# Curiously Fast Convergence of some Stochastic Gradient Descent Algorithms

#### Léon Bottou

#### January 23, 2009

### 1 Context

Given a finite set of m examples  $z_1, \ldots, z_m$  and a strictly convex differentiable loss function  $\ell(z, \theta)$  defined on a parameter vector  $\theta \in \mathbb{R}^d$ , we are interested in minimizing the cost function

$$\min_{\theta} \quad C(\theta) = \frac{1}{m} \sum_{i=1}^{m} \ell(z_i, \theta).$$

One way to perform such a minimization is to use a stochastic gradient algorithm. Starting from some initial value  $\theta[1]$ , iteration t consists in picking an example z[t] and applying the stochastic gradient update

$$\theta[t+1] = \theta[t] - \eta_t \frac{\partial \ell}{\partial \theta} \ell(z[t], \theta[t])$$

where the sequence of positive scalars  $\eta_t$  satisfies the well known Robbins-Monro conditions  $\sum_t \eta_t = \infty$  and  $\sum_t \eta_t^2 < \infty$ . We consider three ways to pick the example z[t] at each iteration:

- *Random* Examples are drawn uniformly from the training set at each iteration.
- Cycle Examples are picked sequentially from the randomly shuffled training set, that is,  $z[km + t] = z_{\sigma(t)}$ , where  $\sigma$  is a random permutation of  $\{1, \ldots, m\}$ , and k is a nonnegative integer, and  $t \in \{1, \ldots, m\}$ .
- Shuffle Examples are still picked sequentially but the training set is shuffled before each pass, that is,  $z[km + t] = z_{\sigma_k(t)}$ , where the  $\sigma_k$  are random permutations of  $\{1, \ldots, m\}$ , and k is a nonnegative integer, and  $t \in \{1, \ldots, m\}$ .

With suitable assumptions on the function  $\ell$ , the random case can be treated with well known stochastic approximation results [1, 5]. With gains of the form  $\eta_t = c/(t+t0)$  and sufficiently large values of the constant c, one obtains results such as

$$\mathbb{E}\left[C(\theta[t]) - \min_{\theta} C(\theta)\right] \sim \frac{1}{t}$$

where the expectation is taken over the random choice of examples at each iteration. Various theoretical works [2, 3, 6] indicate that no choice of  $\eta_t$  can lead to faster convergence rates than  $t^{-1}$ .

#### 2 Experiments

We report now empirical results obtained with the three method.

The task is the classification of RCV1 documents belonging to class CCAT [4]. Each of the 781,265 examples is a pair composed of a 47,152 dimensional vector  $x_i$  representing a document and a variable  $y_i = \pm 1$  representing its appartenance to the class CCAT. The parameter vector  $\theta$  is also a 47,152 dimensional vector and the loss function is

$$\ell(x, y, \theta) = \log\left(1 + e^{-y(\theta.x)}\right)$$
.

All experiments were achieved using a variant of the svmsgd2 program and datasets.<sup>1</sup> The only modification consists in implementing our three schemes for selecting examples at each iteration.

Figure 1 shows log-log plots of the evolution of  $C(\theta[t])$  as a function of the number of iterations. The slope of the curve indicates the exponent of the convergence of the algorithm.

- The random case displays a  $t^{-1}$  convergence as predicted by the stochastic approximation theory.
- The cycle case displays a  $t^{-\alpha}$  convergence with  $\alpha$  significantly greater than one. This means that this example selection strategy leads to a faster convergence. The exact value of  $\alpha$  changes when we consider different permutations of the examples.
- The *shuffle* case displays a more chaotic convergence. A linear interpolation of the curve leads to an exponent  $\alpha$  that is curiously close to

<sup>&</sup>lt;sup>1</sup>http://leon.bottou.org/projects/sgd.

two, suggesting that we have an average  $t^{-2}$  convergence. This result is stable when we repeat the experiment with different permutations of the training set.

## 3 The Question

In light of the theoretical works associated with stochastic approximations, stochastic algorithms that converge faster than  $t^{-1}$  are very surprising.

In fact, the stochastic approximation results rely on randomness assumption on the successive choice of examples are independent. Both the *cycle* and the *shuffle* break these assumptions but provide a more even coverage of the training set.

What can we prove for the *cycle* and the *shuffle* cases?

### References

- [1] A. Benveniste, M. Metivier, and P. Priouret. Algorithmes adaptatifs et approximations stochastiques. Masson, 1987.
- [2] K. Chung. On a stochastic approximation method. Annals of Mathematical Statistics, 25(3):463–484, 1954.
- [3] V. Fabian. On asymptotic normality in stochastic approximation. Annals of Mathematical Statistics, 39(4):1327–1332, 1968.
- [4] D. D. Lewis, Y. Yang, T. G. Rose, and F. Li. RCV1: A new benchmark collection for text categorization research. J. Machine Learning Research, 5:361–397, 2004.
- [5] L. Ljung and T. Söderström. Theory and Practice of recursive identification. MIT Press, Cambridge, MA, 1983.
- [6] P. Major and P. Revesz. A limit theorem for the robbins-monro approximation. Zeitschrift f
  ür Wahrscheinlichkeitstheorie und verwandte Gebiete, 27:79–86, 1973.



Figure 1: Evolution of  $C(\theta[t])$  for our three example selection strategies. The horizontal axe counts the number of epoch. One epoch represents 781,265 iterations, that is, one pass over the training set.