# Notes on solver

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#### Abstract

Solver

## 1 Log barrier sub-problems

This paper is concerned with the following problem:

$$\min f(x) - \mu \log(s) + \frac{1}{2} d_x^T D_x d_x + \frac{1}{2} d_s^T D_s d_s$$
 (1a)

$$a(x) - s = r\mu \tag{1b}$$

$$s \ge 0 \tag{1c}$$

The KKT conditions for (1) are:

$$\nabla_x \mathcal{L}(x, y) = \nabla f(x) + D_x d_x - \nabla a(x)^T y = 0$$
(2a)

$$C_{\mu}(s,y) = Ys - \mu e = 0 \tag{2b}$$

$$\mathcal{P}_{\mu}(x,s) = a(x) - s - \mu r = 0 \tag{2c}$$

$$s, y \ge 0 \tag{2d}$$

Where the Lagrangian  $\mathcal{L}(x,y) := f(x) - y^T a(x)$ .

We combine the log barrier merit function and the complementary conditions as follows:

$$\phi(x,y) = \psi(x) + \zeta(x,y) \tag{3}$$

With:

$$\zeta(x,y) = \frac{\|\mathcal{C}(x,y)\|_{\infty}^{3}}{\mu^{2}}$$

We now introduces models to locally approximate these merit functions  $\nabla_x \mathcal{L}(x,y)$ ,  $\psi$ ,  $\mathcal{C}$  and  $\phi$  respectively. To describe our approximations of a function f around the point (x,y) we use the function  $\tilde{\Delta}_{(x,y)}^f(u,v)$  to denote the predicted increase in the function f at the new point (x+u,y+v). Observe that we use different approximations depending on the choice of function f.

We use a typical linear approximate of  $\nabla_x \mathcal{L}(x,y)$  as follows:

$$\tilde{\Delta}_{(x,y)}^{\nabla_x \mathcal{L}}(d_x, d_y) = \nabla_{x,x} L(x,y) d_x + \nabla a(x) d_y \tag{4}$$

The following function  $\tilde{\Delta}_{(x,y)}^{\psi}(u)$  is an approximation of the function  $\psi(x)$  at the point (x,y) and predicts how much the function  $\psi$  changes as we change the current from x to x+u.

$$\tilde{\Delta}_{(x,y)}^{\psi}(u) = \frac{1}{2}u^{T}M(x,y)u + \nabla\psi(x)^{T}u$$
(5)

With:

$$M(x,y) = \nabla^2 \mathcal{L}(x,y) + \sum_i \frac{y_i}{a(x)} \nabla a(x)^T \nabla a(x)$$
 (6)

Note that if we set  $y_i = \frac{\mu}{s_i}$  then  $M(x,y) = \nabla^2 \psi(x)$  and  $\tilde{\Delta}^{\psi}_{(x,y)}$  becomes the second order taylor approximation of  $\psi$  at the point x. Thus we can think of  $\tilde{\Delta}^{\psi}_{(x,y)}(u)$  as a primal-dual approximation of the function  $\psi$ .

We can also build a model of the  $\zeta(x,y)$  as follows:

$$\tilde{\Delta}_{(x,y)}^{\zeta}(d_x, d_y) = \frac{\|Sy + Yd_s + Sd_y - \mu e\|_{\infty}^3 - \|\mathcal{C}(x,y)\|_{\infty}^3}{\mu^2}$$
(7)

With S a diagonal matrix containing entries of a(x) and  $d_s = \nabla a(x)d_x$ . This model  $\tilde{\Delta}_{(x,y)}^{\mathcal{C}}$  corresponds to the typical primal-dual linear model of  $\mathcal{C}$  i.e.  $C(x+d_x,y+d_y)\approx Sy+Yd_s+Sd_y-\mu e$ .

With S and Y contain the diagonal elements of a(x) and y respectively.

This allows us to approximate the change in the function  $\phi$  at the point (x,y) as follows:

$$\tilde{\Delta}_{(x,y)}^{\phi}(d_x, d_y) = \tilde{\Delta}_{(x,y)}^{\psi}(d_x) + \tilde{\Delta}_{(x,y)}^{\zeta}(d_x, d_y)$$
(8)

We say an iterate (x,y) satisfies approximate complementary if  $(x,y) \in \mathcal{Q}_{\mu}$  where  $\mathcal{Q}_{\mu}$  is defined as follows:

$$Q_{\mu} = \left\{ (x, y) \in \mathbb{R}^{n} \times \mathbb{R}^{m} : a(x) > 0, y > 0, \|\mathcal{C}(x, y)\|_{\infty} \le \frac{\mu}{2} \right\}$$
 (9)

We say the point (x, y) is a  $\mu$ -scaled KKT point if  $(x, y) \in \mathcal{T}_{\mu}$  where:

$$\mathcal{T}_{\mu} = \{ (x, y) \in \mathcal{Q}_{\mu} : \|\nabla \mathcal{L}(x, y)\| \le \mu(\|y\|_1 + 1) \}$$
(10)

In which case the algorithm terminates.

## 2 Algorithm

Let S, Y denote the diagonal matrices with entries of s and y respectively. We can linearize (2) at the iterate (x, y, s) as follows:

$$\begin{bmatrix} \nabla^{2} \mathcal{L}(\hat{x}, \hat{y}) + D_{x} & -\nabla a(\hat{x})^{T} & 0 \\ \nabla a(\hat{x}) & 0 & -I \\ 0 & \hat{S} & \hat{Y} + D_{s} \end{bmatrix} \begin{bmatrix} d_{x} \\ d_{y} \\ d_{s} \end{bmatrix} = - \begin{bmatrix} \nabla \mathcal{L}(x, y) \\ \mathcal{P}_{\mu}(x, s) \\ \mathcal{C}_{\mu}(s, y) \end{bmatrix}$$
(11)

Which is equivalent to solving:

$$\begin{bmatrix} \nabla^2 \mathcal{L}(\hat{x}, \hat{y}) + \nabla a(x)^T D_s \nabla a(x) + D_x & \nabla a(\hat{x})^T \\ \nabla a(\hat{x}) & -(\hat{Y} + D_s)^{-1} \hat{S} \end{bmatrix} \begin{bmatrix} d_x \\ -d_y \end{bmatrix} = -\begin{bmatrix} \nabla \mathcal{L}(x, y) \\ \mathcal{P}_{\mu}(x, s) + (\hat{Y} + D_s)^{-1} \mathcal{C}_{\mu}(s, y) \end{bmatrix}$$
(12)

One can also solve this system by solving the Schur complement:

$$(\nabla^{2} \mathcal{L}(\hat{x}, \hat{y}) + \nabla a(\hat{x})^{T} (\hat{Y} + D_{s}) \hat{S}^{-1} \nabla a(\hat{x}) + D_{x}) d_{x} = -\nabla \mathcal{L}(x, \mu S^{-1} e) - \nabla a(\hat{x})^{T} \hat{Y} \hat{S}^{-1} \mathcal{P}_{\mu}(x, s)$$

Observe that (??) may be singular or correspond to a direction that makes the log barrier objective worse. To rectify this problem we compute the direction as follows:

$$d_x = \arg\min_{\|u\|_2 \le r} \tilde{\Delta}_{(x,y)}^{\psi}(u) \tag{13a}$$

$$d_s = \nabla a(x)d_x \tag{13b}$$

$$d_{y} = -S^{-1}(Yd_{s} + C(x, y))$$
(13c)

$$(M(x,y) + \delta I)d_x = -\nabla \psi(x) \tag{14}$$

Furthermore, by re-arranging this equation we can deduce that  $(d_x, d_y, d_s)$  satisfies a perturbed version of (??):

$$\begin{bmatrix} \nabla^2 \mathcal{L}(x,y) + \delta I & -\nabla a(x)^T & 0 \\ -\nabla a(x) & 0 & I \\ 0 & S & Y \end{bmatrix} \begin{bmatrix} d_x \\ d_y \\ d_s \end{bmatrix} = - \begin{bmatrix} \nabla \mathcal{L}(x,y) \\ 0 \\ \mathcal{C}(x,y) \end{bmatrix}$$
(15)

#### Algorithm 1 Primal-dual trust region step

function Primal-dual-trust-region(x, y, r) \*\*\*\*  $\in$  \*\*\*\*

$$d_x \in \arg\min_{\|u\| \le r} \tilde{\Delta}_{(x,y)}^{\psi}(u) \tag{16a}$$

$$d_s = \nabla a(x)d_x \tag{16b}$$

$$S = \operatorname{Diag}(a(x)) \tag{16c}$$

$$d_{y} = -S^{-1}(Yd_{s} + \mathcal{C}(x, y)) \tag{16d}$$

$$(x^+, y^+) \leftarrow (x + d_x, y + d_y)$$
  
return $(x^+, y^+, d_x, d_y)$   
end function

Our complete algorithm is summarized as follows:

#### Algorithm 2 Primal-dual non-convex interior point algorithm

```
\begin{array}{l} \textbf{function Non-convex-IPM}(x^1,y^1) \\ \textbf{for } k=1,...,\infty \ \textbf{do} \\ r \leftarrow R(y^k) \\ \textbf{repeat} \\ (x^+,y^+,d_x,d_y) \leftarrow \texttt{Primal-dual-trust-region}(x^k,y^k,r) \\ \textbf{if } (x^+,y^+) \in \mathcal{Q}_{\mu} \ \textbf{then} \\ \textbf{if } (x^+,y^+) \in \mathcal{T}_{\mu} \ \textbf{then} \\ \textbf{return } (x^+,y^+) \\ \textbf{end if} \\ \textbf{end if} \\ r \leftarrow r/2 \\ \textbf{until } \phi(x^+) > \phi(x^k) + \frac{1}{2} \tilde{\Delta}_{(x^k,y^k)}^{\phi}(d_x,d_y) \\ x^k \leftarrow x^+ \\ y^k \leftarrow y^+ \\ \textbf{end for} \\ \textbf{end function} \end{array}
```

# 3 Delta computation

#### Algorithm 3 Delta

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\lambda_{lb} = 0, \lambda_{ub} = \delta_{\max} = ||H||_F^2, \delta_{k-1} blower and upper bounds on minimum Try \delta = 0, if succeeds, trial solve with this delta. If step size is small skip to trust region step.
                                                                            ▷ lower and upper bounds on minimum eigenvalue
\delta = \delta_{k-1}
if \delta = 0 then
     \delta = \delta_{\min}
end if
for i = 1, ..., \infty do
    Break if inertia correct and update \lambda_{lb} and \lambda_{ub}.
end for
Trust region
R = \|d_x^{k-1}\|_2
for i = 1, ..., \infty do
     Compute trust region with R
    If trust region is too accurate increase radius size
    If step unsuccessful decrease radius size
    Prevent oscillation
end for
```