### Undergraduate Research Opportunity Programme in Science

# ADMM-type methods for linear programming problems

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#### Abstract

We reviewed and implemented several ADMM-type methods and conducted numerical experiments on large instances of the assignment problem. The methods are (1) classical ADMM with LP reformulation proposed by He and Yuan in [5], (2) Bregman ADMM (BADMM) proposed by Wang and Banerjee in [8] and our new semi-proximal ALM adopted from [4]. In order to achieve efficient updates, different methods require different formulations of the LP problem, with BADMM exploiting the special structure of the assignment problem to achieve fast updates while maintaining feasibility throughout. Meanwhile, the other two algorithms work for general standard form LP problems. For implementation of the semi-proximal ALM, sparse matrix representation, compressed matrix operations and numerical linear algebra routines written in C are exploited to boost performance. Gurobi, a commercial optimization software, is used as a benchmark for numerical experiments. Results of the experiments suggest that BADMM and semi-proximal ALM converge universally faster than the classical ADMM. For instances with small dimensions, BADMM marginally outperforms the semi-proximal ALM; when the dimension is large, the semi-proximal ALM significantly outperforms BADMM in terms of both iteration count and time per iteration, while the former also gives more accurate solutions within a fixed time.

### Chapter 1

### Introduction

We will introduce the standard form LP problem, the convex composite minimization problem with linear constraints and derive their respective optimality conditions. In Chapter 2 we discuss how to apply ADMM-type methods to LP with a short discussion on their convergence. In Chapter 3, we briefly discuss how the algorithms are implemented in Matlab. Numerical experiments are conducted to test and compare the algorithms. Chapter 4 summarizes the work done.

#### 1.1 Standard form LP problem

Consider the standard form LP

$$\min \left\{ \langle c, x \rangle \mid Ax = b, x \in \mathbb{R}^n_+ \right\} \tag{1.1}$$

where  $\mathbb{R}_+ = [0, \infty)$ ,  $A \in \mathbb{R}^{m \times n}$ ,  $b \in \mathbb{R}^n$  and  $\langle \cdot, \cdot \rangle$  denotes the usual inner product. A wide range of real-world problems can be modeled as LP and there has been extensive research on algorithms for LP. The most widely used algorithms for solving LP problems include the *simplex method* (with numerous enhancements) and the *interior point methods*. The alternating direction method of multipliers (ADMM) for convex composite optimization has been extensively researched and developed since 1970. In this paper, we study how ADMM-type methods can be used to solve LP problems efficiently.

#### 1.2 2-block convex minimization problem and classical ADMM

The classical ADMM solves the following convex composite optimization problem

$$\min \{ f_1(\mathbf{x}_1) + f_2(\mathbf{x}_2) \mid A_1\mathbf{x}_1 + A_2\mathbf{x}_2 = \mathbf{b}, \mathbf{x}_1 \in \mathcal{X}_1, \mathbf{x}_2 \in \mathcal{X}_2 \},$$
 (1.2)

where  $\mathcal{X}_i \subseteq \mathbb{R}^{d_i}$  are nonempty closed convex sets,  $\mathcal{A}_i \in \mathbb{R}^{m \times d_i}$  (or more generally, a linear operator from  $\mathbb{R}^{d_i}$  to  $\mathbb{R}^m$ ),  $\mathbf{b} \in \mathbb{R}^m$  and  $f_i : \mathbb{R}^{d_i} \to \mathbb{R} \cup \{\infty\}$  are closed *proper* convex functions for i = 1, 2.

We say f is proper if  $f(x) < \infty$  for at least one x.

**Definition** The augmented Lagrangian of (1.2) is defined as <sup>[5]</sup>

$$\mathcal{L}_{\sigma}(\boldsymbol{x}_{1}, \boldsymbol{x}_{2}, \boldsymbol{\lambda}) = f_{1}(\boldsymbol{x}_{1}) + f_{2}(\boldsymbol{x}_{2}) + \langle \boldsymbol{\lambda}, \mathcal{A}_{1}\boldsymbol{x}_{2} + \mathcal{A}_{2}\boldsymbol{x}_{2} - \boldsymbol{b} \rangle + \frac{\sigma}{2} \|\mathcal{A}_{1}\boldsymbol{x}_{1} + \mathcal{A}_{2}\boldsymbol{x}_{2} - \boldsymbol{b}\|^{2}$$
(1.3)

where  $\lambda \in \mathbb{R}^m$  is the Lagrangian multiplier and  $\sigma > 0$  is the penalty parameter.

**Definition** For any function  $f: \Omega \to \mathbb{R}$ , the set of minimizers of f on  $\Omega$  is denoted by

$$\arg\min\ \{f(x)\mid x\in\Omega\}=\Big\{y\in\Omega\mid f(y)=\inf\{f(x)\mid x\in\Omega\}\Big\}.$$

The classical ADMM consists of choosing an initial iterate  $(\boldsymbol{x}_1^0, \boldsymbol{x}_2^0, \boldsymbol{\lambda}^0)$  and the following updating steps [5]

$$\boldsymbol{x}_{1}^{k+1} \in \arg\min\{\mathcal{L}_{\sigma}(\boldsymbol{x}_{1}, \boldsymbol{x}_{2}^{k}, \boldsymbol{\lambda}^{k}) \mid \boldsymbol{x}_{1} \in \mathcal{X}_{1}\},$$
 (1.4a)

$$\boldsymbol{x}_{2}^{k+1} \in \arg\min\{\mathcal{L}_{\sigma}(\boldsymbol{x}_{1}^{k+1}, \boldsymbol{x}_{2}, \boldsymbol{\lambda}^{k}) \mid \boldsymbol{x}_{2} \in \mathcal{X}_{2}\},$$
 (1.4b)

$$\boldsymbol{\lambda}^{k+1} = \boldsymbol{\lambda}^k + \tau \sigma (\mathcal{A}_1 \boldsymbol{x}_1^{k+1} + \mathcal{A}_2 \boldsymbol{x}_2^{k+1} - \boldsymbol{b}), \tag{1.4c}$$

where the step length  $\tau \in (0, (1+\sqrt{5})/2)$ , [5]  $(\boldsymbol{x}_1^k, \boldsymbol{x}_2^k, \boldsymbol{\lambda}^k)$  are the current variables and  $(\boldsymbol{x}_1^{k+1}, \boldsymbol{x}_2^{k+1}, \boldsymbol{\lambda}^{k+1})$  the updated ones after the k-th iteration. The promising global convergence property of ADMM has been established, for example, in [6]. In theory, for the classical ADMM to converge to an optimal solution of (1.2), there is no restriction on the choice of  $(\boldsymbol{x}_1^0, \boldsymbol{x}_2^0, \boldsymbol{\lambda}^0)$ . However, only for problems with additional structures can we further characterize it (e.g., linear or superlinear rate of convergence). In addition, in [3], Wang and Banerjee have provided sufficient conditions for the convergence of the (generalized) BADMM, which implies convergence of the classical ADMM.

#### 1.3 Duality theories and optimality conditions

#### 1.3.1 Preliminaries

We introduce several concepts and theorems on characterization of optimal solutions via the KKT conditions, which are essential in theoretical convergence analysis and implementation of the algorithms. The following definitions and theorems are adopted from Section 5.5.3 in [2] and Appendix B.2 in [1].

**Definition** Let  $\Omega$  be a convex open subset of  $\mathbb{R}^n$ . A subgradient of a convex function  $f:\Omega\to\mathbb{R}\cup\{\infty\}$  at y is a vector  $v\in\mathbb{R}^n$  such that

$$f(x) \ge f(y) + \langle v, x - y \rangle$$

for any  $x \in \Omega$ . The subdifferential  $\partial f(y)$  is the set of all subgradients of f at y.

When f is differentiable at y,  $\partial f(y) = {\nabla f(y)}$ , the singleton set containing the gradient of f at y.

**Definition** Given a general minimization problem

$$\min \{ f(x) \mid x \in \mathbb{R}^n, h_i(x) \le 0, i = 1, ..., m, \ l_j(x) = 0, j = 1, ..., r \},$$
(1.5)

where  $h_i, l_j : \mathbb{R}^n \to \mathbb{R}$ , its Lagrangian is defined as

$$\mathcal{L}(x, u, v) = f(x) + \sum_{i=1}^{m} u_i h_i(x) + \sum_{j=1}^{r} v_i l_i(x).$$

The Lagrangian dual function is

$$g(u, v) = \min \{ \mathcal{L}(x, u, v) \mid x \in \mathbb{R}^n \}.$$

The  $dual\ problem$  is

$$\max \{ g(u, v) \mid u \in \mathbb{R}^m_+, \ v \in \mathbb{R}^n \}. \tag{1.6}$$

**Remark** In general, the dual problem is always convex since g is always concave. In addition, weak duality always holds:  $f^* \geq g^*$ , where  $f^*$  and  $g^*$  are primal and dual optimal values respectively. Additional conditions are needed for strong duality  $(f^* = g^*)$  to hold.

**Theorem 1.3.1** (Slater's Condition) If the primal problem is convex and there exists x such that  $h_i(x) < 0, i = 1, ..., j$  and  $l_j(x) = 0, j = 1, ..., r$ , then strong duality holds:  $f^* = g^*$ .

**Definition** The Karush-Kuhn-Tucker (KKT) conditions for problem (1.5) are

$$0 \in \partial f(x) + \sum_{i=1}^{m} u_i \partial h_i(x) + \sum_{i=1}^{r} v_j \partial l_j(x), \tag{1.7a}$$

$$u_i \cdot h_i(x) = 0, \ i = 1, ..., m$$
 (1.7b)

$$h_i(x) \le 0, \ i = 1, ..., m, \ l_j(x) = 0, \ j = 1, ..., r,$$
 (1.7c)

$$u_i \ge 0, \ i = 1, ..., m.$$
 (1.7d)

**Remark** Conditions (1.7a) - (1.7c) are frequently referred to as stationarity, complementarity, primary feasibility and dual feasibility respectively.

**Theorem 1.3.2** If  $x^*$  and  $(u^*, v^*)$  are primal and dual optimal solutions respectively with  $f(x^*) = g(u^*, v^*)$  (strong duality), then  $(x^*, y^*, z^*)$  must satisfy the KKT conditions (1.7a) - (1.7d).

In the above theorem we do not assume convexity of the problem. Furthermore, it can be shown that

**Theorem 1.3.3** If the problem is convex (i.e., f,  $h_i$  and  $l_i$  are convex) and  $(x^*, u^*, v^*)$  satisfies the KKT conditions (1.7a) - (1.7d), then  $x^*$  is an optimal solution to the primal problem (1.5) and  $(u^*, v^*)$  is an optimal solution to the dual problem (1.6).

The tuple  $(x^*, u^*, v^*)$  is known as a KKT point of problem (1.5) with dual problem (1.6). For convex problems, KKT conditions are always sufficient for primal and dual optimality. However, it does *not* imply strong duality, and is in fact a necessary condition for optimality when strong duality indeed holds.

In the following analysis, all problems we consider will be convex (convex objective functions with convex feasible set), for which KKT conditions are always sufficient for optimality.

#### 1.3.2 Optimality conditions for the standard form LP problem

As a special case of the general optimality conditions for convex optimization, we have

**Proposition 1.3.4** The KKT conditions for LP

$$Ax - b = 0, (1.8a)$$

$$A^T y + z - c = 0, (1.8b)$$

$$x_{(i)} \cdot z_{(i)} = 0, \ i = 1, ..., n$$
 (1.8c)

$$x, z \ge 0, \tag{1.8d}$$

which are necessary and sufficient for (primal and dual) optimality.

For the standard form LP problem, strong duality holds when the primal problem (1.1) has a finite optimal objective  $\langle c, x^* \rangle$ . In this case, its dual problem also has a finite optimal objective value  $b^T y^*$  with  $\langle c, x^* \rangle = b^T y^*$ .

#### 1.3.3 Optimality conditions for the 2-block convex minimization problem

We can derive a set of sufficient optimality conditions for problem (1.2) based on previous discussion. Consider the problem without the convex constraint sets  $\mathcal{X}_1$  and  $\mathcal{X}_2$ , whose Lagrangian is

$$\mathcal{L}\left(egin{pmatrix} oldsymbol{x}_1 \ oldsymbol{x}_2 \end{pmatrix}, oldsymbol{\lambda} 
ight) = f_1(oldsymbol{x}_1) + f_2(oldsymbol{x}_2) + \langle oldsymbol{\lambda}, \mathcal{A}_1 oldsymbol{x}_1 + \mathcal{A}_2 oldsymbol{x}_2 - oldsymbol{b} 
angle,$$

where  $(x_1, x_2)$  is viewed as one vector and there is no slack variables since the problem does not carry any inequality constraint. As such, its KKT conditions consist of only stationarity and primal feasibility conditions, namely,

$$0 \in \partial \begin{pmatrix} f_1(\boldsymbol{x}_1) \\ f_2(\boldsymbol{x}_2) \end{pmatrix} + \begin{pmatrix} \mathcal{A}_1^T \boldsymbol{\lambda} \\ \mathcal{A}_2^T \boldsymbol{\lambda} \end{pmatrix},$$
$$\mathcal{A}_1 \boldsymbol{x}_1 + \mathcal{A}_2 \boldsymbol{x}_2 - \boldsymbol{b} = 0.$$

The above conditions are sufficient for the optimality of the solutions to the problem without the constraints  $x_1 \in \mathcal{X}_1$  and  $x_2 \in \mathcal{X}_2$ . Hence, for the original problem (1.2) with the constraints, the sufficient optimality conditions are

$$-\mathcal{A}_1^T \boldsymbol{\lambda} \in \partial f_1(\boldsymbol{x}_1), \quad -\mathcal{A}_2^T \boldsymbol{\lambda} \in \partial f_2(\boldsymbol{x}_2), \tag{1.9a}$$

$$\mathcal{A}_1 \boldsymbol{x}_1 + \mathcal{A}_2 \boldsymbol{x}_2 - \boldsymbol{b} = 0, \tag{1.9b}$$

$$\boldsymbol{x}_1 \in \mathcal{X}_1, \quad \boldsymbol{x}_2 \in \mathcal{X}_2.$$
 (1.9c)

The above conditions (1.9a) - (1.9a) will be used in the convergence analysis of the generalized BADMM and to derive residuals used for implementation. Here we view  $A_i$  as matrices, although in general they can be linear operators on any Euclidean space.

#### 1.4 The assignment problem

The special type of LP problems used as test cases are known as assignment problems, which are special cases of transportation problems, which is defined as

$$\min\left\{\langle C, X \rangle \mid X \in \mathbb{R}_{+}^{m \times n}, X \mathbf{1}_{n} = a, X^{T} \mathbf{1}_{m} = b\right\},\tag{1.10}$$

where  $a \in \mathbb{R}^m$ ,  $b \in \mathbb{R}^n$ ,  $C \in \mathbb{R}^{m \times n}$ ,  $\mathbf{1}_k \in \mathbb{R}^k$  is a column vector of 1's for any  $k \in \mathbb{N}$ ,  $\langle \cdot, \cdot \rangle$  is the Euclidean inner product of two matrices of the same dimension,  $\langle A, B \rangle = \text{Tr}(A^T B) = \sum_{i=1}^m \sum_{j=1}^n A_{ij} B_{ij}$  for  $A, B \in \mathbb{R}^{m \times n}$ . Note that we must have  $a \geq 0$ ,  $b \geq 0$  and  $\mathbf{1}_m^T a = \mathbf{1}_n^T b$  for (1.10) to be feasible. The assignment problem is the transportation problem with m = n and  $a = b = \mathbf{1}_m$ .

It is not difficult to rewrite (1.10) into (1.1). Let  $C = (C_1, ..., C_n)$  and  $X = (X_1, ..., X_n)$  and

$$c = \begin{pmatrix} C_1 \\ \vdots \\ C_n \end{pmatrix} \in \mathbb{R}^{mn}, \quad x = \begin{pmatrix} X_1 \\ \vdots \\ X_n \end{pmatrix} \in \mathbb{R}_+^{mn}. \tag{1.11}$$

The constraints  $X\mathbf{1}_n = a$  and  $X^T\mathbf{1}_m = b$  then translate to

$$\begin{pmatrix} I_m \dots I_m \end{pmatrix} x = a, \begin{pmatrix} \mathbf{1}_m^T \\ & \ddots \\ & & \mathbf{1}_m^T \end{pmatrix} x = b, \tag{1.12}$$

where both block matrices consist of  $I_m$  and  $\mathbf{1}_m^T$  repeated n times respectively. To rewrite (1.10)

into (1.1), let

$$A = \begin{pmatrix} I_m & \cdots & I_m \\ \mathbf{1}_m^T & & \\ & \ddots & \\ & & \mathbf{1}_m^T \end{pmatrix} \in \{0, 1\}^{(m+n) \times (mn)}, \quad \boldsymbol{b} = \begin{pmatrix} a \\ b \end{pmatrix} \in \mathbb{R}^{m+n}. \tag{1.13}$$

Note that A in (1.13) is rank deficient, namely, the first (m+n-1) rows are linearly independent and the last row is a linear combination of them. Demote the first (m+n-1) rows of A as  $\tilde{A}$ . By Theorem 13.3 in [7] on a sufficient condition for total unimodularity, we see that  $\tilde{A}$  is totally unimodular. As such, all basic feasible solutions of the polyhedral  $\{x: \tilde{A}x = b, x \geq 0\}$  are integer vectors if A and b are defined as in (1.13) and b is a integer vector. In the case of an assignment problem where  $m=n, a=b=1_m$  and  $X\geq 0$ , we see that every basic feasible solution of  $\{x: \tilde{A}x = b, x \geq 0\}$  is a  $\{0,1\}$ -vector. Since the problem is bounded feasible, there exists an optimal basic feasible solution  $x^* \in \{0,1\}^{mn}$ . Since  $\{0,1\}^{mn}$  contains finitely many vectors, for any C that is randomly generated from a continuous distribution on  $\mathbb{R}_+^{m\times n}$ , the optimal solution  $x^*$  is unique almost surely. Converting  $x^*$  back to  $X^*$ , with the constraints in (1.10) in mind, this implies that each row or column of  $X^*$  has exactly one nonzero entry which equals 1. In other words, there exists an optimal solution  $X^*$  that is a permutation matrix; for randomly generated C, the optimal solution is almost always unique.

### Chapter 2

## Algorithms

#### 2.1 ADMM with reformulation

In [1], He and Yuan (2015) reformulate the LP problem with only inequality constraints into a form that is suitable for efficient ADMM iterations. In this section, we will extend the idea to the standard form LP problem.

We consider an equivalent form of the standard form LP problem (1.1)

$$\min \left\{ \sum_{i=1}^{m} \theta_i(x_i) \mid a_i^T x_i = b_i, x_i = \bar{x} \in \mathbb{R}_+^n, i = 1, 2, ..., m \right\},$$
(2.1)

where  $\theta_i : \mathbb{R}^n \to \mathbb{R}$ ,  $\theta_i(x) = c^T x$ , i = 1, ..., m. Note that the objective function in (2.1) is in fact scaled by m, the number of rows of A. We further define  $X_i = \{x \in \mathbb{R}^n \mid a_i^T x = b_i\}$ , i = 1, ..., m,  $\mathcal{X}_1 = X_1 \times ... \times X_m$ ,  $\mathcal{X}_2 = \bar{\mathcal{X}} = \mathbb{R}^n_+$ , and

$$egin{aligned} oldsymbol{x}_1 = egin{pmatrix} x_1 \ x_2 \ dots \ x_m \end{pmatrix} \in \mathcal{X}_1, \ oldsymbol{x}_2 = ar{x} \in \mathcal{X}_2 = ar{\mathcal{X}}, \ f_1 = \sum_{i=1}^m heta_i, \ f_2 = 0. \end{aligned}$$

As such, we can formulate (1.1) into the framework (1.2). In order to derive efficient updating formulas for the standard LP problem from (1.4a) - (1.4c), we need the following linear algebra facts.

**Lemma 2.1.1** For  $q \in \mathbb{R}^n$ , we have

$$q + \frac{b - \langle a, q \rangle}{\|a\|^2} \cdot a \in \arg\min\{\|x - q\|^2 \mid x \in \mathbb{R}^n, a^T x = b\},$$
 (2.2)

$$\{q^+\} = \arg\min\{\|x - q\|^2 \mid x \in \mathbb{R}_+^n\},$$
 (2.3)

where  $y^+ := \left(\max\{y_{(1)}, 0\}, ..., \max\{y_{(n)}, 0\}\right)^T$  for  $y = \left(y_{(1)}, ..., y_{(n)}\right)^T \in \mathbb{R}^n$ .

**Proof** A minimizer of  $\{\|x-q\|^2 \mid x \in \mathbb{R}^n, a^Tx = b\}$  is the projection of q onto the hyperplane  $S = \{x \in \mathbb{R}^n \mid a^Tx = b\}$ . Let  $x_0 \in S$ ,  $a^Tx_0 = b$ , the projection can be expressed as

$$q - \left\langle q - x_0, \frac{a}{\|a\|} \right\rangle \frac{a}{\|a\|} = q + \frac{b - \left\langle a, q \right\rangle}{\|a\|^2} \cdot a,$$

which gives (2.2). For (2.3), notice that

$$||x - q||^2 = \sum_{i=1}^{n} (x_{(i)} - q_{(i)})^2,$$

and each  $\left(x_{(i)}-q_{(i)}\right)^2$  attains its minimum at  $x_{(i)}=q_{(i)}$  if  $q_{(i)}\geq 0$  or  $x_{(i)}=0$  if  $q_{(i)}<0$ .

Let  $\lambda = \begin{pmatrix} \lambda_1 \\ \vdots \\ \lambda_m \end{pmatrix}$  with  $\lambda_i \in \mathbb{R}^n$ , i = 1, ..., m be the Lagrangian multiplier. From (1.4a), we have

$$\begin{aligned} \boldsymbol{x}_1^{k+1} &\in & \arg\min\left\{\mathcal{L}_{\sigma}(\boldsymbol{x}_1, \bar{\boldsymbol{x}}^k, \boldsymbol{\lambda}^k) \mid \boldsymbol{x}_1 \in \mathcal{X}_1\right\} \\ &= & \arg\min\left\{\sum_{i=1}^m c^T x_i + \sum_{i=1}^m (\lambda_i^k)^T x_i + \frac{\sigma}{2} \sum_{i=1}^m \|x_i - \bar{\boldsymbol{x}}^k\|^2 \mid \begin{pmatrix} x_1 \\ x_2 \\ \vdots \\ x_m \end{pmatrix} \in \mathcal{X}_1\right\} \end{aligned}$$

By (2.2), for i = 1, ..., m, we have

$$\begin{split} x_i^{k+1} &\in & \arg\min\left\{c^T x_i + (\lambda_i^k)^T x_i + \frac{\sigma}{2}\|x_i - \bar{x}^k\|^2 \mid x_i \in X_i\right\} \\ &= & \arg\min\left\{\|x - (\bar{x}^k - \frac{\sigma}{2}(c + \lambda_i^k))\|^2 \mid x \in \mathbb{R}^n, a_i^T x = b_i\right\} \\ &= & \left\{q_i^k + \frac{b_i - \langle a_i, q_i^k \rangle}{\|a_i\|^2} \cdot a_i\right\}, \end{split}$$

where  $q_i^k = \bar{x}^k - \frac{1}{\sigma}(c + \lambda_i^k) \in \mathbb{R}^n$ . Hence

$$\boldsymbol{x}_{1}^{k+1} = \begin{pmatrix} x_{1}^{k+1} \\ \vdots \\ x_{m}^{k+1} \end{pmatrix} \in \arg\min\left\{\mathcal{L}_{\sigma}(\boldsymbol{x}_{1}, \bar{x}^{k}, \boldsymbol{\lambda}^{k}) \mid \boldsymbol{x}_{1} \in \mathcal{X}_{1}\right\}, \tag{2.4}$$

where

$$x_i^{k+1} = q_i^k + \frac{b_i - \langle a_i, q_i^k \rangle}{\|a_i\|^2} \cdot a_i.$$
 (2.5)

From (1.4b), we have

$$\begin{aligned} \boldsymbol{x}_{2}^{k+1} &= \bar{x}^{k+1} &\in & \arg\min\left\{\mathcal{L}_{\sigma}(\boldsymbol{x}_{1}^{k+1}, \bar{x}, \boldsymbol{\lambda}^{k}) \mid \bar{x} \in \bar{\mathcal{X}}\right\} \\ &= & \arg\min\left\{-\sum_{i=1}^{m} (\lambda_{i}^{k})^{T} \bar{x} + \frac{\sigma}{2} \sum_{i=1}^{m} \|\boldsymbol{x}_{i}^{k+1} - \bar{x}\|^{2} \mid \bar{x} \in \bar{\mathcal{X}}\right\} \\ &= & \arg\min\left\{\|\bar{x} - \frac{1}{m} \sum_{i=1}^{m} (\boldsymbol{x}_{i}^{k+1} + \frac{1}{\sigma} \lambda_{i}^{k})\|^{2} \mid \bar{x} \in \mathbb{R}_{\geq 0}^{n}\right\}. \end{aligned}$$

Therefore, by (2.3),

$$\boldsymbol{x}_{2}^{k+1} = \bar{\boldsymbol{x}}^{k+1} = \left(\frac{1}{m}\sum_{i=1}^{m}(\boldsymbol{x}_{i}^{k+1} + \frac{1}{\sigma}\boldsymbol{\lambda}_{i}^{k})\right)^{+} \in \arg\min\left\{\mathcal{L}_{\sigma}(\boldsymbol{x}_{1}^{k+1}, \boldsymbol{x}_{2}, \boldsymbol{\lambda}^{k}) \mid \boldsymbol{x}_{2} \in \mathcal{X}_{2}\right\}. \tag{2.6}$$

The updating formula for the Lagrange multiplier  $\lambda$  (1.4c) is

$$\lambda_i^{k+1} = \lambda_i^k + \tau \sigma(x_i^{k+1} - y^k), \ i = 1, ..., m.$$
 (2.7)

#### 2.2 Generalized BADMM

The (generalized) BADMM was first introduced by Wang and Banerjee in [8]. The idea is to replace the quadratic penalty term in ADMM by the more general *Bregman divergence*. An immediate generalization is to further allow variable-specific Bregman divergence terms in order to achieve more efficient updating.

**Definition** Suppose  $\Omega \in \mathbb{R}^n$  is a closed convex subset set of a Euclidean space and  $\phi : \Omega \to \mathbb{R}$  is continuously differentiable and strictly convex on ri  $(\Omega)$ , the relative interior of  $\Omega$ . Define  $B_{\phi} : \Omega \times \text{ri}(\Omega) \to \mathbb{R}_+$ , the Bregman divergence induced by  $\phi$ , as

$$B_{\phi}(x,y) = \phi(x) - \phi(y) - \langle \nabla \phi(y), x - y \rangle, \tag{2.8}$$

where  $\nabla \phi(y)$  denotes the gradient of  $\phi$  at y.

For example, the square of the Euclidean 2-norm  $B_{\phi}(x,y) = ||x-y||^2$  can be obtained by taking  $\phi: z \mapsto ||z||^2$ , and the Kullback-Leibler (KL) divergence

$$B_{\phi}(x,y) = \sum_{i=1}^{n} x_{(i)} \log \frac{x_{(i)}}{y_{(i)}} - \sum_{i=1}^{m} x_{(i)} + \sum_{i=1}^{m} y_{(i)}$$
(2.9)

can be obtained by taking  $\phi: z \mapsto \sum_{i=1}^n z_{(i)} \log z_{(i)} - \sum_{i=1}^n z_{(i)}$ . Note that in both cases  $\phi$  is continuously differentiable and strictly convex on  $\mathbb{R}^n_{>0}$  (since their Hessians are well-defined and positive

definite). An important property of the Bregman divergence is  $B_{\phi}(x,y) \geq 0$  where the equality holds if and only if x = y. This is because  $\phi$  is strictly convex.

Under the framework (1.2), the generalized BADMM consists of the 3 updating steps

$$\mathbf{x}_{1}^{k+1} \in \arg \min\{f_{1}(\mathbf{x}_{1}) + \langle \boldsymbol{\lambda}^{k}, \mathcal{A}_{1}\mathbf{x}_{1} + \mathcal{A}_{2}\mathbf{x}_{2}^{k} - \boldsymbol{b} \rangle 
+ \sigma B_{\phi}(\boldsymbol{b} - \mathcal{A}_{1}\mathbf{x}_{1}, \mathcal{A}_{2}\mathbf{x}_{2}^{k}) + \sigma_{1}B_{\phi_{1}}(\mathbf{x}_{1}, \mathbf{x}_{2}^{k}) \mid \mathbf{x}_{1} \in \mathcal{X}_{1} \}, \qquad (2.10a)$$

$$\mathbf{x}_{2}^{k+1} \in \arg \min\{f_{2}(\mathbf{x}_{2}) + \langle \boldsymbol{\lambda}^{k}, \mathcal{A}_{1}\mathbf{x}_{1}^{k+1} + \mathcal{A}_{2}\mathbf{x}_{2} - \boldsymbol{b} \rangle 
+ \sigma B_{\phi}(\boldsymbol{b} - \mathcal{A}_{1}\mathbf{x}_{1}^{k+1}, \mathcal{A}_{2}\mathbf{x}_{2}) + \sigma_{2}B_{\phi_{2}}(\mathbf{x}_{1}^{k+1}, \mathbf{x}_{2}) \mid \mathbf{x}_{2} \in \mathcal{X}_{2} \}, \qquad (2.10b)$$

$$\boldsymbol{\lambda}^{k+1} = \boldsymbol{\lambda}^{k} + \tau \sigma(\mathcal{A}_{1}\mathbf{x}_{1}^{k+1} + \mathcal{A}_{2}\mathbf{x}_{2}^{k+1} - \boldsymbol{b}), \qquad (2.10c)$$

where  $\phi$ ,  $\phi_1$  and  $\phi_2$  are continuously differentiable and strictly convex functions on some convex domains depending on the problem,  $\sigma > 0$ ,  $\tau > 0$ ,  $\sigma_1 \ge 0$ ,  $\sigma_2 \ge 0$ .

The ordinary BADMM is obtained by setting  $\sigma_1 = \sigma_2 = 0$  in (2.10a) - (2.10c) and the classical ADMM (1.4a) - (1.4c) can be obtained by further setting  $\phi : z \mapsto \frac{1}{2}||z||$ , which gives  $B_{\phi}(x,y) = \frac{1}{2}||x-y||^2$ .

#### 2.2.1 Convergence analysis

The following discussion is based on from Section 3 in [8]. We establish the convergence of the generalized BADMM, which them gives the convergence of BADMM without variable-specific Bregman divergence terms and ADMM (perhaps with different restrictions on step lengths).

Referring to problem (1.2) and the iterative steps (2.10a) - (2.10c), in additions to the properties of  $f_1$ ,  $f_2$ ,  $\mathcal{X}_1$  and  $\mathcal{X}_2$  specified by (1.2), for a chosen  $\phi$ , additional technical conditions on  $\sigma$ ,  $\sigma_1$ ,  $\sigma_2$  and  $\tau$  are needed for (2.10a) - (2.10c) to generate a sequence that converges to an optimal solution.

Let  $(\boldsymbol{x}_1^*, \boldsymbol{x}_2^*, \boldsymbol{\lambda}^*)$  satisfies the optimality conditions (1.9a) - (1.9c). Consider the subproblems (2.10a) and (2.10b) (finding  $\boldsymbol{x}_1^{k+1}$  and  $\boldsymbol{x}_2^{k+1}$  as minimizers of two modified augmented Lagrangian functions with Bregman terms). For (2.10a), the optimality conditions are

$$egin{array}{lll} m{x}_1^{k+1} & \in & \mathcal{X}_1, \\ 0 & \in & \partial f_1(m{x}_1^{k+1}) + \mathcal{A}_1^T m{\lambda} + \sigma 
abla_{m{x}_1} B_{\phi}(m{b} - \mathcal{A}_1 m{x}_1^{k+1}, \mathcal{A}_2 m{x}_2^k) + \sigma_1 
abla_{m{x}_1} B_{\phi_1}(m{x}_1^{k+1}, m{x}_2^k), \end{array}$$

which can be written as

$$\boldsymbol{x}_{1}^{k+1} \in \mathcal{X}_{1}, \quad -\mathcal{A}_{1}^{T} \boldsymbol{\lambda}^{k} - \sigma \mathcal{A}_{1}^{T} \left( \nabla \phi(\boldsymbol{b} - \mathcal{A}_{1} \boldsymbol{x}_{1}^{k+1}) - \nabla \phi(\mathcal{A}_{2} \boldsymbol{x}_{2}^{k}) \right) - \sigma_{1} \left( \nabla \phi_{1}(\boldsymbol{x}_{1}^{k+1}) - \nabla \phi_{1}(\boldsymbol{x}_{1}^{k}) \right) \in \partial f_{1}(\boldsymbol{x}_{1}^{k+1}). \tag{2.11}$$

Note that we have used  $\nabla_x B_{\phi}(x,y) = \nabla \phi(x) - \nabla \phi(y)$ . Similarly, for (2.10b), the optimality condi-

tions are

$$\boldsymbol{x}_2^{k+1} \in \mathcal{X}_2, \ -\mathcal{A}_2^T \boldsymbol{\lambda}^k - \sigma \mathcal{A}_2^T \left( \nabla \phi(\mathcal{A}_2 \boldsymbol{x}_2^{k+1}) - \nabla \phi(\boldsymbol{b} - \mathcal{A}_1 \boldsymbol{x}_1^{k+1}) \right) - \sigma_2 \left( \nabla \phi_2(\boldsymbol{x}_2^{k+1}) - \nabla \phi_2(\boldsymbol{x}_2^k) \right) \in \partial f_2(\boldsymbol{x}_2^{k+1}).$$

**Lemma 2.2.1** If  $(x_1^k, x_2^k, \lambda^k)$  and  $(x_1^{k+1}, x_2^{k+1}, \lambda^{k+1})$  satisfy

$$B_{\phi_1}(\boldsymbol{x}_1^{k+1}, \boldsymbol{x}_1^k) = 0 \ B_{\phi_2}(\boldsymbol{x}_2^{k+1}, \boldsymbol{x}_2^k) = 0,$$
 (2.12)

$$\mathcal{A}_2(\boldsymbol{x}_2^{k+1} - \boldsymbol{x}_2^k) = 0, \quad \mathcal{A}_1 \boldsymbol{x}_1^{k+1} + \mathcal{A}_2 \boldsymbol{x}_2^{k+1} - \boldsymbol{b} = 0, \tag{2.13}$$

$$\boldsymbol{x}_1 \in \mathcal{X}_1, \ \boldsymbol{x}_2 \in \mathcal{X}_2, \tag{2.14}$$

then  $(x_1^k, x_2^k, \lambda^k) = (x_1^{k+1}, x_2^{k+1}, \lambda^{k+1})$  is a KKT point.

**Proof** First, (2.17c) and the second equation in (2.13) imply that  $\boldsymbol{\lambda}^{k+1} = \boldsymbol{\lambda}^k$ . We will show that  $(\boldsymbol{x}_1^k, \boldsymbol{x}_2^k, \boldsymbol{\lambda}^k)$  satisfies (1.9a) - (1.9c). In fact, the two equations in (2.13) imply that  $\mathcal{A}_1 \boldsymbol{x}_1^{k+1} + \mathcal{A}_2 \boldsymbol{x}_2^k = \boldsymbol{b}$ , while (2.12) implies  $\boldsymbol{x}_1^{k+1} = \boldsymbol{x}_1^k$ . Hence by (2.12), we have

$$-\mathcal{A}_1 \boldsymbol{\lambda}^k \in \partial f_1(\boldsymbol{x}_1^k).$$

Similarly, (2.12) implies  $\boldsymbol{x}_2^{k+1} = \boldsymbol{x}_2^k$ , and by (2.12) we have

$$-\mathcal{A}_2 \boldsymbol{\lambda}^k \in \partial f_1(\boldsymbol{x}_2^k).$$

Note that in the case of BADMM with  $\sigma_1 = \sigma_2 = 0$ , (2.13) and (2.14) serve as sufficient optimality conditions for optimality for the general problem (1.2).

**Definition** (See (31) in [8]) The residual of the optimality condition (2.12) and (2.13) is defined as

$$R(k) = B_{\phi}(\boldsymbol{b} - \mathcal{A}_{1}\boldsymbol{x}_{1}^{k+1}, \mathcal{A}_{2}\boldsymbol{x}_{2}^{k}) + \frac{\sigma_{1}}{\sigma}B_{\phi_{1}}(\boldsymbol{x}_{1}^{k+1}, \boldsymbol{x}_{1}^{k}) + \frac{\sigma_{2}}{\sigma}B_{\phi_{2}}(\boldsymbol{x}_{2}^{k+1}, \boldsymbol{x}_{2}^{k}) + \gamma \|\mathcal{A}_{1}\boldsymbol{x}_{1}^{k+1} + \mathcal{A}_{2}\boldsymbol{x}_{2}^{k+1} - \boldsymbol{b}\|,$$
(2.15)

where  $\gamma > 0$ .

If R(k) = 0 and  $\mathbf{x}_1^k \in \mathcal{X}_1$ ,  $\mathbf{x}_2^k \in \mathcal{X}_2$ , (2.12) - (2.14) are satisfied and an optimal solution is obtained. Hence, in order to establish convergence of the generalized BADMM, it is sufficient to show R(k) converges to 0 under certain conditions on the parameters. The following theorem in [2] summarizes the above discussion.

**Theorem 2.2.2** Suppose (i) an optimal solution of (1.2) exists and (ii) the Bregman divergence  $B_{\phi}$  is defined on  $\phi: \Omega \to \mathbb{R}$  which is  $\alpha$ -strongly convex with respect to a p-norm. In other words, there exists  $\alpha > 0$  such that for any  $x \in \Omega$ ,  $y \in \Omega \setminus \Omega$ ,

$$B_{\phi}(x,y) \ge \frac{\alpha}{2} \|x - y\|_{p}^{2}.$$
 (2.16)

Let the sequence  $(\boldsymbol{x}_1^k, \boldsymbol{x}_2^k, \boldsymbol{\lambda}^k)$  be generated by the generalized BADMM iterative steps (2.10a) - (2.10c). Let  $(\boldsymbol{x}_1^k, \boldsymbol{x}_2^k, \boldsymbol{\lambda}^k)$  be a KKT point. If  $\delta = \min\{1, m^{\frac{2}{p}-1}\}$ ,  $\tau \leq (\alpha\delta - 2\gamma)$  and  $0 < \gamma < \frac{\alpha\delta}{2}$  for some  $\gamma > 0$ , then  $R(k) \to 0$  and  $(\boldsymbol{x}_1^k, \boldsymbol{x}_2^k, \boldsymbol{\lambda}^k) \to (\boldsymbol{x}_1^k, \boldsymbol{x}_2^k, \boldsymbol{\lambda}^k)$  as  $k \to \infty$ .

**Remark** When  $p \in (0,2]$ , we have  $\delta = 1$ , and  $\tau \leq \alpha - 2\gamma$ . In fact, for  $\phi(x) = \frac{1}{2}||x||^2$  and  $B_{\phi}(x,y) = \frac{1}{2}||x-y||^2$ , (2.16) holds with  $\alpha = 1$  and p = 2; for  $B_{\phi}$  being the KL divergence, we have (2.16) holds with  $\alpha = p = 1$  in the interior of the unit simplex.

#### 2.2.2 BADMM for the assignment problem

We then derive the updating formulas for the transportation problem (1.10) from (2.10a) - (2.10c). Instead of converting (1.10) into (1.1), we will convert its original matrix form into (1.2). In fact, this can be done by setting

$$\boldsymbol{x}_1 = X \in \mathcal{X} = \{X \in \mathbb{R}_+^{m \times n} \mid X \mathbf{1}_n = a\},$$

$$\boldsymbol{x}_2 = Z \in \mathcal{Z} = \{Z \in \mathbb{R}_+^{m \times n} \mid X^T \mathbf{1}_m = b\},$$

$$\boldsymbol{\lambda} = Y,$$

$$\boldsymbol{\lambda}_1 = I, \ \boldsymbol{\lambda}_2 = -I, \boldsymbol{b} = 0,$$

$$f_1(X) = \langle C, X \rangle, \ f_2(X) = 0.$$

Again, to have a feasible problem, it is necessary that  $a \ge 0$ ,  $b \ge 0$  and  $\mathbf{1}_m^T a = \mathbf{1}_n^T b$ . Let  $\sigma > 0$ ,  $\sigma_1 = \sigma_2 = 0$  and  $B_{\phi} = B_{KL} : (A, B) \mapsto \sum_{i=1}^m \sum_{j=1}^n (A_{ij} \log \frac{A_{ij}}{B_{ij}} - A_{ij} + B_{ij})$ , the KL divergence (2.9). By (2.10a) - (2.10c), one iteration consists of

$$X^{k+1} \in \arg\min\{\langle C, X \rangle + \langle Y^{k+1}, X \rangle + \sigma B_{KL}(X, Z^k) \mid X \in \mathcal{X}\}, \tag{2.17a}$$

$$Z^{k+1} \in \arg\min\{\langle Y^k, -Z \rangle + \sigma B_{KL}(Z, X^{k+1}) \mid Z \in \mathcal{Z}\}, \tag{2.17b}$$

$$Y^{k+1} = Y^k + \frac{1}{2}\sigma(X^{k+1} - Z^{k+1}). \tag{2.17c}$$

Note that we set the step length  $\tau=1/2$  in (2.17c), where we take  $\gamma=1/4$  and  $\alpha=p=1$  in Theorem 2.2.2. Setting  $\alpha=p=1$  makes (2.16) hold in the interiors of  $\mathcal X$  and  $\mathcal Z$ .

**Proposition 2.2.3** The updating formulas (2.17a) and (2.17b) both have closed from solutions

$$X_{ij}^{k+1} = \frac{Z_{ij}^k \exp(-\frac{C_{ij} + Y_{ij}^k}{\sigma})}{\sum_{j=1}^n Z_{ij}^k \exp(-\frac{C_{ij} + Y_{ij}^k}{\sigma})} a_i,$$
 (2.18a)

$$Z_{ij}^{k+1} = \frac{X_{ij}^{k+1} \exp(\frac{Y_{ij}^k}{\sigma})}{\sum_{i=1}^m X_{ij}^{k+1} \exp(\frac{Y_{ij}^k}{\sigma})} b_j.$$
 (2.18b)

**Proof** We show that  $X^{k+1}$  given by (2.18a) satisfies (2.17a). Denote

$$h(X) = \langle C, X \rangle + \langle Y^{k+1}, X \rangle + \sigma B_{KL}(X, Z^k).$$

Notice that

$$h(X) = \sum_{i=1}^{m} h_i(X_{i1}, \dots, X_{in}),$$

where

$$h_i(X_{i1}, \dots, X_{in}) = \sum_{j=1}^n \left( (C_{ij} + Y_{ij}^{k+1}) X_{ij} + \sigma(X_{ij} \log \frac{X_{ij}}{Z_{ij}^k} - X_{ij} + Z_{ij}^k) \right).$$

Let

$$g_i(X_{i1}, \dots, X_{in}) = \sum_{i=1}^n X_{ij} - a_i.$$

The problem is reduced to minimizing  $h_i(X_{i1}, \dots, X_{in})$  subject to the constraint  $g_i(X_{i1}, \dots, X_{in}) = 0$  for each i. Consider the Lagrangian

$$\mathcal{L}_i(X_{i1},\cdots,X_{in},\lambda) = h_i(X_{i1},\cdots,X_{in}) + \lambda g_i(X_{i1},\cdots,X_{in}).$$

Since  $h_i$  is strictly convex (positive definite Hessian on  $\mathbb{R}^n_{>0}$ ), a possible set of optimality conditions for the *i*-th subproblem is

$$g_i(X_{i1}, \dots, X_{in}) = 0,$$

$$\frac{\partial}{\partial X_{ii}} \mathcal{L}_i(X_{i1}, \dots, X_{in}, \lambda) = 0, \quad j = 1, ..., n$$

and

$$\frac{\partial}{\partial \lambda} \mathcal{L}_h(X_{i1}, \cdots, X_{in}, \lambda) = 0.$$

In other words,

$$(C_{ij} + Y_{ij}^{k+1}) + \sigma (\log X_{ij}^{k+1} - \log Z_{ij}^{k}) + \lambda = 0, \ j = 1, ..., n$$
 (2.19)

$$X_{i1}^{k+1} + \dots + X_{in}^{k+1} = a_i. {(2.20)}$$

From (2.19) we have

$$X_{ij} = \exp(-\lambda^*) \cdot Z_{ij}^{k+1} \exp\left(-\frac{C_{ij} + Y_{ij}^k}{\sigma}\right)$$
 (2.21)

Substitute (2.21) into (2.20) we obtain

$$\exp(-\lambda) = \frac{a_i}{\sum_{j=1}^n Z_{ij}^k \exp\left(-\frac{C_{ij} + Y_{ij}^k}{\sigma}\right)}.$$

Hence (2.18a) is the solution to (2.19) and solves (2.17a). A similar argument proves (2.18b).

One desirable property of the above iterative scheme (2.17a) - (2.17c) is that the feasibility of  $X^k$  and  $Z^k$  is maintained throughout, as long as we choose  $X^0 \in \mathcal{X}$  and  $Z^0 \in \mathcal{Z}$  such that  $X^0, Z^0 > 0$ , which is possible when a, b > 0.

#### 2.2.3 ADMM for the assignment problem in matrix form

Note that the classical ADMM is a special case of the generalized BADMM with  $\sigma_1 = \sigma_2 = 0$  and  $B_{\phi}$  being the squared Euclidean norm. As such, the assignment problem (1.10) can be written in the form for classical ADMM:  $\min\{\langle C, X \rangle \mid X - Z = 0, X \in \mathcal{X}, Z \in \mathcal{Z}\}$ . The updating formulas are therefore, by (1.4a) - (1.4c) or (2.17a) - (2.17c),

$$X^{k+1} \in \arg\min\{\langle C, X \rangle + \langle Y^k, X \rangle + \frac{\sigma}{2} \|X - Z^k\|^2 \mid X \in \mathcal{X}\}, \tag{2.22a}$$

$$Z^{k+1} \quad \in \quad \arg\min\{\langle Y^k, -Z\rangle + \frac{\sigma}{2}\|X^{k+1} - Z\|^2 \mid Z \in \mathcal{Z}\}, \tag{2.22b}$$

$$Y^{k+1} = Y^k + \tau \sigma(X^{k+1} - Z^{k+1}), \tag{2.22c}$$

where  $\tau \in (0, (1+\sqrt{5})/2)$  is the step length. Using (2.2), (2.22a) translates to for every i,

$$(X_{i1}^{k+1}, \dots, X_{in}^{k+1}) \in \arg\min \left\{ \sum_{j=1}^{n} \left( (C_{ij} + Y_{ij}^{k}) X_{ij} + \frac{\sigma}{2} (X_{ij} - Z_{ij}^{k})^{2} \right) \mid X_{i1} + \dots + X_{in} = a_{i}, X_{ij} \ge 0 \right\}$$

$$= \arg\min \left\{ \sum_{j=1}^{n} \left( X_{ij} - (Z_{ij}^{k} - \frac{1}{\sigma} \left( C_{ij} + Y_{ij}^{k} \right) \right) \right)^{2} \mid X_{i1} + \dots + X_{in} = a_{i}, X_{ij} \ge 0 \right\}.$$

$$(2.23)$$

Solving (2.23) is equivalent to finding the "projection" of a vector in  $\mathbb{R}^n$  onto a simplex. An efficient method in O(n) time (where n is the dimension of the Euclidean space) is proposed in [3]. It has been shown in [8] that this iterative scheme (2.22a) - (2.22c) converges slower than Gurobi and BADMM when solving assignment problems and does not converge after a maximum number of iterations when the dimension of input instances is large.

#### 2.3 Semi-proximal augmented Lagrangian method

We will content ourselves by considering the special case of the general semi-proximal ALM for standard form LP problems. The general algorithm can be found in Appendix B in [4]. Consider the dual of the standard form LP in minimization form

$$\min \{-b^T y \mid A^T y + z = c, \ z \ge 0\}$$
 (2.24)

and its augmented Lagrangian function

$$\mathcal{L}_{\sigma}(y, z, x) = -b^{T}y + \langle x, A^{T}y + z - c \rangle + \frac{\sigma}{2} ||A^{T}y + z - c||^{2}.$$
(2.25)

The semi-proximal augmented Lagrangian method consists of the following 4 updating steps

$$\bar{y}^{k+1} \in \arg\min\{\mathcal{L}_{\sigma}(y, z^k, x^k) \mid y \in \mathbb{R}^m\},$$
 (2.26a)

$$z^{k+1} \in \arg\min\{\mathcal{L}_{\sigma}(\bar{y}^{k+1}, z, x^k) \mid z \in \mathbb{R}^n_+\}, \tag{2.26b}$$

$$y^{k+1} \in \arg\min\{\mathcal{L}_{\sigma}(y, z^{k+1}, x^k) \mid y \in \mathbb{R}^m\}, \tag{2.26c}$$

$$x^{k+1} = x^k + \tau \sigma (A^T y^{k+1} + z^{k+1} - c), \tag{2.26d}$$

where in order to establish convergence we need to set  $\tau \in (0,2)$ .

We show that (2.26a) - (2.26c) all have closed form solutions. For (2.26a),  $\mathcal{L}_{\sigma}(y, z^k, x^k)$  is a quadratic function of  $y \in \mathbb{R}^m$ . Since there is no constraint on y, a possible optimality condition is

$$\nabla_{y} \mathcal{L}_{\sigma}(\bar{y}^{k+1}, z^{k}, x^{k}) = 0,$$

which gives

$$-b + Ax^{k} + \sigma A(A^{T}\bar{y}^{k+1} + z^{k} - c) = 0,$$

or

$$AA^{T}\bar{y}^{k+1} = \frac{1}{\sigma}(b - Ax^{k}) - A(z^{k} - c), \tag{2.27}$$

which has a unique solution when rank(A) = m. For (2.26b), notice that

$$\arg \min \{ \mathcal{L}_{\sigma}(\bar{y}^{k+1}, z, x^{k}) \mid z \in \mathbb{R}_{+}^{n} \}$$

$$= \arg \min \left\{ (x^{k})^{T} z + \frac{\sigma}{2} \|A^{T} \bar{y}^{k+1} + z - c\|^{2} \mid z \in \mathbb{R}_{+}^{n} \right\}$$

$$= \arg \min \left\{ \|z - (c - A^{T} \bar{y}^{k+1} - \frac{1}{\sigma} x^{k})\| \mid z \in \mathbb{R}_{+}^{n} \right\}.$$

By (2.3), we have

$$z^{k+1} = \left(c - A^T \bar{y}^{k+1} - \frac{1}{\sigma} x^k\right)^+ \in \arg\min\{\mathcal{L}_{\sigma}(\bar{y}^{k+1}, z, x^k) \mid z \in \mathbb{R}_+^n\}.$$
 (2.28)

Similar to the derivation of (2.27),

$$AA^{T}y^{k+1} = \frac{1}{\sigma}(b - Ax^{k}) - A(z^{k+1} - c).$$
 (2.29)

### Chapter 3

# Numerical experiments and results

We tested the above algorithms with their respective problem formulations in Matlab. Gurobi, a commercial optimization software, is used to verify the optimality of answers and as a performance benchmark.

Unlike the simplex method which, in theory, gives an exact optimal basic feasible solution to a bounded feasible LP problem, the ADMM-type methods only guarantee, in theory, a sequence of points that converge to an optimal solution. As such, for implementation of the algorithms and numerical experiments, we need to set termination conditions. Note that all algorithms of interest consist of initialization and iteratively applying a set of updating formulas. Hence, we may set the terminal conditions as either (a) the total number of iteration exceeds a predetermined threshold or (b) the residual of the optimality conditions is sufficiently small.

#### 3.1 Residuals

In this section, we derive formulas for residuals mentioned above from both theoretical and computational perspectives.

**Definition** In the context of the standard form LP problem (1.1) with dual problem (??), consider its KKT conditions (1.8a) - (1.8d), for  $\{(x^k, y^{k+1}, z^k)\}_{k>0}$ , define

$$R_{LP}(k) = \max \left\{ \frac{\|Ax^k - b\|}{1 + \|b\|}, \frac{\|A^Ty^k + z^k - c\|}{1 + \|c\|} \right\}.$$
 (3.1)

The denominators are included to enhance numerical stability in computation: to normalize the magnitudes of the norms in the numerators and make the program less sensitive to scaling of input data. Since the standard form LP problem is convex and strong duality holds if the problem is feasible and bounded,  $x^*$  solves the primal problem and  $(y^*, z^*)$  solves the dual problem (with the same optimal objective value) if and only if  $(x^*, y^*, z^*)$  is a KKT point. By the continuity of  $R_{LP}(k)$  in  $(x^k, y^k, z^k)$ , we have

**Proposition 3.1.1** for any sequence  $(x^k, y^k, z^k)$  that converges to a KKT point, we have  $R_{LP}(k) \to 0$ . Conversely, if  $(x^k, y^k, z^k)$  converges to  $(x^*, y^*, z^*)$ ,  $R_{LP}(k) \to 0$  and for each k, we have  $x^k \ge 0$ ,

 $z^{k+1} \geq 0$  and  $(z^k)_{(i)} \cdot (x^k)_{(i)} \rightarrow 0$  for any i = 1, ..., n (the dual feasibility and complementarity conditions), then  $(x^*, y^*, z^*)$  is a KKT point.

For semi-proximal augmented Lagrangian method,  $R_{LP}(k)$  can be easily computed since the tuple  $(x^k, y^k, z^k)$  is maintained during each iteration. For the classical ADMM with reformulation, since the dual and slack variables are not maintained, we need an alternative but similar definition of residual, which involves constructing the slack variable z from primal, dual feasibility and complementarity.

**Definition** In the context of Section 2.1, let  $\eta^k \in \mathbb{R}^n$  with

$$(\eta^k)_{(i)} = \frac{b_i - a_i^T q_i}{a_i^T a_i},$$

where  $q_i^k = \bar{x}^k - \frac{1}{\sigma}(c + \lambda_i^k) \in \mathbb{R}^n$ . Define the residual as

$$R_{LP}^{ADMM}(k) = \max\left\{\frac{\|A\bar{x}^k - b\|}{1 + \|b\|}, \frac{\|A^Ty^k + z^{k+1} - c\|}{1 + \|c\|}\right\},\tag{3.2}$$

where  $\bar{x}^k$  is maintained in every iteration,  $y^k = \sigma \eta^k/m$  and  $z^{k+1} = \sigma(\bar{x}^{k+1} - \bar{x}^k) + c - A^T y$  are the modified dual and slack variables respectively.

Note that from the above definition we have  $A^Ty^k + z^{k+1} - c = \sigma(\bar{x}^{k+1} - \bar{x}^k)$ , which is consistent with (2.13) and (2.14). In other words,  $R_{LP}^{ADMM}(k) = 0$  implies that a solution to the optimality conditions (2.13) - (2.14) is obtained.

**Proposition 3.1.2** If a sequence  $(\boldsymbol{x}_1^k, \bar{x}^k, \boldsymbol{\lambda}^k)$  generated by ADMM with reformulation (2.4) - (2.7) converges to a point  $(\boldsymbol{x}_1^*, \bar{x}^*, \boldsymbol{\lambda}^*)$  with  $R_{LP}^{ADMM}(k) \to 0$ , then we have  $(\boldsymbol{x}_1^*, \bar{x}^*, \boldsymbol{\lambda}^*)$  is a KKT point of the problem.

For the BADMM, we will use the residual defined in (2.15), which is a generalization of the residual of the optimality condition for the 2-block convex minimization problem with linear constraints. In the case of KL divergence and the assignment problem, we have

$$R_{BADMM}(k) = B_{KL}(X^{k+1}, Z^k) + \frac{1}{4} ||X^k - Z^k||^2.$$

By (2.12) - (2.14), the definition of the residual (2.15), it is clear that

**Proposition 3.1.3** If a sequence  $(X^k, Y^k, Z^k)$  generated by BADMM (2.17a) - (2.17c) converges to  $(X^*, Y^*, Z^*)$ , with  $X^* \in \mathcal{X}$ ,  $Z^* \in \mathcal{Z}$  and  $R_{BADMM}(k) \to 0$ , then  $(X^k, Y^k, Z^k)$  converges to a KKT point of the problem.

#### 3.2 Implementation and numerical experiments

We discuss briefly how the algorithms are implemented from a computational perspective. For all algorithms, the termination condition is set to be either (a) the total number of iterations exceeds

10000 or (b) the residual is less than  $10^{-5}$ . Note that for experiment purpose, we let each program calculates the residual during every iteration, which is time-consuming and may not be necessary in practice.

For ADMM with reformulation, we set  $\tau = 1.618$ . The vector  $\boldsymbol{x}_1^k$ ,  $\boldsymbol{\lambda}^k$  are stored as matrices of dimension  $n \times m$ , with each column representing  $x_i, \lambda_i \in \mathbb{R}^n$  respectively. In this way, updating all  $x_i$ 's can be done in efficient matrix operations without any inner loop.

For BADMM, with careful rearrangement, the updates can be done efficiently in compressed matrix operations. There is no inner loop inside the main loop.

For semi-proximal ALM, we set  $\tau = 1.9$ . The Cholesky factorization of  $AA^T$  is computed before the main loop for efficiently solving the linear systems during each iteration. In addition, several external mex files for efficient numerical linear algebra routines such as matrix-vector multiplication are called.

#### 3.2.1 Running statistics

Below is the running statistics for the 3 algorithms as input dimension ranges from m=n=10 to m=n=500. We omit the execution of ADMM with reformulation when  $m=n\geq 100$  since the patterns suggest that it does not converge after the maximum number of iterations. For each input dimension the main program generates the input data and calls the 3 algorithms (as Matlab functions). See Appendix A for the programs.

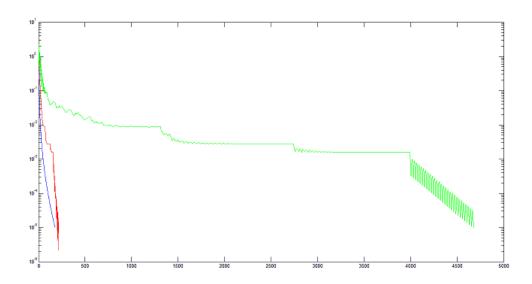
| Problem   | Gurobi optimal | BADMM opt_obj           | SPALM opt_obj           | ADMM LP reformulation     |
|-----------|----------------|-------------------------|-------------------------|---------------------------|
| dimension | objective      | it_count   time         | it_count   time         | opt_obj   it_count   time |
| (m=n)     |                |                         |                         |                           |
| 10        | 0.1337202778   | 0.133793   176   0.05   | 0.133676   216   0.03   | 0.133721   4676   0.78    |
| 20        | 0.1608133806   | 0.161327   520   0.13   | 0.160761   444   0.08   | 0.162288   10000   4.22   |
| 30        | 0.1856271350   | 0.186206   590   0.17   | 0.185599   473   0.08   | 0.187329   10000   9.51   |
| 40        | 0.1705273270   | 0.171129   506   0.16   | 0.170447   912   0.20   | 0.171518   10000   23.85  |
| 50        | 0.1576763389   | 0.158461   609   0.28   | 0.157586   805   0.27   | 0.159822   10000   76.39  |
| 75        | 0.1647900104   | 0.166104   855   0.61   | 0.164747   823   0.51   | 0.168612   10000   276.50 |
| 100       | 0.1605997205   | 0.161848   969   0.94   | 0.160652   1168   1.03  | 0.166659   10000   630.13 |
| 150       | 0.1464856269   | 0.148102   1265   2.25  | 0.146557   930   1.56   | -                         |
| 200       | 0.1603708148   | 0.162349   1518   4.49  | 0.160448   979   2.67   | -                         |
| 300       | 0.1658212714   | 0.168437   1950   15.97 | 0.165877   1085   6.68  | -                         |
| 500       | 0.1648862238   | 0.168237   2622   69.61 | 0.164768   1100   20.27 | -                         |

#### 3.2.2 Residual plots

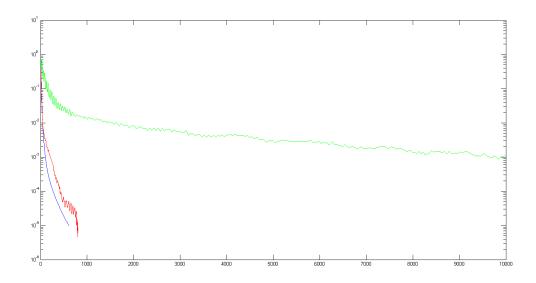
During each execution, the residuals are tracked and plotted in a log scale. Notice that for instances with small dimensions, BADMM converges slightly faster than semi-proximal ALM with respect to iteration count. When the dimension is large, semi-proximal ALM converges significantly faster.

ADMM with reformulation does not display favorable residual convergence pattern even after the maximum number of iterations.

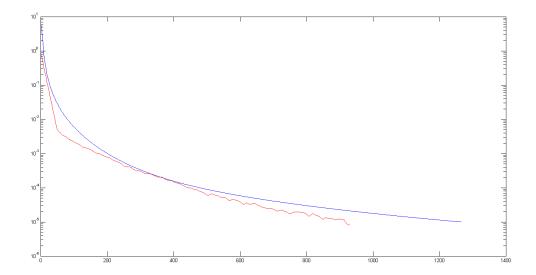
m=n=10: (Blue: BADMM; Red: SPALM; Green: ADMM with reformulation)



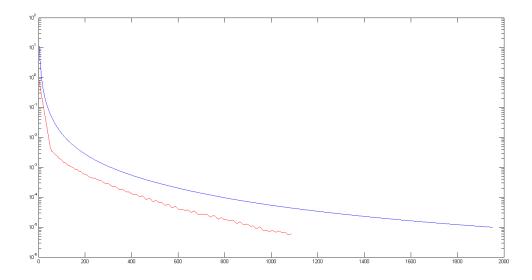
m=n=50



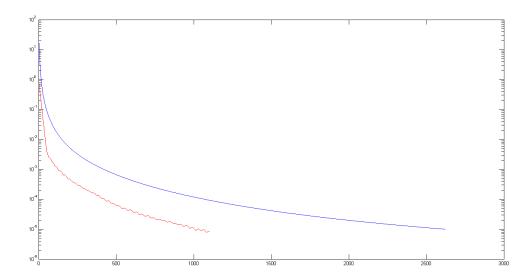
m=n=150 (ADMM with reformulation is not executed)



m=n=300 (ADMM with reformulation is not executed)



m=n=500 (ADMM with reformulation is not executed)



#### 3.2.3 Discussion

Experiment results suggest that the classical ADMM with LP reformulation is not a practical approach even on instances with small dimensions. Besides the observed sub-linear rate of convergence, the residual plot shows an undesirable oscillating pattern near the end of the main loop (m = n = 10, after the 4000-th iteration). In addition, from the running statistics table in 3.2.1 it can be observed that the time for each iteration increases at a rate higher than  $O(n^2)$ . This is primarily due to the

reformulation scheme, which constructs  $x_1 \in \mathbb{R}^{nm^2}$ , where m, n are the dimensions of the input cost matrix  $C \in \mathbb{R}^{m \times n}$ .

When the dimension is large, semi-proximal ALM outperforms BADMM both in terms of total number of iterations and total running time. Besides an observed faster rate of convergence (after the same number of iterations, the residual of semi-proximal ALM is much smaller), semi-proximal ALM also exploits efficient linear algebra routines during each iteration, although the program for BADMM is also highly optimized via using compressed matrix operations to replace inner loops. In theory, both algorithms require  $O(n \times m)$  floating point operations for each iteration.

As mentioned above, in practice, it is not necessary to calculate the objective value and residual during every iteration. When the dimension of the input instance is large, calculating the residual and objective value becomes relatively more time-consuming, since they always have the same order of time complexity as the main iterative schemes. An alternative is to calculate the residual and objective value after every fixed number of iterations.

One drawback of the above experiment is that the residuals of the algorithms are not consistently defined. The residual defined for BADMM measures the violation to the optimality condition (2.12) - (2.14), while the other two measure the violation to the optimality condition for the standard form LP problem, namely the KKT system (1.8a) - (1.8d). As such, setting the same terminal residual threshold may lead to biased results. However, for high-dimensional input instances, the experiment results suggest that semi-proximal ALM gives more accurate optimal objective as compared to BADMM in less number of iterations and less total running time. As such, the conclusion is not affected by these differences.

### Chapter 4

# Conclusion

We have shown that ADMM-type methods can be used for solving LP problems, with semi-proximal ALM and BADMM significantly outperforms the classical ADMM with reformulation for instances with dimensions ranging from m=n=10 to m=n=100. For instances with large dimensions (m=n=100 to m=n=500), semi-proximal ALM converges faster than the BADMM and terminates faster given the same terminal residual threshold of  $10^{-5}$ . Under this terminal residual threshold, semi-proximal ALM also gives more accurate objective values as compared to BADMM, relative to the corresponding optimal objectives given by Gurobi.

One thing to note is that the above BADMM iterative scheme (2.17a) - (2.17c) only works for the assignment problem, while semi-proximal ALM works on any standard form LP problem and does not depend on the structure of the assignment problem to achieve the demonstrated efficiency.

There have been numerous open problems on both theories and and applications of ADMM, including convergence analysis, alternative (changing) step lengths and multi-block extensions of the 2-block ADMM. Specific to the above algorithms, possible future work directions include formulating more general types of problems into the linearly constrained convex optimization framework (1.2) with generic iterative scheme (1.4a) - (1.4c), deriving efficient updating formulas that lead to efficient implementation. In addition, as pointed out in [8], for a specific optimization problem, the Bregman divergence chosen may lead to large p and hence a step length  $\tau$  that is too small to be practical. In fact, experiments therein show that setting a larger  $\tau$  which does not satisfy the conditions in Theorem 2.2.2 might still lead to a working algorithm. Hence it is reasonable to conjecture that the conditions can be relaxed via alternative proving techniques.

# References

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#### Appendix A: Matlab programs for the algorithms

Please refer to the files attached for the Matlab programs used for numerical experiments.

admm\_reformulated.m

This is the program for the classical ADMM with reformulation proposed by He and Yuan in [1] in Section 2.1.

badmm\_assignment\_problem\_solver.m

This is the program for the BADMM for assignment problem in Section 2.2.

semi\_proximal\_alm\_lp\_solver.m

This is the program for the semi-proximal ALM in Section 2.3.

main\_script\_comparing\_the\_algorithms.m

This is the script that generates random input instances, calculates the optimal objective value using Gurobi, calls the 3 programs and plots the residuals in log-scale.