

# Adaptive Consistency Management for Distributed Machine Learning

by

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Submitted to the Department of Electrical Engineering and Computer  
Science

in partial fulfillment of the requirements for the degree of

Master of Science in Computer Science and Engineering

at the

TECHNISCHE UNIVERSITÄT BERLIN

April 2017

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## Abstract

In recent years, machine learning emerged as the most important part of many successful applications and businesses. While the rise of artificial intelligence continues, the amount of data to be processed grows even faster. Unsurprisingly a lot of research has since focused on methods to parallelize commonly used algorithms and new systems and frameworks have been published to improve the performance of distributed machine learning. Even though there has been a lot of progress towards more efficient systems, many state-of-the-art systems still have limitations. Current frameworks are neither expressible nor flexible enough to allow efficient development of distributed machine learning algorithms, which makes them unsuited for experimentation and quick prototyping even though this is an essential part to optimal performance. On the other hand, most parallelization strategies exploit the algorithms inherent stochastic nature to allow for parallel execution at the expense of lowered consistency among the distributed data structures. Even though this does not necessarily affect quality of the results, an improper chosen level of consistency can severely affect algorithm performance, resulting in a non optimal convergence rate and therefore increased runtime.

This thesis introduces a novel framework for distributed machine learning, which is based on a state centric programming model with yield semantics. The programming model allows the user to focus on the key elements of developing distributed machine learning algorithms, namely update communication, update merging and consistency management among distributed workers, while the system takes care of distributing the computation in an optimal fashion. The experiments using elastic-net regularized linear regression show an increased performance compared to state-of-the-art data processing systems like Apache Spark and at the same time reduce the effort of developing, parallelizing and experimenting with distributed machine learning algorithms at scale.

Thesis Supervisor: Prof. Dr. Odej Kao  
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# Acknowledgments

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# Chapter 1

## Introduction

One of the most challenging tasks in computer science and engineering resolves around improving algorithm performance. In general this has been done by making hardware faster and inventing new strategies and algorithms to parallelize work more efficiently. Since it is clear that Moore's Law will not hold anymore, a lot of effort has been spent to horizontally scale algorithm computation across multiple machines. Machine learning is no exception and efficient parallelization is a key aspect towards more intelligent systems. By now, many general purpose frameworks for large scale data processing have been developed and published. Many of those are used for running more complex machine learning algorithms at scale as well. Unfortunately, the performance often is not satisfying due to the architecture and programming model not reflecting the underlying structure of most commonly used machine learning algorithms. Common data processing tasks can be represented as an extract-transform-load (ETL) pipeline, which is often easily parallelizable. This does not hold for machine learning, where algorithms are mostly sequential in nature and the only way of enabling parallel computation is by exploiting their inherent stochastic properties. This allows to break the sequential execution in favor of parallel learning on subsets, which then needs to be combined in order to obtain a global solution to the task. While this can lead to a great speedup in terms of the amount of data processed, it can have a negative affect on the learning progress. Therefore a key part of horizontally scaling machine learning algorithms is to ensure all participating learners have a consistent



view on each others progress while at the same time maintaining a tradeoff between communicating progress and spending time on their own learning task.

## 1.1 MapReduce and Beyond

Many of todays successful businesses throughout fields like finance, e-commerce and healthcare rely heavily on the ability to process vast amounts of user or sensorical data, collected to make services smarter and user experience better. In order to learn from the collected data, discover patterns and ultimately gain new insights, it needs to be processed by an algorithm. It is not uncommon that the input ranges somewhere between hundred gigabytes and tenth of petabytes. In the past, processing this much data required either a supercomputer, which was only available to large institutions or government entities or some proprietary compute cluster. All this changed when Google introduced MapReduce [1] in 2004. The MapReduce framework made it possible to process data in a distributed and fault tolerant way with the help of a compute cluster formed by hundredth or thousandth of machines. Instead of using a single, expensive, special hardware supercomputer the framework provides the foundation to assemble commodity hardware machines into a compute cluster. The framework takes care of all necessary aspects to ensure a fault tolerant and parallel execution of a task submitted to the cluster. The advantage compared to previous approachs is that the framework can be run entirely on top of machines using commodity hardware, which does not require special hardware and therefore equals low cost.

MapReduce esentially led the path to a convenient and widespread use of big data processing, which found its open source implementation in the Apache Hadoop project [2]. The project quickly gained traction and has spawned many business grade platforms, which quickly gained widespread adoption and by now provides a whole ecosystem around big data processing. The software stack includes a fault tolerant distributed filesystem (HDFS) a MapReduce framework and a cluster resource manager (YARN) [3]. On the other hand, MapReduce suffers from some practical

limitations that lead to the development of new, more sophisticated and specialised big data frameworks. With the most widely used frameworks being Apache Spark [4], Apache Flink [5] and GraphLab [6]. The first two frameworks use at its core a dataflow pipeline based architecture, whileas the latter uses a graph abstraction to model particular algorithms. All this works well for algorithms that can be expressed as an extract-transform-load (ETL) pipeline and are often embarrassingly parallel in nature. On the other hand, machine learning algorithms often rely heavily on many, computationally light, iterations to iteratively update a shared state (model) such as logistic regression or latent dirichlet allocation (LDA). These so called iterative-convergent algorithms required a change in how systems for distributed machine learning operate at its core.

## 1.2 Distributed Machine Learning

This limitation essentially sparked the development of specialized frameworks for distributed execution of iterative-convergent algorithms used in common machine learning tasks. The most widely recognized systems are Petuum [7], ParameterServer [8] and MALT [9]. Different from the MapReduce paradigm, these frameworks, instead of using a pipeline to transform an immutable dataset into another immutable dataset, operate on a fixed but mutable state, which is held by a single machine or distributed over multiple machines. This state can then be updated by workers that have computed an update locally and send it to these state keepers. Which act as surrogates, taking state updates and incorporating these into the state by some user defined function. While these systems can increase the performance on machine learning algorithms by an order of magnitude [7] compared to dataflow systems, most systems come with either limited usability, which makes it difficult to implement additional algorithms, are tied to a specific algorithm or are very low level frameworks. Efficiently distributing machine learning algorithms while at the same time provide the ability to conscisely express machine learning algorithms remains an extremely challenging problem. A system targeting the execution of those algorithms at scale must

therefore provide the ability to concisely express the underlying algebraic structure and at the same time be flexible enough to allow experimentation and fine tuning. Where consistency management is an essential part to ensure that algorithms are executed fast and efficient.

## 1.3 Thesis Outline

In this thesis we introduce a novel framework for large scale distributed machine learning. It improves upon currently available systems by providing a powerful programming abstraction that can concisely express state of the art machine learning algorithms and at the same time minimizes the effort necessary to move from a single machine to a cluster. The framework design is optimized for efficient parallel asynchronous execution of iterative-convergent algorithms in a cluster and ensures the required consistency is enforced among parallel learners, depending on the algorithm properties. By allowing the user to decide how to maintaining the best tradeoff between algorithm execution and progress communication the performance is improved as well. All of this can be easily customized for quick prototyping and finetuning, which makes the system suited for developers as well as researchers. The goal of this thesis is to implement the state centric programming model and show its performance in comparison with Apache Spark on an example implementation of the CoCoA [10] framework. I start off by providing a background on the architecture and inner workings of state of the art frameworks for big data processing and distributed machine learning in Chapter 2. Furthermore the most commonly used algorithms and optimization techniques are introduced. The majority of those algorithms can be classified into the group of so called iterative-convergent algorithms for which I am going to provide a more formal description. I introduce the common algorithm parallelization strategies in distributed machine learning. I conclude the chapter by providing an overview over the challenges and issues that need to be addressed and considered when developing a distributed machine learning system and how this is achieved by current frameworks. This will give rise to understanding why a differ-

ent framework architecture and abstraction is necessary to improve the performance and expressability of large scale distributed machine learning applications. Chapter 3 therefore introduces the state centric programming model, which treats the state as a mutable first class citizen, which can be distributed and altered by updates that result from distributed computation. This allows the system to reason about the most optimal distribution of state in the cluster which is then scheduled with computation that can update the state. I further describe the architecture of the system and how the essential components are implemented. When updating a shared state from multiple different locations, consistency must be maintained in order to ensure the algorithm progresses as expected. The system is responsible for managing a state's consistency among all of its instances across the cluster. Chapter 4 therefore describes several schemes for ensuring consistency and at the same time optimizing for bandwidth and computational resource usage. In order to show the system and its consistency management in action, Chapter 5 compares the system against Apache Spark by running the CoCoA framework on two datasets with elastic-net regularized linear regression as the chosen algorithm. Chapter 6 summarize the experiments with the lessons learned and provides suggestions on how to further improve systems for large scale distributed machine learning.

# Chapter 2

## Background

This section provides the necessary background to follow the argumentation in the following chapters regarding state centric programming model and consistency management. This includes an understanding of algorithms and optimization techniques commonly used in practice (not limited to distributed machine learning), the current state-of-the-art in dataflow systems and how dataflow is used to provide a fault tolerant and distributed framework for large scale data processing and machine learning. Furthermore the field of distributed machine learning is explained in more detail, including the current state-of-the-art frameworks used for this purpose, their limitations and challenges that arise when machine learning algorithms are parallelized in a distributed fashion among multiple physical machines.

### 2.1 Algorithms and Optimization

#### 2.1.1 Iterative Convergent Algorithms

Consider a supervised learning setup with a dataset  $D = \{z_1, \dots, z_n\}$  with each example  $z_i$  being represented by a pair  $(x_i, y_i)$  consisting of an input  $x_i$  and a scalar output  $y_i$ . Consider also a loss function  $\ell(\hat{y}, y)$  quantifying the cost of predicting  $\hat{y}$  when the true output is  $y$ . As a model, a family  $F$  of functions  $f_w(x)$  parameterized by a weight vector  $w$  is chosen. The goal is to find a function  $f \in F$  that minimizes the

loss  $Q(z, w) = \ell(f_w(x), y)$ . Empirical risk  $E_n(f) = \frac{1}{n} \sum_{i=0}^n \ell(f(x_i), y_i)$  performance on training set, expected risk generalization performance.

$$E_n(f_w) = \frac{1}{n} \sum_{i=0}^n \ell(f_w(x_i), y_i) \quad (2.1)$$

In order to find an optimal solution many algorithms used in large scale machine learning such as regression, topic models, matrix factorization or neural networks employ either gradient based methods or markov chain monte carlo methods. To obtain the optimal solution those algorithms try to iteratively update the weight vector  $w$ . At each iteration  $t$  an updated weight vector  $w^t$  is computed based on the vector of the previous iteration  $w^{(t-1)}$  and the data  $D$ . The resulting model  $f_{w^t}$  is again a better summary of the data  $D$  under the objective  $Q$ . 2.2 shows the process of refining the model, with  $\Delta$  being an arbitrary update function.

$$w^t = w^{(t-1)} + \Delta(w^{(t-1)}, D) \quad (2.2)$$

The update function depends on the algorithm employed and can be viewed as a procedure of obtaining a step towards a better model. At each iteration an update  $\Delta w$  is computed and applied to the previous weight vector until a stopping condition is satisfied. E.g. the distance to the optimal solution or the objective difference between two iterations is monitored. When the difference is below a certain threshold the computation stops and the algorithm is said to be converged.

### 2.1.2 Convex Optimization

In order to estimate the optimal parameters  $w^*$  of a function belonging to class  $f_{w^*} \in F$ , numerous techniques can be employed to estimate said parameters. In many cases, especially large scale machine learning methods such as (stochastic) gradient descent and coordinate ascent are used to iteratively optimized the parameterization of the chosen function class. Both techniques represent different rules of computing the update shown in 2.2. Gradient descent updates the weights  $w$  at each iteration  $t$

on the basis of the gradient of  $E_n(f_w)$ ,

$$w^t = w^{(t-1)} - \eta \frac{1}{n} \sum_{i=0}^n \nabla_w Q(z_i, w^{(t-1)}) \quad (2.3)$$

where  $\eta$  is a chosen gain, often referred to as learning rate. While this achieves linear convergence under sufficient regularity assumptions and a sufficiently small learning rate  $\eta$  [11] [12] a more simplified version is commonly used in practice, called stochastic gradient descent (SGD). Instead of computing the gradient  $\nabla_w E_n(f_w)$  exactly, the gradient is estimated at each iteration  $t$  based on a single randomly picked example  $z_t$ .

$$w^t = w^{(t-1)} - \eta_t \nabla_w Q(z_t, w^{(t-1)}) \quad (2.4)$$

The assumption is that the gradient obtained by 2.4 behaves similar to its expectation in 2.3. The convergence properties have been studied extensively and under mild conditions an almost sure convergence can be established when the learning rate satisfies the conditions  $\sum_t \eta_t^2 < \infty$  and  $\sum_t \eta_t = \infty$  [12]. The general structure of stochastic gradient descent is described in Algorithm 1.

---

**Algorithm 1** Stochastic Gradient Descent

---

- 1:  $k \leftarrow 1$  and initialize  $w^0 \in \mathbb{R}^d$
  - 2: **repeat**
  - 3:     **for do**
  - 4:          $w^t \leftarrow w^{(t-1)} - \eta_t \nabla_w Q(z_t, w^{(t-1)})$
  - 5:     **until** termination criteria satisfied
- 

Coordinate descent on the other hand iteratively tries to optimize a given objective by successively performing approximate minimization along a coordinate direction while keeping the other directions fixed. TODO: give a more in detail explanation and provide algorithm description

### 2.1.3 Regularization

$$Q(z, w) = \ell(f_w(x), y) + r(w) \quad (2.5)$$

---

**Algorithm 2** Stochastic Coordinate Ascent

---

```
1:  $k \leftarrow 1$  and initialize  $w^0 \in \mathbb{R}^d$ 
2: repeat
3:   for do
4:
5: until termination criteria satisfied
```

---

$$Q(z, w) = \ell(f_w(x), y) + r(w) \quad (2.6)$$

$$Q(z, w) = \ell(f_w(x), y) + r(w) \quad (2.7)$$

NOTES: - regularization in general - L1, L2 regularization - elastic-net as special case - convexity properties - mini-batch setup algorithm

### 2.1.4 CoCoA

Due to their widespread application in large scale machine learning and recent advances in the field of distributed optimization the thesis focuses on linear regularized objectives. The theoretical contemplation as well as the experiments in Section 5 focus on a framework for convex optimization problems called CoCoA (Communication-efficient distributed dual Coordinate Ascent) [10] and its successors CoCoA<sup>+</sup>[13] and PROXCoCoA<sup>+</sup>[14]. CoCoA as described in Algorithm 3 provides a communication-efficient framework for solving convex optimization problems of the following form

$$\min_{\alpha} Q(z, \alpha) = \ell(f_{\alpha}(x), y) + r(\alpha) \quad (2.8)$$

in a distributed setting. Where  $\alpha$  denotes the weight vector,  $\ell$  is convex and smooth and  $r$  is assumed to be separable, which in this context means  $r(x) = \sum_{i=0}^n r_i(x_i)$ . Commonly the term  $\ell$  is an empirical loss over the data of the form  $\sum_i \ell(f_w(x_i), y_i)$  and the term  $r$  is a regularizer, e.g.  $r(w) = \lambda \|w\|_p$  where  $\lambda$  is a regularization parameter. Many algorithms in machine learning can be expressed in this form, such as logistic and linear regression, lasso and sparse logistic regression and support vector



machines.

CoCoA leverages the primal-dual relation which allows for solving the problem in either the primal or dual formulation. For some application where the number of examples  $n$  is much smaller than the number of features  $d$ , *nwaylessratherthan*, solving the problem in the dual may be easier because this problem has  $n$  variables to optimize, compared to  $d$  for the primal. The CoCoA framework leverages Fenchel-Rockafellar duality to quadratically approximate the global objective in 2.8. This leads to separability of the problem over the coordinates of  $\alpha$  and the partitions, where the local subproblems are similar in structure to the global problem and also exploit second order information within the local data partition. Therefore the dataset  $D \in \mathbb{R}^{d \times n}$  can be distributed either example-wise or feature-wise over  $K$  physical machines according to the partitioning  $\{P\}_{k=1}^K$ , depending which is more efficient to solve. The size of the partition on machine  $k$  is denoted by  $n_k = |P_k|$ . The key to efficient distributed optimization is that the local subproblems can be solved independently on each worker in parallel and only a single parameter vector  $v = \nabla f(D\alpha) \in \mathbb{R}^n$  needs to be shared after each round in order to communicate the progress of each local worker on its subproblem. As the data stays local and only a single parameter vector of dimension  $n$  needs to be exchanged, CoCoA is very communication efficient and as the local subproblems are very similar to the global problem, arbitrary solvers can be employed as well. The local quadratic subproblem has the following form

$$\min_{\alpha_{[k]} \in \mathbb{R}^n} \zeta_k^{\sigma'}(\Delta\alpha_{[k]}, v, \alpha_{[k]}), \quad (2.9)$$

where

$$\zeta_k^{\sigma'}(\Delta\alpha_{[k]}, v, \alpha_{[k]}) = \frac{1}{K} f(v) + w^T A_{[k]} \Delta\alpha_{[k]} + \frac{\sigma'}{2\tau} \|A_{[k]} \Delta\alpha_{[k]}\|^2 + \sum_{i \in P_k} g_i(\alpha_i + \Delta\alpha_{[k]_i}), \quad (2.10)$$

with  $A_{[k]}$  referring to the the local data and  $w = \nabla f(D\alpha) \in \mathbb{R}^n$ .  $\alpha_{[k]_i}$  denotes the local weight vector stored on machine  $k$  with  $\alpha_{[k]_i} = \alpha_i$  if  $i \in P_k$  otherwise  $\alpha_{[k]_i} = 0$ . In summary the framework provides a procedure to effectively combine the results

---

**Algorithm 3** CoCoA Framework

---

**Data:** Data matrix  $D$ , distributed column-wise according to partition  $\{P_k\}_{k=1}^K$

**Input:** aggregation parameter  $\gamma \in (0, 1]$ , subproblem parameter  $\sigma'$

**Initialize:**  $t \leftarrow 0$ ,  $\alpha^{(0)} \leftarrow 0 \in \mathbb{R}^n$ ,  $v^{(0)} \leftarrow 0 \in \mathbb{R}^d$

```
1: repeat  
2:    $t \leftarrow t + 1$   
3:   for  $k \in \{1, \dots, K\}$  do  
4:     obtain an approximate solution  $\Delta\alpha_{[k]}$  for the local subproblem 2.9  
5:     update local weights  $\alpha_{[k]}^t = \alpha_{[k]}^{t-1} + \gamma\Delta\alpha_{[k]}$   
6:     update shared parameter vector  $\Delta v_k = A_{[k]}\Delta\alpha_{[k]}$   
7:   compute  $v^t = v^{t-1} + \gamma \sum_{k=1}^K \Delta v_k$   
8: until termination criteria satisfied
```

---

from local computation without having to deal with conflicts resulting from similar updates computed on other machines.

NOTES: - local solver (batch coordinate descent) - elastic-net in the cocoa setup  
- algorithm parameters sigma and gamma

## 2.2 Dataflow Systems

NOTES: - description of dataflow in general - architecture of dataflow systems, Apache Spark and Flink - description of RDDs

## 2.3 The Challenges of Distributed Machine Learning

While the execution of iterative-convergent algorithms on a single machine is straightforward, time constraints and the ever growing amount of data to be processed require these algorithms to be executed in parallel. This poses a number of challenges which are often observed when parallelizing algorithms, such as partitioning the state used in the algorithm as well as communication and consistency management. In this context, state refers to a structure containing arbitrary data e.g. an array, tensor or list. While intra-node parallelism in multi-core and multi-processor systems can mitigate these problems, it is not satisfying in terms of cost and scalability. On the other hand, inter-node parallelism has the desired properties but can not be easily parallelized.

Therefore in the past decade a lot of research has focused on inventing new systems to deal with those challenges and make distributed machine learning more efficient and scalable.

### 2.3.1 State Partitioning

Depending on the algorithm and optimization technique employed to solve for the optimal solution, there exist multiple approaches to distribute the state used in the algorithm (e.g. input data, model parameters). As depicted in Figure 2-1, an initial approach by Zinkevich et. al [15] introduced a data-parallel approach for computing stochastic gradient descent (SGD) via the MapReduce framework. In this context,

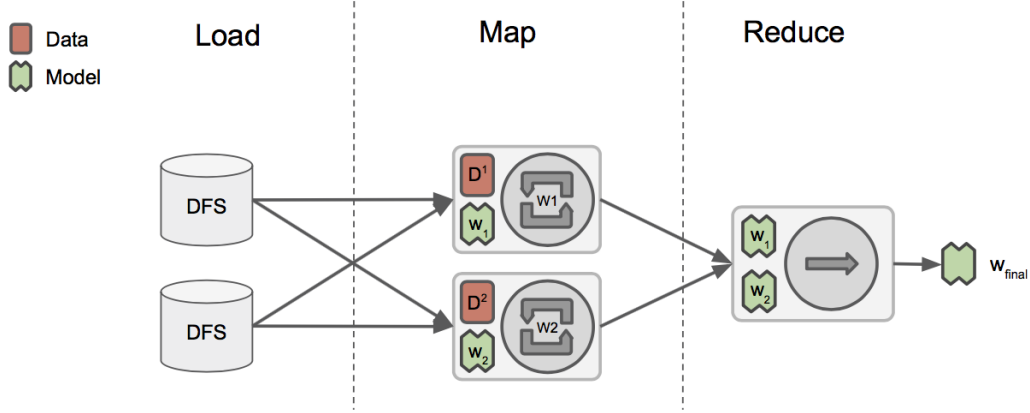


Figure 2-1: Data-parallelism via MapReduce

data-parallelism means that  $K$  machines work in parallel on the input data  $D$ , hence the data is distributed according to partition  $\{P_k\}_{k=1}^K$  into local parts  $D^k$ . Each machine maintains a local model  $w_k$  that is iteratively refined based on 2.2 until convergence, using only the local part of the input data  $D^k$ . The local update is of the form

$$w_k^t = w_k^{t-1} + \Delta(w_k^{t-1}, D^k) \quad (2.11)$$

The final model  $w_{final}$  is then obtained in the reduce step by averaging over all local models  $w_k$ .

$$w_{final} = \frac{1}{K} \sum_{k=1}^K w_k \quad (2.12)$$

Even though this approach works well in practice and gives considerable good results, it suffers from two limitations. First, if the size of a local model  $w_k$  exceeds the available memory on a single machine, the algorithm can not work properly. This is the case e.g. for topic models at web scale or deep neural networks used in the Google Brain project [16], consisting of billions of parameters. Second, even though the scalability improved by introducing data-parallelism the lack of parameter exchange during runtime can lead to suboptimal performance [17] as the algorithm essentially resembles batch gradient descent, which is known to have suboptimal convergence properties compared to mini-batch or stochastic gradient descent [12] [14]. In order to improve the performance, a system must be able to communicate more frequently and it must also be able to partition a model accross multiple machines to scale with the size of the model. Following those requirements resulted in the publication of the parameter server [8], which provides a framework for inter-node parallelism of iterative convergent algorithms.

### 2.3.2 Communication

As depicted in Figure 2-2, the parameter server is a group of an arbitrary number of machines  $\{S\}_{l=1}^L$ , e.g.  $S = \{S_1, S_2\}$  where each member of the group is responsible for storing a part of the model  $\{w\}_{l=1}^L$  and making it accessible to the workers  $\{W\}_{k=1}^K$ , e.g.  $W = \{W_1, W_2, W_3\}$ , via a defined interface that is similar to a key-value store. The model is partitioned among the machines of the server to provide optimal throughput, fault tolerance and to mitigate the effect of a model exceding the memory of a single machine. Each of the workers maintains a local partition of the input data  $\{D\}_{k=1}^K$ , which is used to iteratively compute updates for the parameters  $w$  according to

$$\Delta w_{k_i}^t = \Delta(w_{k_i}^{t-1}, D^k). \quad (2.13)$$

In the parameter server setup, the local state  $w$  acts as a cache for global parameters in order to reduce network usage. Therefore, depending on the caching policy,  $w_{k_i}$  can either be directly read either the local cache or must be retrieved from the parameter

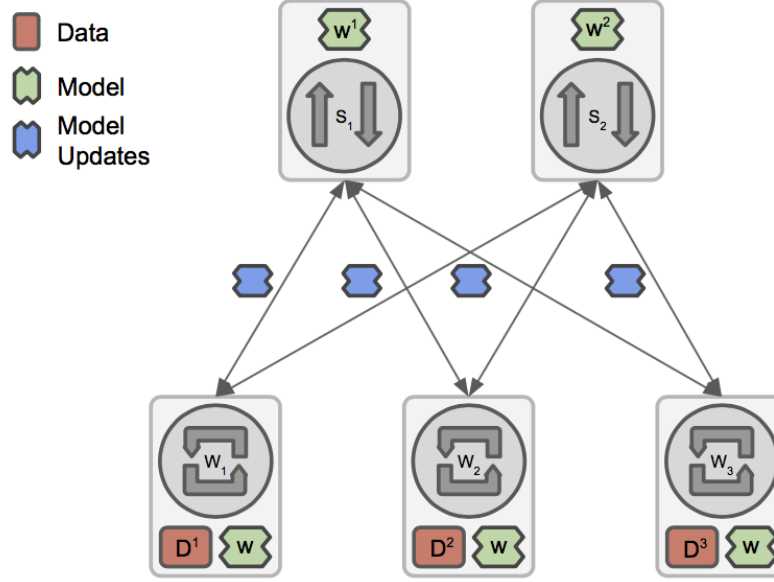


Figure 2-2: Parameter Server

server. Additionally to applying the update to the local model  $w$ , the difference  $\Delta w_{k_i}^t$  is published to the parameter server as well, which takes care of applying it to the corresponding entry  $w_i$  to make the update available to all other workers  $W_i$ ,  $i \neq k$ . In case multiple updates for the same parameter  $w_i$  arrive at the same time, a user defined function (UDF) needs to be provided to the server, which takes care of combining those updates so it can be applied to the parameter stored on the server. The procedure of retrieving, updating and publishing is executed concurrently on all workers. This enables all workers to work in parallel on the iterative parameter refinement while asynchronously updating and retrieving the parameters necessary for computing the next update. The procedure of updating the model in parallel is also known as model-parallelism.

While this schema has been proven to work well in practice (some numbers), a couple of issues still remain when running iterative-convergent algorithms at scale using the parameter server. According to [18] this can be viewed as finding the trade-off between algorithm throughput and data throughput. In other words the challenge is to find a balance between the quality of parameter refinements and the quantity at which they are generated. As with distributed systems in general, network com-

communication is the bottleneck in distributed machine learning as well. Even though due to the stochastic nature of many machine learning algorithms, the communication can be reduced compared to the exact serial algorithm, it has been proven that fresher model parameters increase the algorithm throughput per iteration [19]. Therefore, in order to guarantee an optimal algorithm throughput, a worker should always work with the latest parameters. On the other hand, exchanging parameter updates over the network more frequently is time consuming, which leaves less time to run local computation and therefore essentially decreases the data throughput due to increased time spent on network management. The second issue concerns the relaxed consistency among participating workers due to reduced communication compared to the serial algorithm. Though this is what makes the parameter server concept so powerful because it increases the data throughput by relaxing the consistency requirements when updating parameters on the server. By decoupling the progress of workers it is possible to minimize the effect of stragglers and synchronization delay between workers [20]. However, as discussed in the next section, combining model parameters obtained from workers with greatly differing algorithm progress can have a detrimental effect on algorithm throughput. Finding the balance between algorithm throughput and data throughput can therefore be seen as consistency management.

### 2.3.3 Consistency

The most important part of any distributed system is the synchronization strategy used to ensure consistency among multiple machines concurrently accessing and updating some shared state. In distributed machine learning the shared state is the model, which is for example stored in a parameter server and continuously refined by updates locally computed by a worker and then published to the server. There are three schemes used to synchronize workers during the iterative parameter refinement. *Bulk synchronous parallelization (BSP)* leads to the best algorithm throughput (convergence achieved per number of examples processed). In this scheme, all workers are required to finish their current iteration and at the end successfully publish all updates to the parameter server. The server then computes a refined model  $w^t$  ac-

cording to 2.14 and each worker retrieves the updated parameters before beginning the next iteration. This synchronization scheme guarantees consistency among all nodes at all times.

$$w^t = w^{t-1} + \frac{1}{K} \sum_{k=1}^K \Delta(w_k^{t-1}, Dk) \quad (2.14)$$

While this synchronization strategy essentially recovers the sequential algorithm for a single machine and has the same convergence properties and guarantees, it suffers from a severe limitation when used in a distributed setup [19]. In case one of the workers is for some reason a lot slower than the others the synchronization barrier imposed by BSP forces all workers to wait for this particular worker in the group. This is well known as the straggler problem [20] and can seriously affect performance in a distributed environment, because the progress is limited by the slowest node in the cluster. BSP is commonly used in MapReduce frameworks such as Hadoop and data-flow systems like Apache Spark and Apache Flink to ensure correct program execution. The second strategy is known as *total asynchronous parallelization (TAP)*. Similar to BSP, all workers publish their locally computed parameter updates to the server after each iteration but in this case the changes are applied to the model immediately. Therefore no waiting for other workers is required, resulting in a very high data throughput. The straggler problem can be mitigated by this synchronization scheme as well, as depicted in Figure 2-3. Even though worker  $W_3$  is a

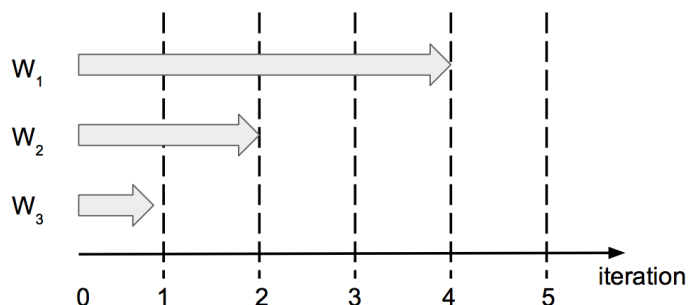


Figure 2-3: Straggler in TAP

straggler, which would have prevented the remaining workers  $\{W_2, W_3\}$  from proceeding beyond the synchronization barrier of iteration 1, the workers can continue with their next iterations without waiting for the slower worker. Although this consistency

scheme seems to work quite well in practice [8], it lacks formal convergence guarantees and can even diverge [21]. This stems from the fact that no theoretical convergence bound can be established as the divergence in iteration between workers is unbound. A middle ground between bulk synchronous parallelization and total asynchronous

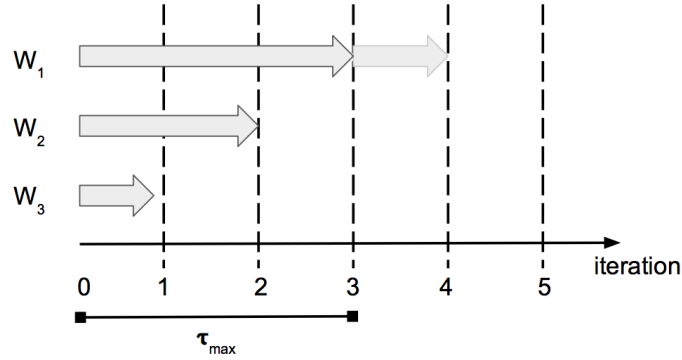


Figure 2-4: Straggler in SSP

parallelization is *stale synchronous parallel (SSP)* [22] or *bounded staleness (BS)*. As shown in Figure 2-3, SSP introduces a fixed maximum delay, or staleness threshold, of  $\tau_{max}$  between the slowest and fastest node. In the example, for  $\tau_{max} = 3$  worker  $W_1$  is blocked and can not proceed beyond iteration 3 as the slowest worker  $W_3$  has not finished its first iteration. As soon as worker  $W_3$  has completed its first iteration,  $W_1$  is unblocked and can proceed as long as the difference in iteration stays below  $\tau_{max}$ . SSP overcomes some of the limitations of TAP by introducing a bound on divergence in number of iterations between workers. The staleness threshold resembles a bound which can be used to restore formal convergence guarantees while still maintaining the flexibility of asynchronous parallelization and limiting but not completely preventing the straggler problem [23]. In general this helps to compensate e.g. update related communication between iterations or fluctuations in worker performance, which explains why SSP works so well in practice.



## Chapter 3

# State Centric Programming Model

Most state-of-the-art frameworks for distributed machine learning like Petuum [7] and Parameter Server [8] are based on the parameter server concept introduced in the previous section. The framework essentially provides a low level API<sup>1</sup> for publishing and retrieving values similar to a distributed key-value store, where the key  $i$  is for example the index of a weight vector  $w$  stored on the server and the value is the weight  $w_i$ . Implementing an algorithm that relies on a parameter server requires to incorporate the publishing and retrieval of parameters deeply into the algorithm. This contrasts the general workflow of developing and testing an algorithm locally on a single machine and then transition to a distributed environment. The parameter server paradigm provides only a minor abstraction, leaving the developer with the task of distributing state, scheduling distributed computation, consistency management and managing cluster resources. A developer should be able to focus on the key elements to efficient parallel execution of distributed machine learning algorithms. In the following section these key elements of a framework that improves upon currently available systems are derived by the help of an example based on elastic-net regularized linear regression. The result is a state centric programming model, treating state as a first class citizen that can be distributed and altered by local and remote transformation.

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<sup>1</sup>Application programming interface

### 3.1 Control Flow

Figure 3-1 shows the logical control flow of a general pipeline for training an arbitrary iterative-convergent machine learning algorithm. Usually training starts with load-

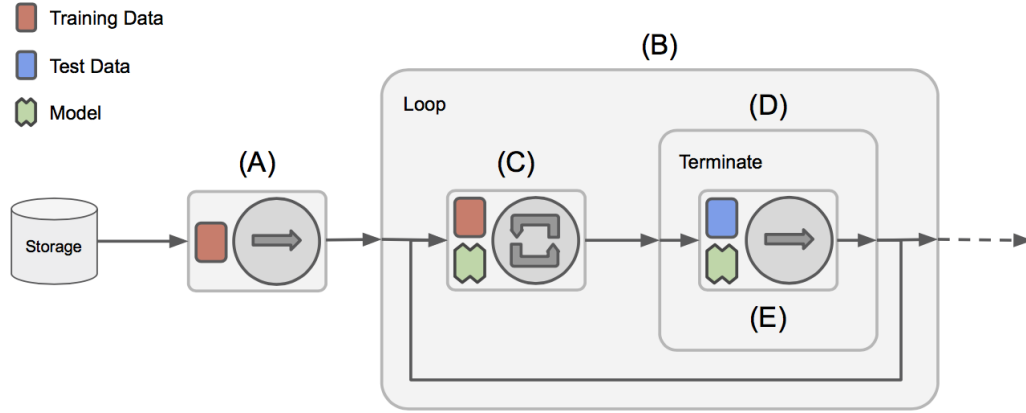


Figure 3-1: Machine Learning Pipeline for iterative-convergent algorithms

ing the input data  $D$  from the storage followed by one or more preprocessing steps (e.g. normalization or standardization) as well as splitting the data into training and test set **(A)**. A square represents a step of an algorithm, which contains an arbitrary number of input states (e.g. data, model) and some kind of transformation applied to the input states. The transformation process is depicted as a circle and can either be applied once (arrow) or multiple times (cyclic arrows) to the input state(s) during this particular step. After applying the preprocessing the actual training process is triggered, which is commonly contained in a loop **(B)**. A loop symbolises that the containing steps are executed repeatedly until some termination criterion is satisfied, which is computed in **(C)**. For most machine learning algorithms the termination criterion can either be a fixed number of iterations, the change in objective  $Q$  between iterations or the generalization performance **(E)**. **(D)** is the actual training step which iteratively refines the model by updates computed from the input data according to 2.2. It can already be seen from the example that a framework for distributed machine learning must be capable of concisely expressing a complex sequence of arbitrary control flow operators and transformation steps. An algorithm expressed in this form can be executed as is on a single machine without modification but in order to distribute

it among a cluster of machines each transformation step must be parallelized as can be seen in Figure 3-2. **NOTE:** add caption for state, make termination more generic, it's better to call this control flow The pipeline described in Figure 3-1 is parallelized

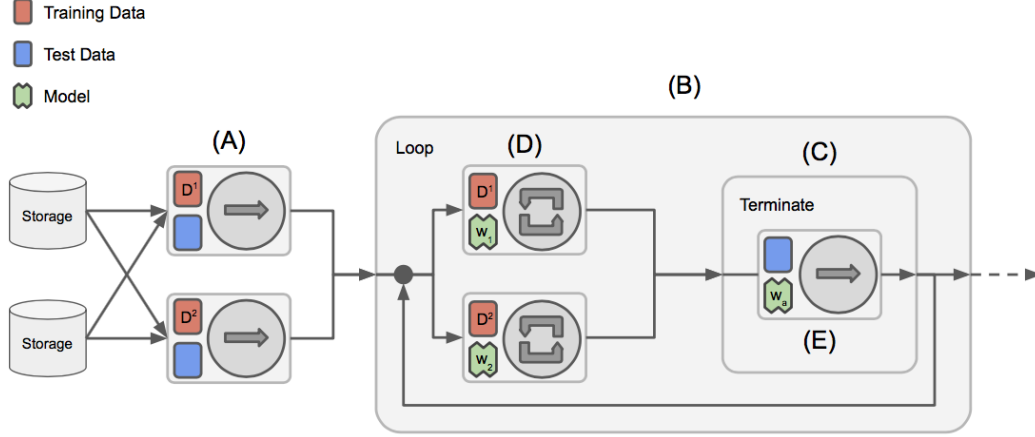


Figure 3-2: Distributed Machine Learning Pipeline for Iterative-convergent Algorithms

with a degree of parallelism of two. Now each transformation step of the pipeline can be scheduled and executed on two machines in parallel. It is worth noting that the transformation logic stays the same independently of the  $\text{dop}^2$ , including a  $\text{dop}$  of one, which is essentially a single machine. Only the input and output state(s) of each step have changed in a sense that each parallel step only has access to a part of the original state.

## 3.2 Distributed State

Therefore the actual concern in distributed machine learning is not how to distribute computation but how to distribute the state in order to achieve optimal performance. As this thesis focuses on machine learning, a state is in general represented by a tensor (e.g. vector or matrix). The most efficient state distribution depends on properties of the state(s) such as size and sparsity of the input data and model, algorithm used for optimization and infrastructure properties such as available memory and computational resources.

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<sup>2</sup>degree of parallelism

### 3.3 Consistency

NOTES: - motivate the argumentation with an example machine learning pipeline (elastic-net linear regression) which includes preprocessing (e.g. standard-scaling), iterative model refinement which could be data/model parallel and depending on the size of the model it could either be replicated or partitioned among machines/workers, also a ML pipeline involves generalization error assessment via test/validation dataset, which should be included into the stopping criterion this should show: - what kind of workflow is generally executed in practice - what structure a ml pipeline has and that there are parts that can be run in parallel and in parallel with relaxed consistency/synchronization - show what kind of information can be inferred from knowledge about the architecture (computational and memory resources, bandwidth, bandwidth quota) and the problem size (size of input data, size of model) - what part of distributing a machine learning algorithm can be handled by the system (distributing data, distributing model and computation, scheduling of work) and what needs to be specified by the developer (parallelism, control flow and required states (?), )

- describe the limitations of current state-of-the-art distributed machine learning systems - low level primitives - not flexible nor expressible api for developing real word machine learning pipelines - inference of certain properties depending on hardware architecture (GPU, CPU, FPGA) and algorithm properties - distributed control flow - focus on optimizing distributed machine learning performance by quick prototyping - only concerned with the important parts for going from an exact single machine to a distributed algorithm implementation by specifying data partitioning, topology and consistency management properties (synchronization requirements/schemes, filter, update/merging strategies) - dealing with algorithm related hyperparameters

## Chapter 4

# Consistency Management

# Chapter 5

## Experiments

## Chapter 6

## Conclusion

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