists one additional local minimum at $(0.1244, 0.3284)$. As demonstrated in Figure 2, the SWLS method might converge to a local minimum; however, it seems to have a greater chance than the SFP algorithm to avoid local minima; for example, the SWLS converged to the (relatively far) global minimum from the initial starting point $(0.5, 0.1)$, while the SFP converged to the local minimum. We estimated the probability to converge to the global minimum by invoking both methods for 1681 initial starting

TABLE 2
Comparison between the SFP and SWLS methods.

m	#tight	$\#(f(\hat{\mathbf{x}}_{\text{SFP}}) > f(\hat{\mathbf{x}}_{\text{SWLS}}))$	Iter – SFP	Iter – SWLS
3	314	152	207(500.2)	26.2 (5)
4	325	96	124(192.6)	29.9(1.8)
5	259	83	93.6(96.2)	30.9(3.1)
10	278	23	66.5 (35.3)	31.6 (1.3)

points, which are the nodes of a 41×41 grid over the square $[0, 1] \times [0, 1]$. The SFP method converged to the global minimum in 45.87% of the runs, while the SWLS methods converged to the global minimum in 83.28% of the runs. Thus, the SWLS method has a much wider region of convergence to the global minimum. This was our observation in many other examples that we ran, which suggests that the SWLS method has the tendency to converge to the global minimum.

Remark 4.1. As shown in Proposition 2.1, the SFP scheme is just a gradient method with a fixed step size. Thanks to Lemma 2.5, which as shown in section 2.2 can be used in order to avoid the nonsmoothness, we can of course use more sophisticated smooth unconstrained minimization methods. Indeed, we also tested a gradient method with an Armijo step-size rule and a trust region method [8], which uses second order information. Our observation was that, while these methods usually possess an improved rate of convergence in comparison to the SFP method, they essentially have the same region of convergence to the global minimum as the SFP algorithm.

Example 4.2 (comparison of the SFP and SWLS methods). We performed Monte Carlo runs, where in each run the sensor locations \mathbf{a}_j and the true source location were randomly generated from a uniform distribution over the square $[-1000, 1000] \times [-1000, 1000]$. The observed distances d_j are given by (1.1) with ε_j being generated from a normal distribution with mean zero and standard deviation 20. Both the SFP and SWLS methods were employed with (the same) initial point, which was also uniformly randomly generated from the square $[-1000, 1000] \times [-1000, 1000]$. The stopping rule for both the SWLS and SFP methods was $\|\nabla f(\mathbf{x}^k)\| < 10^{-5}$.

The results of the runs are summarized in Table 2. For each value of m , 1000 realizations were generated. The numbers in the first column are the number of sensors, and in the second column we give the number of runs out of 1000 in which the SDR of the ML problem was tight; that is, the matrix which is the optimal solution of the SDR has rank one. We have also compared the SWLS solution with the SDR solution for these “tight” runs (about a quarter of the runs). In all of these runs, the SWLS and SDR solutions coincided; i.e., the SWLS method produced the exact ML solution. The third column contains the number of runs out of 1000 in which the solution produced by the SFP method was worse than the SWLS method. In all of the remaining runs, the two methods converge to the same point; thus, there were no runs in which the SWLS produced worse results. The last two columns contain the mean and standard deviation of the number of iterations of each of the methods in the form “mean (standard deviation).”

As can be clearly seen from the table, the SWLS method requires much less iterations than the SFP method, and in addition it is more robust in the sense that the number of iterations are more or less constant. In contrast, the standard deviations of the number of iterations of the SFP method are quite large. For example, the huge standard deviation 500.2 in the first row stems from the fact that in some of the runs the SFP algorithm required thousands of iterations!

TABLE 3
Mean squared position error of the SDR, LS and SWLS methods.

σ	SDR	LS	SWLS
$1e-3$	$2.4e-6$	$2.7e-6$	$1.5e-6$
$1e-2$	$2.2e-4$	$1.6e-4$	$1.3e-4$
$1e-1$	$2.2e-2$	$1.9e-2$	$1.3e-2$
$1e+0$	$2.2e+0$	$2.7e+0$	$2.0e+0$

From the above examples we conclude that the SWLS method does tend to converge to the global minimum. Of course, we can always construct an example in which the method converges to a local minimum (as was demonstrated in Example 4.1), but it seems that for random instances this convergence to a nonglobal solution is not likely.

We should also note that we also compared the SFP and SWLS methods with the initial point chosen as the solution of the LS problem (1.3). For this choice of the initial point, the SFP and SWLS methods always converged to the same location point⁴ (which is probably the global minimum); however, with respect to the number of iterations, the SWLS method was still significantly superior to the SFP algorithm. We have also compared the SWLS solution with the SDR solution for the runs in which the SDR solution is tight (about a quarter of the runs (cf. column 1 in Table 2)). In all of these runs, the SWLS and SDR solutions coincided; i.e., the SWLS method produced the exact ML solution.

The last example shows the attractiveness of the SWLS method over the LS and SDR approaches.

Example 4.3 (comparison with the LS and SDR estimates). Here we compare the solution of (1.3) and the solution of the SDR with the SWLS solution. The stopping rule for the SWLS method was $\|\nabla f(\mathbf{x}_k)\| < 10^{-5}$. We generated 100 random instances of the source localization problem with five sensors, where in each run the sensor locations \mathbf{a}_j and the source location \mathbf{x} were randomly generated from a uniform distribution over the square $[-10, 10] \times [-10, 10]$. The observed distances d_j are given by (1.1) with ε_j being independently generated from a normal distribution with mean zero and standard deviation σ . In our experiments σ takes four different values: $1, 10^{-1}, 10^{-2}$, and 10^{-3} . The numbers in the three right columns of Table 3 are the average of the squared position error $\|\hat{\mathbf{x}} - \mathbf{x}\|^2$ over 100 realizations, where $\hat{\mathbf{x}}$ is the solution by the corresponding method. The best result for each possible value of σ is marked in boldface. From the table, it is clear that the SWLS algorithm outperforms the LS and SDR methods for all four values of σ .

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⁴Numerically we used the criteria that two vectors \mathbf{x}_1 and \mathbf{x}_2 are “the same” if $\|\mathbf{x}_1 - \mathbf{x}_2\| \leq 10^{-8}$.

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