A Stochastic Quasi-Newton Optimizer for TensorFlow

Jason Chen

Department of Computer Science Stony Brook University Stony Brook, NY 11790 hungrchen@cs.stonybrook.edu

David Kraemer

Department of Applied Mathematics Stony Brook University Stony Brook, NY 11790 davidkraemer@stonybrook.edu

Abstract

We do some shit and it doesn't work.

1 Introduction

Limited memory variations of classical second-order minimization algorithms can prove highly attractive in applications with high-dimensional problems or when the cost of each iteration step is prohibitive under these classical formulations. Specifically, the canonical second-order technique, Newton's method, involves explicit computation of the Hessian matrix of the objective function, whose time complexity grows with the square of the parameter dimension. Even with BFGS, which seeks to reduce the direct computational difficulty of exactly constructing the Hessian itself, encounters similar difficulties of scale. The limited memory approach is to bound this overall cost and ensure that approximate Hessians can be computed on existing hardware.

We implement a variant of stochastic second-order algorithm as described by Wang, Ma, Gold-farb, and Liu [3] for application in the TensorFlow library [2]. TensorFlow is a framework for graph-based computation models. Its primary use within the statistical learning community is in neural networks, which can be readily implemented by the library. After the model is constructed, the actual training is performed by an Optimizer object. In this setting, the Optimizer is presented with a partially observable objective function and gradient from which to perform optimization. Some typical Optimizer subclasses which are available in the Python API include an GradientDescentOptimizer¹ and an AdagradOptimizer²

The presently available optimization algorithms implemented in the Python library are either first or zeroeth-order methods, requiring access to either the objective function or its gradient. The conspicuous absence of second-order methods, such as Newton or quasi-Newton algorithms, has been a subject of much discussion inside the TensorFlow community³

Following the suggestions of the community discussion, we present an implementation of Stochastic damped limited memory BFGS (SdLBFGS) for TensorFlow through the use of an ExternalOptimizerInterface. Presently used to wrap existing minimization routines implemented by outside libraries such as SciPy, it serves as a simple way to engage with the TensorFlow internals while avoiding a deep dive into its architecture.

Our work proceeds as follows. Section 2 reviews the problem proposed in Wang et al and gives an explicit formulation of SdLBFGS according to their paper. We identify small modifications made to the core pseudocode and explain their necessity. Section 3 describes our SdLBFGS implementation

¹See source.

²See source.

³The primary discussion has occurred surrounding Issue #446, in which contributors have noted that the current infrastructure for developing second-order methods is lacking and any implementation would require considerable re-inventing the wheel.

and how it interfaces with TensorFlow. Section 4 shows preliminary experimental results on the correctness of our implementation on simple problems. Importantly, we report that our present implementation is incapable of giving meaningful progress on training practical models on common datasets. In Section 5, we discuss the implications of this negative result on future development and specify areas in which our project may be continued and improved.

Source code for this project can be found at https://github.com/DavidNKraemer/cse592-project. Documentation, cruft removal, and user interface improvements are currently evolving.

2 Review of theory

- 1. Review main results of the Wang, Ma, Goldfarb, Liu paper [3]
- 2. State the SdLBFGS algortihm.
- 3. Mention some details about how the algorithm translates to code.

The SdLBFGS algorithm is proposed by Wang, Ma, Goldfarb, and Liu [3] for solving nonconvex stochastic optimization problems. In particular, it solves

$$\begin{array}{ll}
\text{minimize} \\
x \in \mathbb{R}^n
\end{array} f(x) = \mathbb{E}[F(x,\xi)] \tag{1}$$

for $F \in C^1(\mathbb{R}^n \times \mathbb{R}^d)$ and $\xi \in \mathbb{R}^d$ is a random variable with cumulative distribution function P. Here $\mathbb{E}[\cdot]$ denotes the mathematical expectation. In general F need not be convex.

The function F is readily interpreted in the context of supervised statistical learning. Given a corpus Ξ of data, we put a distribution on Ξ which reflects a sampling regime for a training set. The object of the learning process is, then, to the model parameter $x \in \mathbb{R}^n$ which minimizes F in expectation with respect to the training set. Provided that the sampling maintains some resemblance to the overall corpus, this gives a useful approximate parameter for the whole corpus.

The SdLBFGS algorithm emerges from two concurrent optimization techniques. First, it follows in the tradition of approximate second-order algorithms which seek to perform a version of Newton's method while avoiding the computation of the Hessian $\nabla^2 f(x)$ explicitly. The classical BFGS approach is the canoncal quasi-Newton method. Second, it follows the development of stochastic algorithms which seek to randomly sample the components of the gradient $\nabla f(x)$ so that the essential convergence properties hold in expectation. In this sense it succeeds Stochastic Gradient Descent, among other canonical methods of this class. The limited memory constraint has antecedents in both streams of optimization techniques.

Wang et al. [3] assume that Problem (1) satisfies the following assumptions:

- 1. There is a real lower bound for f. That is, $\inf_{x \in \mathbb{R}^n} f(x) > -\infty$.
- 2. The gradient ∇f is Lipschitz continuous. That is, there exists L>0 such that

$$\|\nabla f(x) - \nabla f(y)\| \le L\|x - y\|$$

for all $x,y\in\mathbb{R}^n$. Moreover, $\nabla_{xx}F(x,\xi)$ exists and is continuous and there exists $\kappa>0$ with $\|\nabla^2_{xx}F(x,\varepsilon)\|\leq \kappa$ for any x,ξ .

3. The sampling regime is unbiased. That is, for the kth iterate the following hold:

$$\mathbb{E}[g(x_k, \xi_k)]_{\xi_k} = \nabla f(x_k),$$

$$\mathbb{E}[\|g(x_k, \xi_k) - \nabla f(x_k)\|^2]_{\xi_k} \le \sigma^2,$$

where $g(x, \xi_k) = \nabla F(x, \xi_k)$ and σ^2 is the noise level of estimating $g(x, \xi_k)$.

For a convex function, $\nabla^2 f(x)$ is always positive semidefinite (see Boyd and Vandenberghe §3.1.4 [1]), but this need not be the case in for nonconvex functions. The work in Wang et al. [3] shows how the BFGS approximate Hessian may be tweaked to ensure that positive semidefiniteness is preserved even in nonconvex cases. The overall stochastic quasi-Newton method is described in Algorithm 2, while the detailed implementation of SdLBFGS is given in Algorithm 2.

Algorithm 1 The high level stochastic quasi-Newton minimization algorithm. Given an initial point $x_1 \in \mathbb{R}^n$, batch sizes $\langle m_k \rangle$, step sizes $\langle \alpha_k \rangle$, and a maximum memory capacity p, perform BFGS with the stochastic dampened direction update.

```
1: function SQN(x_1, \langle m_k \rangle_{k=1}^{\infty}, \langle \alpha_k \rangle_{k=1}^{\infty}, p)
 2:
            data \leftarrow \emptyset
 3:
            for k = 1, 2, ... do
                  g_k \leftarrow \frac{1}{m_k} \sum_{i=1}^{m_k} g(x_k, \varepsilon_{k,i})
 4:
 5:
                  \Delta x_k, s_{k-1}, \bar{y}_{k-1}, \rho_{k-1} \leftarrow \text{SDLBFGS}(g_k, *(\text{data}^\top))
                                                                                                                                 6:
                  x_{k+1} \leftarrow x_k - \alpha_k \Delta x_k
                  append(\mathrm{data},(s_{k-1},\bar{y}_{k-1},\rho_{k-1}))
 7:
 8:
                  if k > p then
                        pop(data)
 9:
                  end if
10:
            end for
11:
12: end function
```

Algorithm 2 The SdLBFGS update step. The resulting output is $H_k g_k = v_p$, where H_k is the approximation of the kth iterate Hessian and g_k is the approximation of the kth iterate gradient. Note that the computation of H_k is implicit, preventing additional storage requirements.

```
1: function SDLBFGS(g_{k-1}, \langle s_j \rangle_{j=k-p}^{k-2}, \langle \bar{y}_j \rangle_{j=k-p}^{k-2}, \langle \rho_j \rangle_{j=k-p}^{k-2})
                     \text{mem} \leftarrow \min(p, k-1)
  2:
                   \begin{aligned} & s_{k-1} \leftarrow x_k - x_{k-1} \\ & y_{k-1} \leftarrow \frac{1}{m_{k-1}} \sum_{i=1}^{m_{k-1}} [g(x_k, \xi_{k-1,i}) - g(x_{k-1}, \xi_{k-1,i})] \\ & \gamma_k \leftarrow \max\left(\frac{y_{k-1}^\top y_{k-1}}{s_{k-1}^\top y_{k-1}}, \delta\right) \geq \delta \\ & \theta \leftarrow \max\left(\frac{3}{4} \frac{\gamma_k^{-1} s_{k-1}^\top s_{k-1}}{\gamma_k^{-1} s_{k-1}^\top s_{k-1} + y_{k-1}}, 1\right) & > \theta \end{cases} \\ & \bar{y}_{k-1} \leftarrow \theta y_{k-1} + (1 - \theta \gamma_k^{-1} s_{k-1}) \end{aligned}
  3:
                                                                                                                                                                         \triangleright \theta preserves positive semidefiniteness.
  7:
                     \rho_{k-1} \leftarrow (s_{k-1}^{\mathsf{T}} \bar{y}_{k-1})^{-1}
  8:
                     for i = 0, ..., \text{mem} - 1 do
  9:
                               \mu_i \leftarrow \rho_{k-i-1} u_i^{\top} s_{k-i-1}
10:
11:
                               u_{i+1} \leftarrow u_i - \mu_i \bar{y}_{k-i-1}
                     end for
12:
                     v_0 \leftarrow {\gamma_k}^{-1} u_{\text{mem}}
13:
                     \begin{array}{l} \textbf{for } i = 0, \dots, \text{mem} - 1 \textbf{ do} \\ \nu_i \leftarrow \rho_{k-\text{mem}+i} v_i^{\top} \bar{y}_{k-\text{mem}+i} \end{array}
14:
15:
                                v_{i+1} \leftarrow v_i + (\mu_{\text{mem}-i-1} - \nu_i) s_{k-\text{mem}+i}
16:
17:
                     end for
18:
                     return v_p, s_{k-1}, \bar{y}_{k-1}, \rho_{k-1}
19: end function
```

3 Implementation

- 1. Describe how TensorFlow algorithms are usually implemented.
- 2. Contrast with our implementation of SdLBFGS.
- 3. Explain some implementation issues.

4 Empirical results

- 1. Results on simple optimization problems.
- 2. Results on potentially more difficult optimization problems.
- 3. Lack of results on MNIST with TensorFlow.

5 Discussion

- Explaining why the basic project implementing a useful practical TensorFlow optimzier
 — failed.
- 2. Takeaways from the results.
- 3. Future work.

References

- 1. Boyd, S. & Vandenberghe, L. *Convex Optimization* ISBN: 0521833787 (Cambridge University Press, New York, NY, USA, 2004).
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- 3. Wang, X., Ma, S., Goldfarb, D. & Liu, W. Stochastic Quasi-Newton Methods for Nonconvex Stochastic Optimization. *SIAM Journal on Optimization* **27**, 927–956 (2017).