Numerical Optimization 08: Quasi-Newton methods

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May 14, 2020

Overview

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- BFGS method
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Quasi-Newton's method

Just as the secant method approximates f " in the univariate case, quasi Newton approximate the inverse Hessian $((\mathbf{H}^k)^{-1})$ which is needed for each step of update

$$oldsymbol{x}^{k+1} \leftarrow oldsymbol{x}^k - lpha^k (oldsymbol{H}^k)^{-1} oldsymbol{g}^k$$

These methods typically set $(\mathbf{H}^k)^{-1}$ (let's call it \mathbf{Q} from now on) to the identity matrix and then apply updates to reflect information learned with each iteration. To simplify the equations for various quasi-Newton methods, we define the following

$$egin{aligned} &m{\gamma}^{k+1} = m{g}^{k+1} - m{g}^k \ &m{\delta}^{k+1} = m{x}^{k+1} - m{x}^k \end{aligned}$$

A new quadratic model

Instead of computing the exact Q, we can update it in a simple manner to account for the curvature measured during the most recent step. Suppose, we have generated x^{k+1} and wish to construct a new quadratic model,

$$m^{k+1}(\boldsymbol{p}) = f(x^{k+1}) + \boldsymbol{g}^{k+1}\boldsymbol{p} + \frac{1}{2}\boldsymbol{p}^T \boldsymbol{Q}^{k+1}\boldsymbol{p}$$

We let the gradient of m^{k+1} match the gradient of f for at least two steps x^{k+1} and x^k .

$$\nabla m^{k+1}(-\alpha^k p^k) = \boldsymbol{g}^{k+1} - \alpha^k \boldsymbol{Q}^{k+1} \boldsymbol{p}^k = \boldsymbol{g}^k$$

Since $\nabla m^{k+1}(0) = g^{k+1}$, the second of these condition is satisfied automatically. Rearranging it, we obtain the so called secant condition.

$$\boldsymbol{Q}^{k+1}\alpha^{k}\boldsymbol{p}^{k} = \boldsymbol{g}^{k+1} - \boldsymbol{g}^{k} \quad \rightarrow \quad \boldsymbol{Q}^{k+1}\boldsymbol{\delta}^{k} = \boldsymbol{\gamma}^{k} \tag{1}$$

A new quadratic model

Given the displacements δ^k and the change of gradients γ^k . It requires that the symmetric positive definite matrix Q^{k+1} , it needs that

$$\delta^k \gamma^k > 0$$

At this stage, there still exists an infinite number of solutions of Q^{k+1} . To determine a unique solution, we impose another condition, which is that Q^{k+1} is close to the current Q^k

$$\label{eq:stable} \begin{split} \min_{oldsymbol{Q}} & \|oldsymbol{Q} - oldsymbol{Q}^k\| \ \mathrm{s.t.} \quad oldsymbol{Q} = oldsymbol{Q}^T, \quad oldsymbol{B} \delta^k = \gamma^k \end{split}$$

Different matrix norms can be applied here to give different quasi-Newton methods.

The Davidon-Fletcher-Powell (DFP) method

Davidon proposed the following relation between $oldsymbol{Q}^k$ and $oldsymbol{Q}^{k+1}$

$$\boldsymbol{Q}^{k+1} = \boldsymbol{Q}^k + a u u^T + b v v^T$$

According to the secant condition

$$oldsymbol{Q}^k \delta^k + extbf{a} extbf{u} extbf{U}^T \delta^k + extbf{b} extbf{v} extbf{V}^T \delta^k = oldsymbol{\gamma}^k$$

An obvious choice for u and v is

$$u = \gamma^k, \qquad v = \boldsymbol{Q}^k \boldsymbol{\delta}^k \quad o \quad a u^T \boldsymbol{\delta}^k = 1, \quad b v^T \boldsymbol{\delta}^k = -1$$

where

$$a = 1/u^{\mathsf{T}} \delta^k = 1/u^{\mathsf{T}} \delta^k \quad b = -1/v^{\mathsf{T}} \delta^k = 1/u^{\mathsf{T}} \delta^k$$

$$oldsymbol{Q}^{k+1} = oldsymbol{Q}^k - rac{oldsymbol{Q}^k \gamma^k (\gamma^k)^T oldsymbol{Q}^k}{(\gamma^k)^T oldsymbol{Q}^k \gamma^k} + rac{\delta(\delta^k)^T}{(\delta^k)^T \gamma^k}$$

W. C. Davidon, Variable Metric Method for Minimization SIAM Journal on Optimization. 1. (1991), 1-17.

The Broyden-Fletcher-Goldfarb-Shanno (BFGS) method

In the BFGS algorithm, it does not approximate ${oldsymbol Q}^k$, but handles ${oldsymbol H}^k = {oldsymbol Q}^{k-1}$

$$oldsymbol{H}^{k+1}oldsymbol{\gamma}^k=oldsymbol{\delta}^k$$

The minimize condition is,

$$\min_{\boldsymbol{H}} ||\boldsymbol{H} - \boldsymbol{H}^{k}||$$
s.t. $\boldsymbol{H} = \boldsymbol{H}^{T}, \quad \boldsymbol{H} \boldsymbol{\gamma}^{k} = \boldsymbol{\delta}^{k}$

$$oldsymbol{Q}^{k+1} = oldsymbol{Q}^k + rac{\delta \gamma^T oldsymbol{Q} + oldsymbol{Q} \gamma \delta^T}{\delta^T \gamma} + igg(1 + rac{\gamma^T oldsymbol{Q} \gamma}{\delta^T oldsymbol{Q}}igg) rac{\delta \delta^T}{\delta^T \gamma}$$

BFGS does better than DFP with approximate line search.

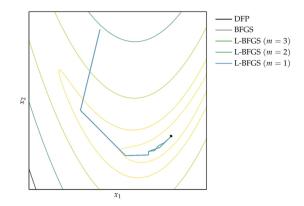
Limited-memory BFGS

BFGS still uses an $n \times n$ dense matrix, which is a problem for storage of the hessian when dealing with very large scale problems. The L-BFGS method can be used to approximate BFGS with a relatively cheaper solution. In L-BFGS, it stores the last *m* values for δ and γ rather than the entire inverse of *H*.

$$oldsymbol{Q} \leftarrow oldsymbol{Q} - rac{\delta \gamma^{oldsymbol{T}} oldsymbol{Q} + oldsymbol{Q} \gamma \delta^{oldsymbol{T}}}{\delta^{oldsymbol{T}} \gamma} + igg(1 + rac{\gamma^{oldsymbol{T}} oldsymbol{Q} \gamma}{\delta^{oldsymbol{T}} oldsymbol{Q}}igg) rac{\delta \delta^{oldsymbol{T}}}{\delta^{oldsymbol{T}} \gamma}$$

BFGS does better than DFP with approximate line search but still uses an $n \times n$ dense matrix.

Comparison of various quasi-Newton algos



Summary

- Quasi-Newton method attempted to approximate the Hessian from function and gradient evaluations.
- The first step approximation of hessian in the quasi-newton methods is usually an identity matrix
- BFGS performs better than DFP, but it still relies on the storage of big Hessian matrix
- L-BFGS is a more scalable approach for large scale problems.
- All quasi-Newton methods can work with the approximate line search.