

# The Augmented Lagrangian Method

Lecture 14, Continuous Optimisation  
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We will now see a related method that does not require  $\mu_k$  to converge to zero, and yet in a neighbourhood of a KKT point  $x^*$  of the nonlinear optimisation problem

$$\begin{aligned} \text{(NLP)} \quad & \min_{x \in \mathbb{R}^n} f(x) \\ \text{s.t.} \quad & g_{\mathcal{E}}(x) = 0 \\ & g_{\mathcal{I}}(x) \geq 0, \end{aligned}$$

the iterates  $x_k$  still converge to  $x^*$  if the LICQ and the second order sufficient optimality conditions hold at this point. In fact,  $\mu$  can even be held constant after a while and the convergence of  $x_k$  continues!

In Lecture 13 we saw that the quadratic penalty method has the disadvantage that the penalty parameter  $\mu$  has to be reduced to very small values before  $x_k$  becomes feasible to high accuracy.

Moreover, we pointed out that reducing  $\mu$  to very small values can lead to numerical instabilities if the method is not implemented very carefully.

## Motivation:

The method is motivated by the observation that if we knew the Lagrange multipliers  $\lambda^*$  such that  $(x^*, \lambda^*)$  is a KKT point for (NLP), then we could find  $x^*$  by solving the unconstrained problem

$$\min_{x \in \mathbb{R}^n} \mathcal{L}(x, \lambda^*). \tag{1}$$

Indeed, as already remarked in Lemma 1.2 i) of Lecture 12, the first set of KKT conditions  $\nabla_x \mathcal{L}(x^*, \lambda^*) = 0$  amount to the first order necessary optimality conditions for (1).

Of course,  $\lambda^*$  is not known, but we know from Lecture 13 that one can obtain estimates  $\lambda^{[k]}$  which can be used to set up the problem

$$\min_{x \in \mathbb{R}^n} \mathcal{L}(x, \lambda^{[k]}).$$

as an approximation of (1).

If the estimates  $\lambda^{[k]}$  can be iteratively improved and made to converge to  $\lambda^*$ , then this can form the basis of an algorithmic framework for solving (NLP).

### Algorithm: Augmented Lagrangian Method (AL)

**S0** Initialisation: choose the following,

$x_0 \in \mathbb{R}^n$  (starting point, not necessarily feasible)

$\lambda^{[0]} \in \mathbb{R}^{|\mathcal{E} \cup \mathcal{I}|}$  (initial "guestimate" of Lagrange multiplier vector)

$\mu_0 > 0$  (initial value of homotopy parameter)

$(\tau_k)_{\mathbb{N}_0} \searrow 0$  (error tolerance)

### The Merit Function:

The merit function used by this algorithm is the *augmented Lagrangian* of (NLP), defined as follows,

$$\begin{aligned} \mathcal{L}_A(x, \lambda, \mu) &= \mathcal{L}(x, \lambda) + \frac{1}{2\mu} \sum_{i \in \mathcal{I} \cup \mathcal{E}} \tilde{g}_i^2(x) \\ &= f(x) - \sum_{i \in \mathcal{I} \cup \mathcal{E}} \lambda_i g_i(x) + \sum_{i \in \mathcal{I} \cup \mathcal{E}} \frac{\tilde{g}_i(x)}{2\mu} g_i(x) \\ &= f(x) + \sum_{i \in \mathcal{I} \cup \mathcal{E}} \left( \frac{\tilde{g}_i(x)}{2\mu} - \lambda_i \right) g_i(x), \end{aligned}$$

where  $\tilde{g}_i$  is defined as in Lecture 13,

$$\tilde{g}_i(x) = \begin{cases} g_i(x) & (i \in \mathcal{E}) \\ \min(g_i(x), 0) & (i \in \mathcal{I}). \end{cases}$$

**S1** For  $k = 0, 1, 2, \dots$  repeat

$$y^{[0]} := x_k, l := 0$$

until  $\|\nabla_x \mathcal{L}_A(y^{[l]}, \lambda^{[k]}, \mu_k)\| \leq \tau_k$  repeat

compute  $y^{[l+1]}$  such that  $\mathcal{L}_A(y^{[l+1]}, \lambda^{[k]}, \mu_k) < \mathcal{L}_A(y^{[l]}, \lambda^{[k]}, \mu_k)$

(using unconstrained minimisation method)

$$l \leftarrow l + 1$$

end

$$x_{k+1} := y^{[l]}$$

$$\lambda_i^{[k+1]} := \lambda_i^{[k]} - \frac{\tilde{g}_i(x_{k+1})}{\mu_k}, \quad (i \in \mathcal{E} \cup \mathcal{I}),$$

$$\lambda_i^{[k+1]} \leftarrow \max(0, \lambda_i^{[k+1]}), \quad (i \in \mathcal{I})$$

choose  $\mu_{k+1} \in (0, \mu_k)$

end

- By arguments similar to those in Theorem 2.2 in Lecture 13,

$$\lambda_i^{[k]} - \frac{\tilde{g}_i(x_{k+1})}{\mu_k} \simeq \lambda_i^*, \quad (i \in \mathcal{E} \cup \mathcal{I}).$$

- Therefore, we have

$$\tilde{g}_i(x_{k+1}) \simeq \mu_k (\lambda_i^{[k]} - \lambda_i^*), \quad (i \in \mathcal{E} \cup \mathcal{I}),$$

which suggests that if  $\lambda^{[k]} \rightarrow \lambda^*$  then all constraint residuals converge to zero like a function  $o(\mu_k)$ , where

$$\lim_{\mu \rightarrow 0} \frac{o(\mu)}{\mu} = 0.$$

That is, the convergence is much faster than the  $O(\mu_k)$  convergence obtained in the quadratic penalty function method.

A quick argument gives insight into why this method can be expected to converge before  $\mu_k$  reaches very small values:

- We have

$$\begin{aligned} \nabla_x \mathcal{L}_A(x_{k+1}, \lambda^{[k]}, \mu_k) &= \nabla f(x_{k+1}) - \sum_{i \in \mathcal{E} \cup \mathcal{I}} \left( \lambda_i^{[k]} - \frac{\tilde{g}_i(x_{k+1})}{\mu_k} \right) \nabla g_i(x_{k+1}). \end{aligned}$$

- Using  $\|\nabla_x \mathcal{L}_A(x_{k+1}, \lambda^{[k]}, \mu_k)\| \leq \tau_k$ , we find

$$\sum_i \left( \lambda_i^{[k]} - \frac{\tilde{g}_i(x_{k+1})}{\mu_k} \right) \nabla g_i(x_{k+1}) = \nabla f(x_{k+1}) + O(\tau_k).$$

**Theorem 1:** Let  $x^*$  be a local minimiser of (NLP) where the LICQ and the first and second order sufficient optimality conditions are satisfied for some Lagrange multiplier vector  $\lambda^*$ . Then there exists a constant  $\bar{\mu} > 0$  such that  $x^*$  is a strict local minimiser of

$$\min_{x \in \mathbb{R}^n} \mathcal{L}_A(x, \lambda^*, \mu)$$

for all  $\mu \in (0, \bar{\mu}]$ .

Theorem 2: For  $(x^*, \lambda^*)$  and  $\bar{\mu}$  as in Theorem 1 there exist constants  $M, \varepsilon, \delta > 0$  such that the following is true:

i) If  $\mu_k \leq \bar{\mu}$  and

$$\|\lambda^{[k]} - \lambda^*\| \leq \frac{\delta}{\mu_k}, \quad (2)$$

then the constrained minimisation problem

$$\begin{aligned} \min_x \mathcal{L}_A(x, \lambda^{[k]}, \mu_k) \\ \text{s.t. } \|x^* - x\| \leq \varepsilon \end{aligned} \quad (3)$$

has a unique minimiser  $x_{k+1}$ ,

and furthermore,

$$\|x^* - x_{k+1}\| \leq M\mu_k \|\lambda^{[k]} - \lambda^*\|, \quad (4)$$

ii) if  $\mu_k$  and  $\lambda^{[k]}$  are as in part i) and if  $\lambda^{[k+1]}$  is chosen as in Algorithm (AL), then

$$\|\lambda^{[k+1]} - \lambda^*\| \leq M\mu_k \|\lambda^{[k]} - \lambda^*\|. \quad (5)$$

Some remarks about this result:

- (3) suggests the use of a trust-region method in the inner loop of Algorithm (AL).
- Without loss of generality, we may assume that  $\bar{\mu} \leq (2M)^{-1}$ . Note that if  $(\lambda^{[k]}, \mu_k)$  satisfy the conditions of part i) of the theorem,

$$\begin{aligned} \text{I)} \quad & \mu_k \leq \bar{\mu}, \\ \text{II)} \quad & \|\lambda^{[k]} - \lambda^*\| \leq \frac{\delta}{\mu_k}, \end{aligned}$$

and if it is also the case that

$$\text{III)} \quad x_k \in B_\varepsilon(x^*),$$

then  $x_k$  is a feasible starting point for the constrained problem

$$\begin{aligned} \min_x \mathcal{L}_A(x, \lambda^{[k]}, \mu_k) \\ \text{s.t. } \|x^* - x\| \leq \varepsilon. \end{aligned}$$

Furthermore, we have

$$\begin{aligned} \text{I')} \quad & \mu_{k+1} \leq \mu_k \leq \bar{\mu}, \\ \text{II')} \quad & \|\lambda^{[k+1]} - \lambda^*\| \stackrel{\text{II), (5)}}{\leq} M\mu_k \frac{\delta}{\mu_k} = \delta M < \frac{\delta}{\bar{\mu}} \leq \frac{\delta}{\mu_{k+1}}, \\ \text{III')} \quad & x_{k+1} \in B_\varepsilon(x^*). \end{aligned}$$

Hence, by induction the relations I), II) and III) hold at every subsequent iteration  $j$  and the assumptions of part i) remain valid.

- Let  $k_0$  be the iteration where (4) and (5) first hold,

$$\begin{aligned}\|x^* - x_{k+1}\| &\leq M\mu_k \|\lambda^{[k]} - \lambda^*\|, \\ \|\lambda^{[k+1]} - \lambda^*\| &\leq M\mu_k \|\lambda^{[k]} - \lambda^*\|.\end{aligned}$$

Then induction on  $k$  shows that

$$\|\lambda^{[k]} - \lambda^*\|, \|x_k - x^*\| \leq (M\bar{\mu})^{k-k_0} \|\lambda^{[k_0]} - \lambda^*\| \leq \frac{1}{2^{k-k_0}} \|\lambda^{[k_0]} - \lambda^*\|.$$

Therefore,  $x_k \rightarrow x^*$  and  $\lambda^{[k]} \rightarrow \lambda^*$  at a Q-linear rate if  $\mu \leq \bar{\mu}$  is held fixed.

**Reading Assignment:** Lecture-Note 14.

**Recommended Additional Reading:** Section 17.4, Nocedal–Wright.