

AI-Systems Distributed Training

Joseph E. Gonzalez
Co-director of the RISE Lab
jegonzal@cs.berkeley.edu

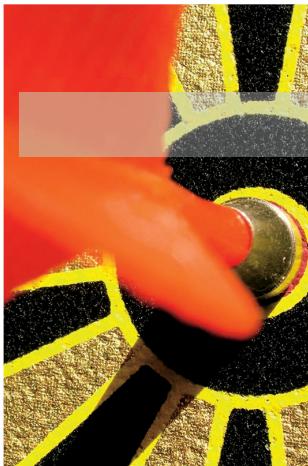
What is? & Why? Distributed Training

- **Distributed Training*** ~ Training across multiple devices
 - Different local and remote memory speeds / network
- Why do we need distributed training?
 - Faster training by leveraging **parallel computation**
 - **Additional memory** (memory bandwidth) for larger model
 - “Need” to store weights + activations
 - Reduce or eliminate **data movement**
 - Privacy → Federated Learning
 - Limited bandwidth to edge devices
 - Need to process all the data?

*Very simplified definition.

On Dataset Size and Learning

- Data is a resource! (e.g., like processors and memory)
 - Is having lots of processors a problem?
- You don't have to use all the data!
 - Though using more data can often help
- More data often* dominates models and algorithms



EXPERT OPINION
Contact Editor: Brian Brannon, bbrannon@computer.org

The Unreasonable Effectiveness of Data

Alon Halevy, Peter Norvig, and Fernando Pereira, Google

*More data also supports more sophisticated models and algorithms.

What are the Metrics of Success?

- **Marketing Team:** Maximize number of GPUs/CPUs used
 - A bad metric ... why?
- **Machine Learning:** Minimize passes through the training data
 - Easy to measure, but not complete ... why?
- **Systems:** minimize time to complete a pass through the training data
 - Easy to measure, but not complete ... why?

Ideal Metric of Success

$$\left(\frac{\text{“Learning”}}{\text{Second}} \right) = \left(\frac{\text{“Learning”}}{\text{Record}} \right) \times \left(\frac{\text{Record}}{\text{Second}} \right)$$

*Convergence
Machine Learning
Property*

*Throughput
System
Property*

Metrics of Success

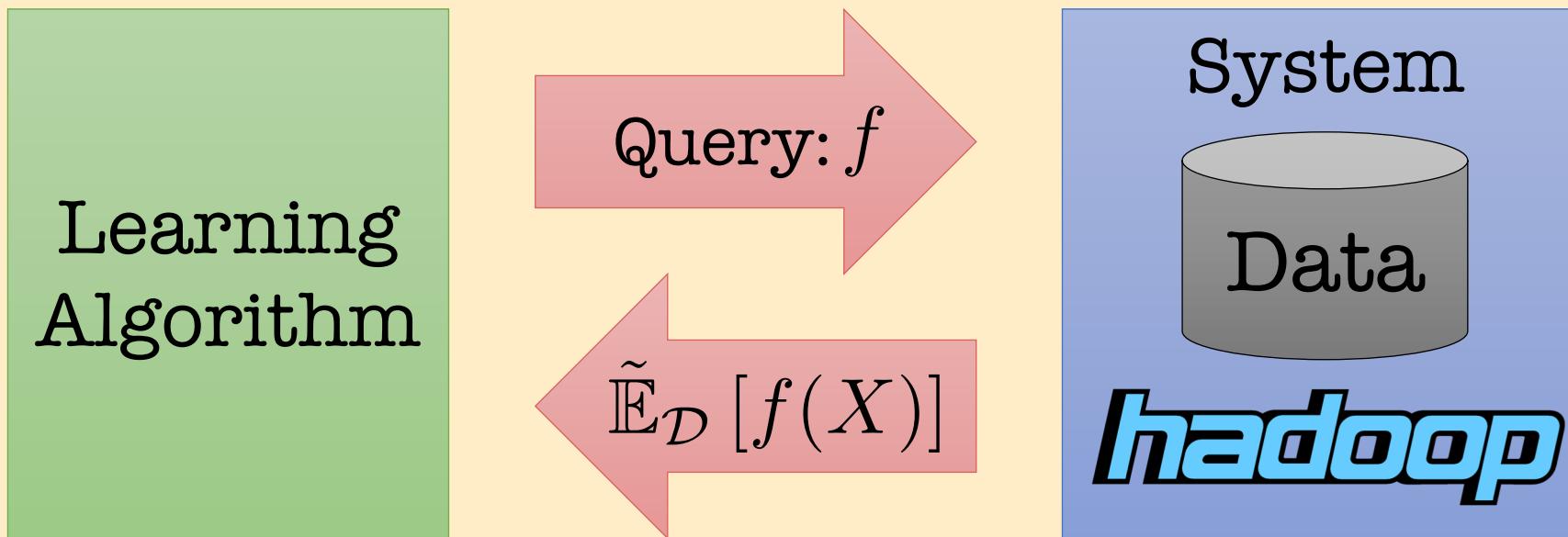
- Minimize training time to “best model”
 - Best model measured in terms of test error
- Other Concerns?
 - **Complexity:** Does the approach introduce additional training complexity (e.g., hyper-parameters)
 - **Stability:** How consistently does the system *train* the model?
 - **Cost:** Will obtaining a faster solution cost more money (power)?

The Early Days....

Map-Reduce for Distributed Training

Learning by Distributed Aggregation

LEARNING FROM STATISTICS (AGGREGATION)*

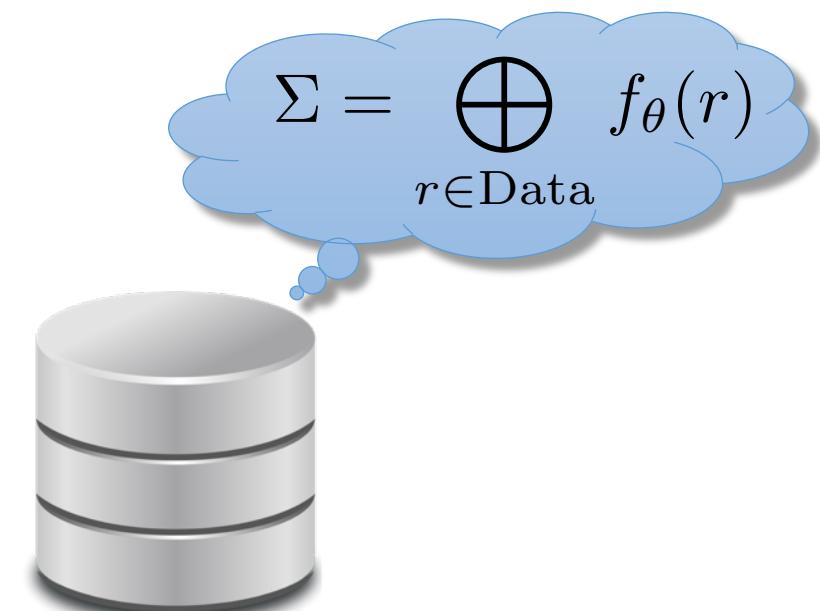
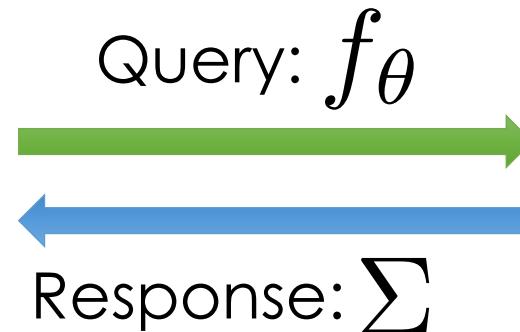
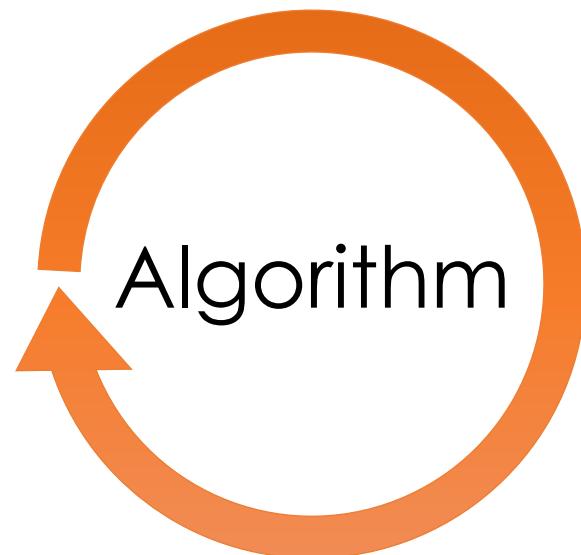


- D. Caragea et al., *A Framework for Learning from Distributed Data Using Sufficient Statistics and Its Application to Learning Decision Trees*. Int. J. Hybrid Intell. Syst. 2004
- Chu et al., *Map-Reduce for Machine Learning on Multicore*. NIPS'06.

Can we compute

$$\hat{\theta} = (X^T X)^{-1} X^T Y$$

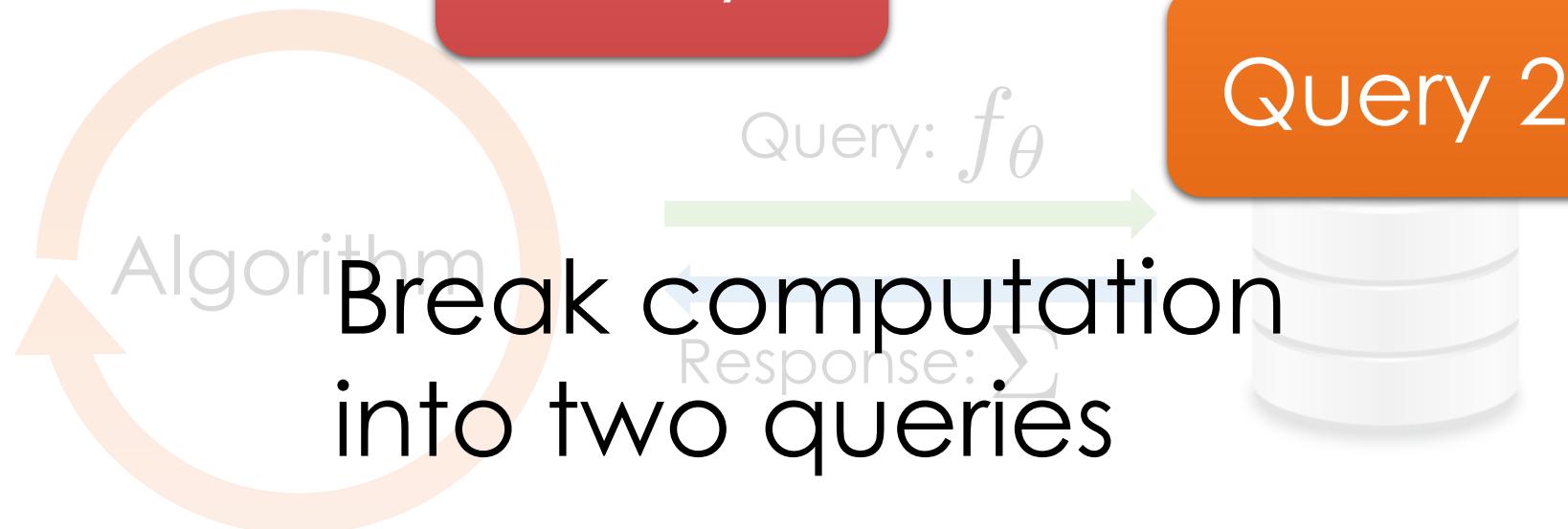
using the statistical query pattern
in map-reduce?



Can we compute

$$\hat{\theta} = \underline{(X^T X)^{-1}} \underline{X^T Y}$$

using the statistical query pattern
in map-reduce?



Cost Analysis

$$\hat{\theta} = (X^T X)^{-1} X^T Y$$

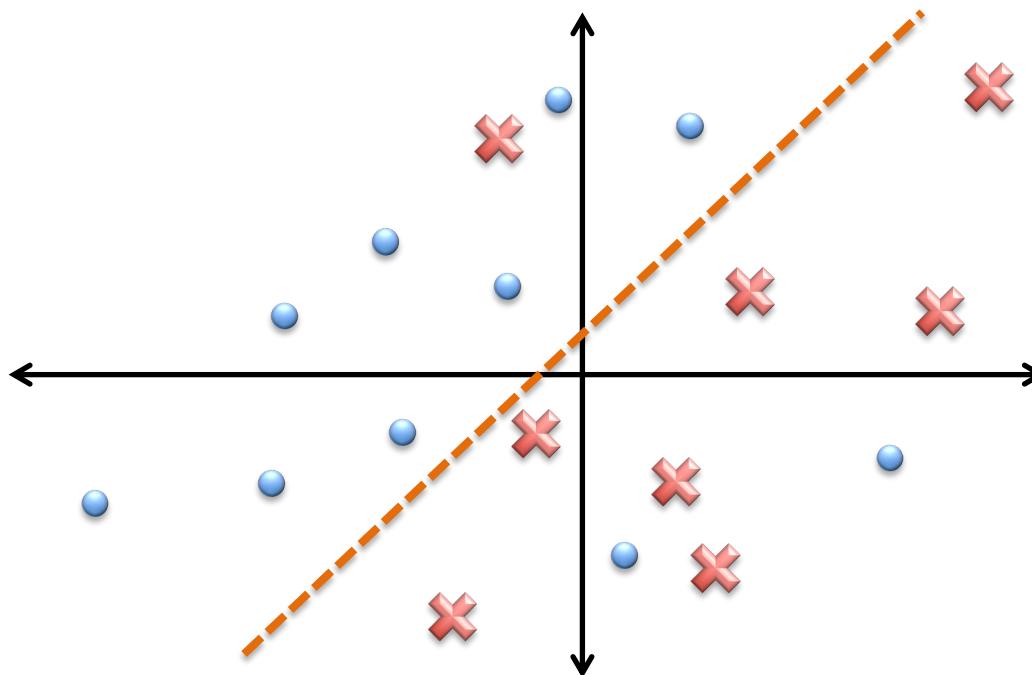
Computation	Cost	
$A = X^T X$	$\Rightarrow O(np^2)$	When $n \gg p$ we want to distribute this computation
$C = X^T Y$	$\Rightarrow O(np)$	

$$B = A^{-1} \Rightarrow O(p^3)$$

$$BC \Rightarrow O(p^2)$$

What about

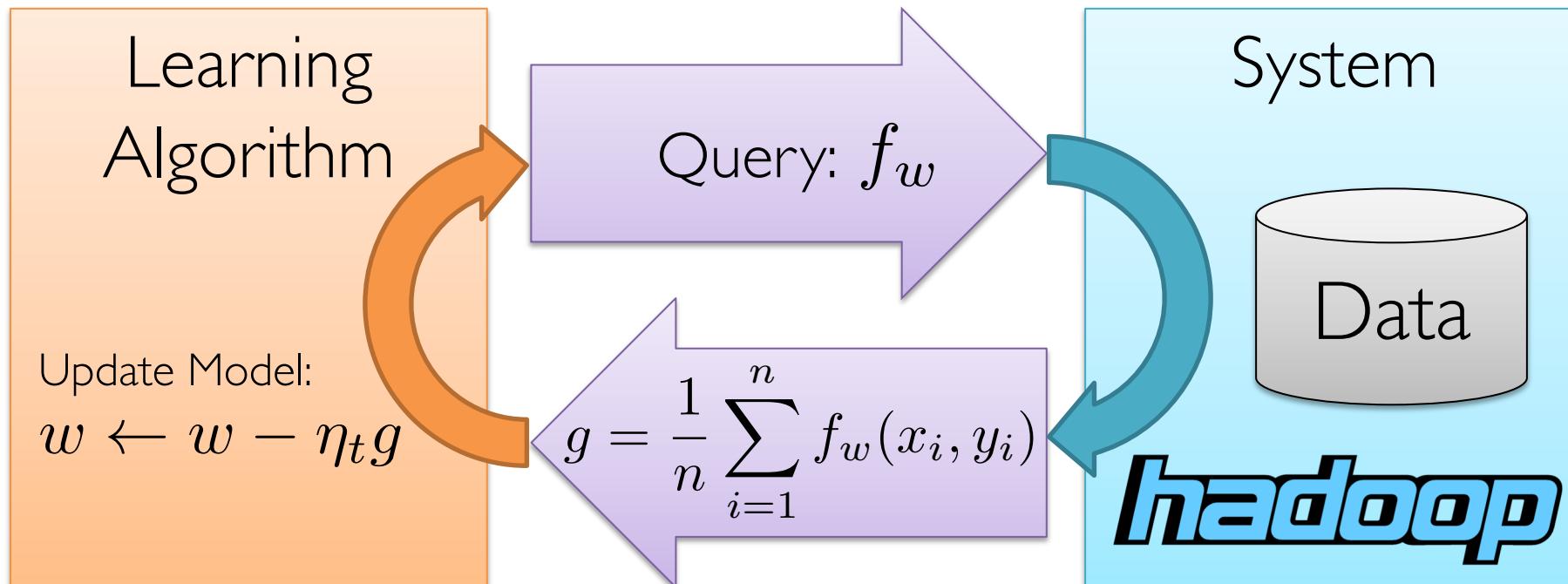
Logistic Regression using Gradient Descent?



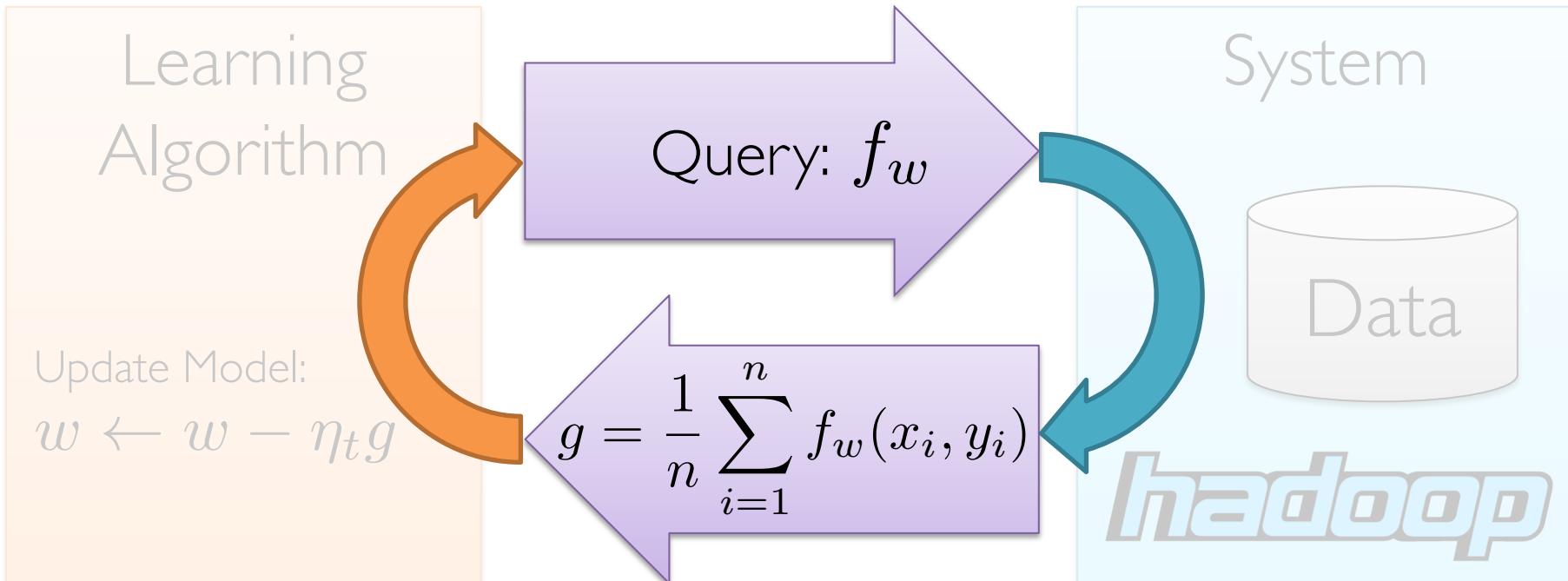
Logistic Regression in Map-Reduce

Gradient descent:

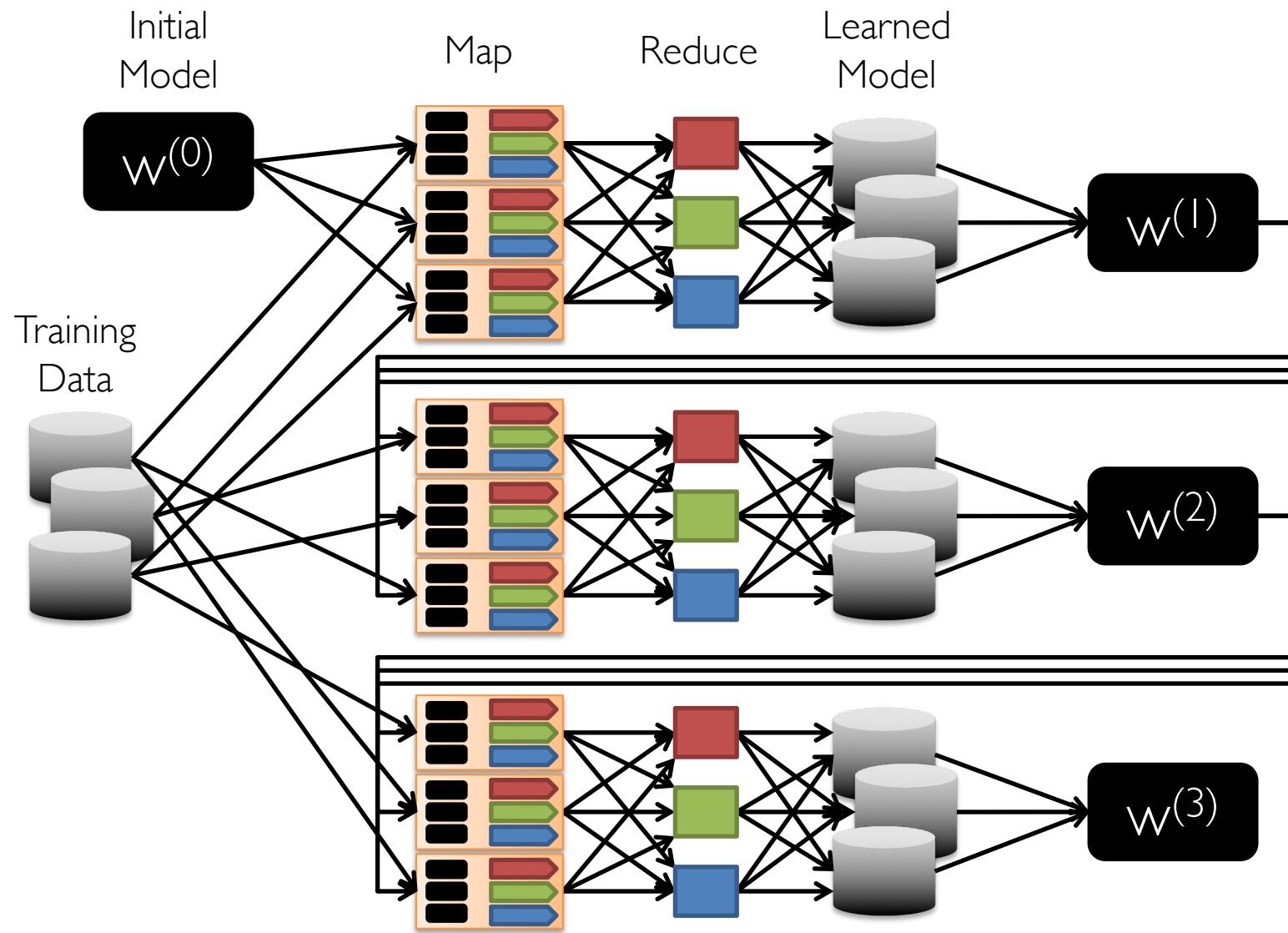
$$f_w(x, y) = \nabla \log L(y, h_w(x))$$



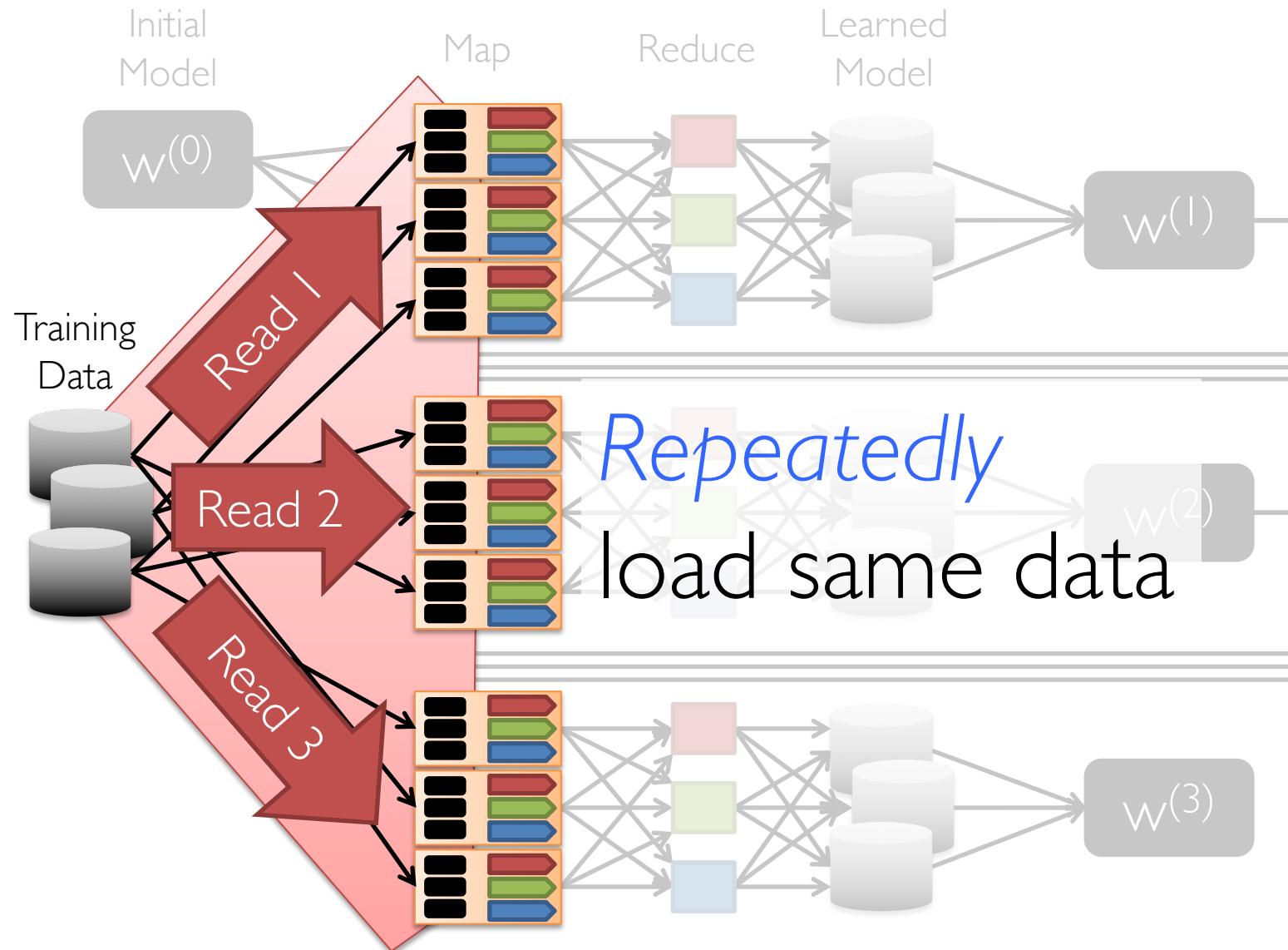
Map-Reduce is not optimized for iteration and multi-stage computation



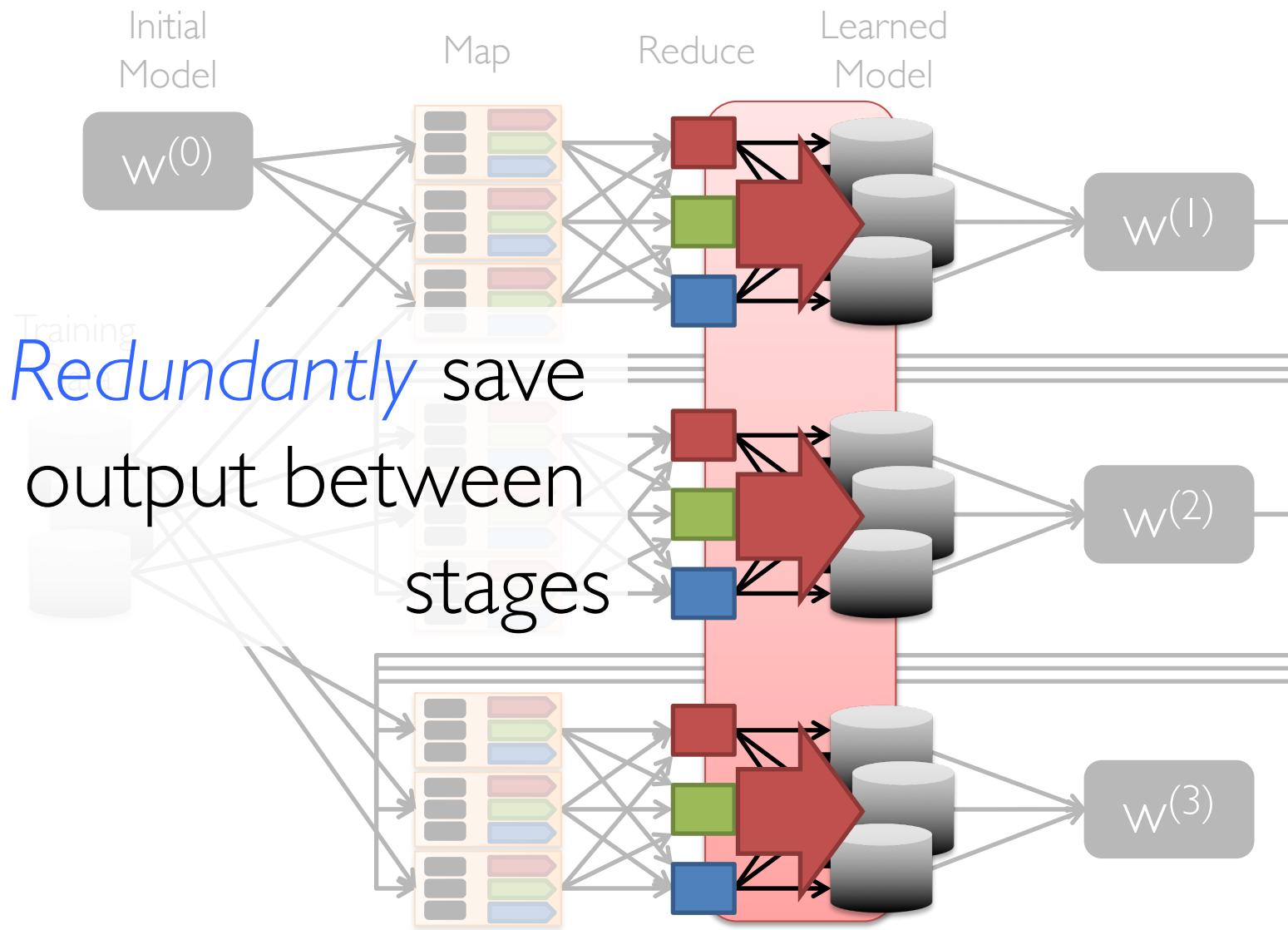
Iteration in Map-Reduce



Cost of Iteration in Map-Reduce



Cost of Iteration in Map-Reduce





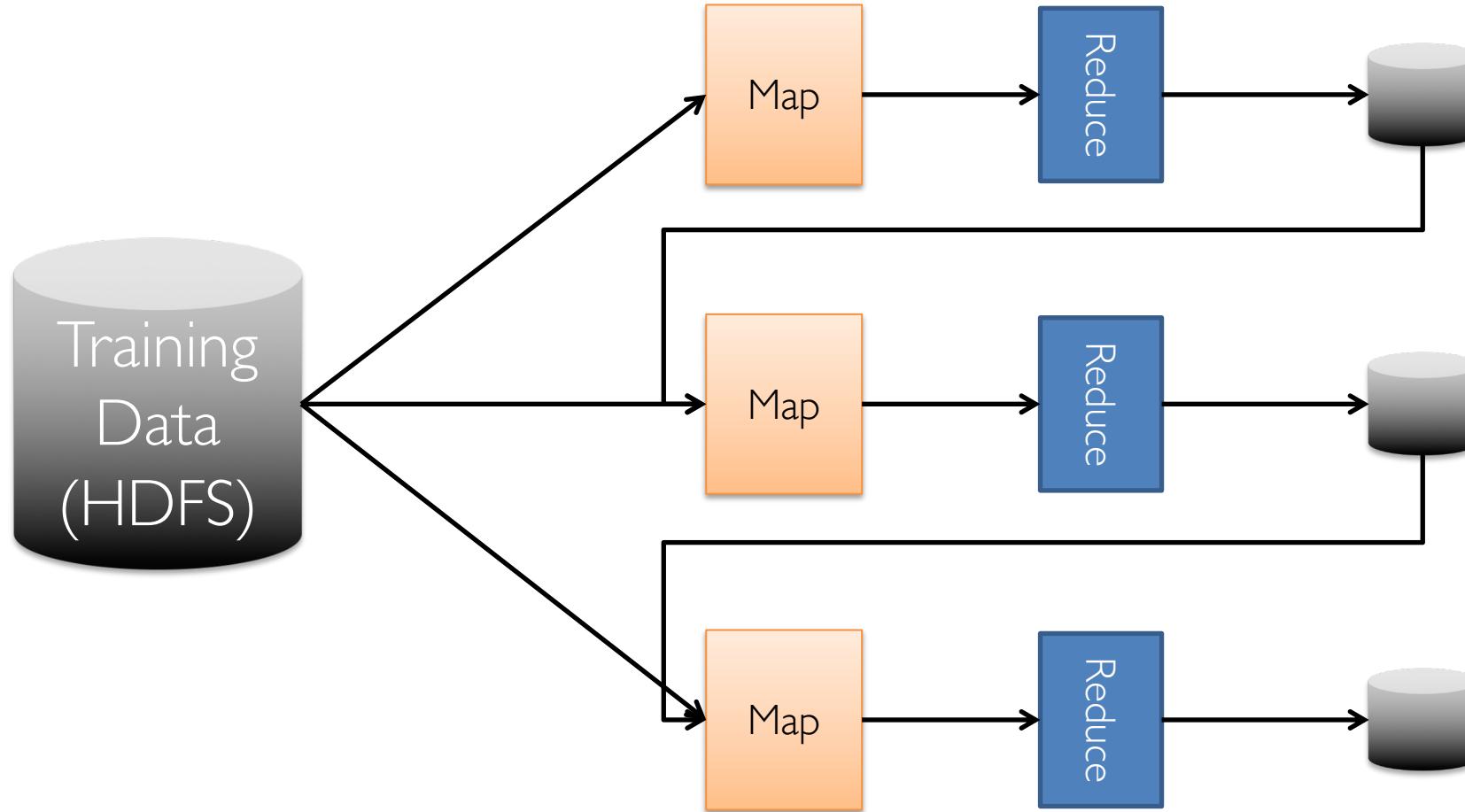
Iteration and Multi-stage computation

In-Memory Dataflow System

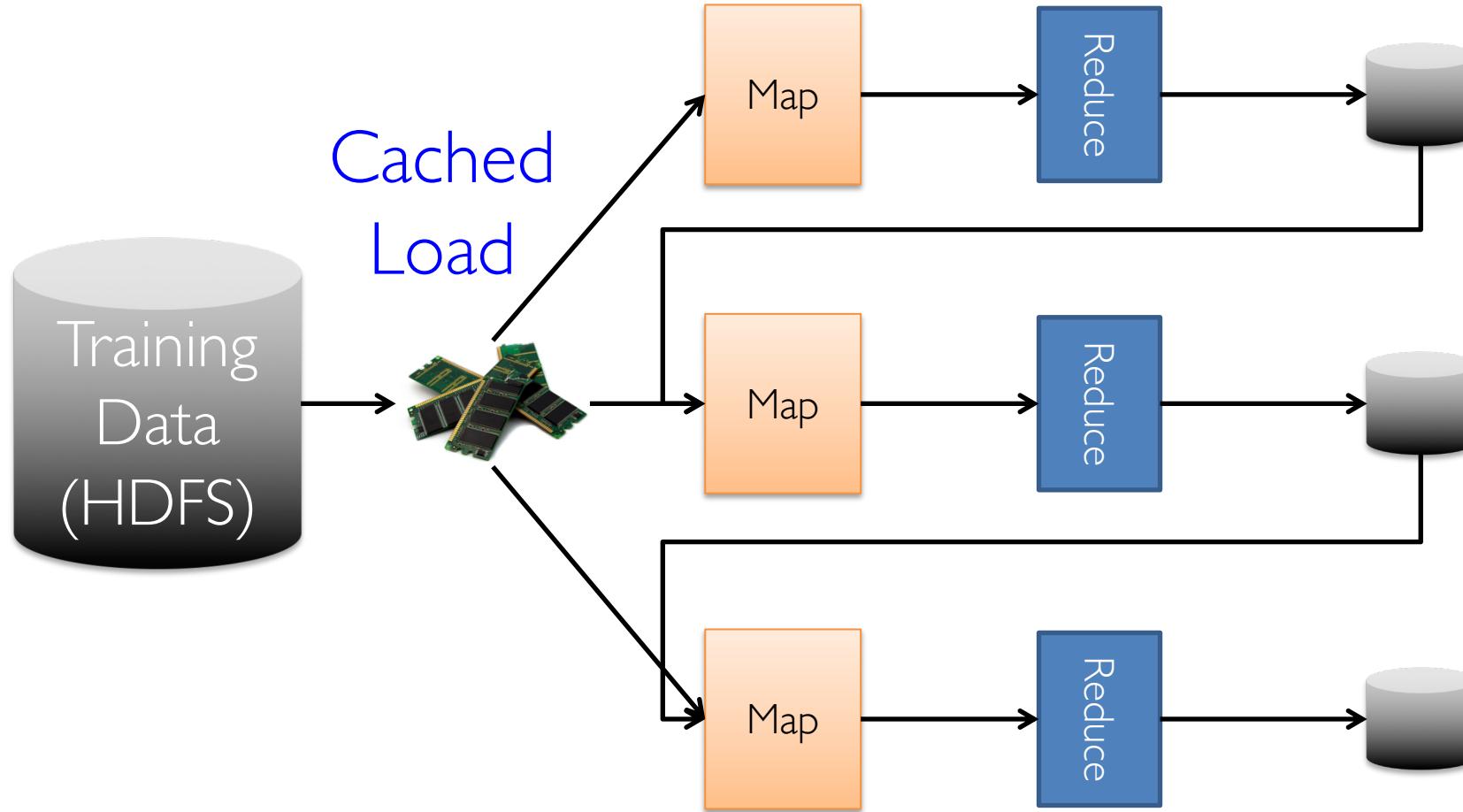
M. Zaharia, M. Chowdhury, M. J. Franklin, S. Shenker, and I. Stoica. *Spark: cluster computing with working sets*. HotCloud'10

M. Zaharia, M. Chowdhury, T. Das, A. Dave, J. Ma, M. McCauley, M.J. Franklin, S. Shenker, I. Stoica.
Resilient Distributed Datasets: A Fault-Tolerant Abstraction for In-Memory Cluster Computing, NSDI 2012

Dataflow View

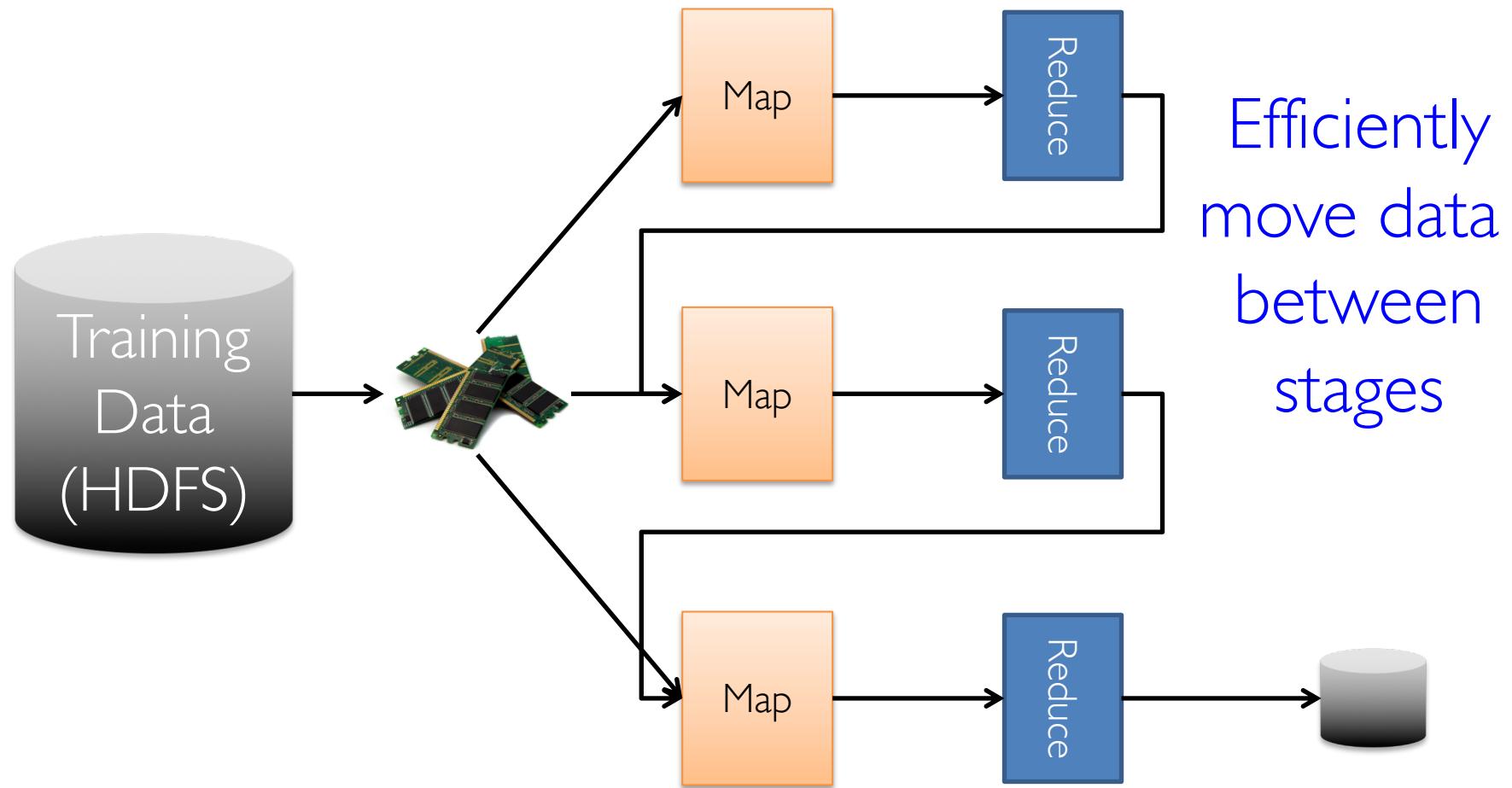


Memory Opt. Dataflow



10-100× faster than network and disk

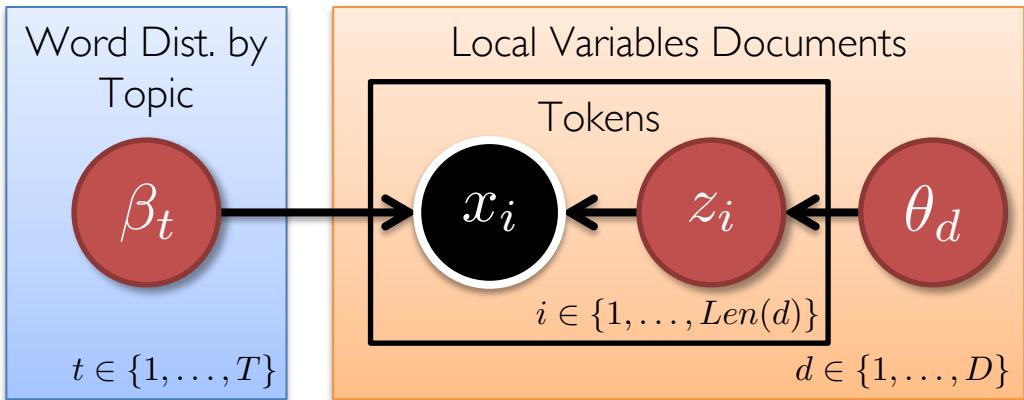
Memory Opt. Dataflow View



Efficiently
move data
between
stages

Statistical Inference in Large Latent Variable Models

- Large topic models associated variables with each word and document



- Not a good fit for BSP model

Scalable Inference in Latent Variable Models

Amr Ahmed, Mohamed Aly, Joseph Gonzalez,^{*} Shravan Narayananmurthy, Alexander Smola
Yahoo Research, Santa Clara, CA, USA
{amahmed, aly, jegonzal, shravanm, smola}@yahoo-inc.com

ABSTRACT

Latent variable techniques are pivotal in tasks ranging from predicting user click patterns and targeting ads to organizing the news and managing user generated content. Latent variable techniques like topic modeling, clustering, and subspace estimation provide substantial insight into the latent structure of complex data with little or no external guidance making them ideal for reasoning about large-scale, rapidly evolving datasets. Unfortunately, due to the data dependencies and global state introduced by latent variables and the iterative nature of latent variable inference, latent-variable techniques are often prohibitively expensive to apply to large-scale streaming datasets.

In this paper we present a scalable parallel framework for efficient inference in latent variable models over streaming, web-scale data. Our framework addresses three key challenges: 1) *synchronizing the global state* which includes global latent variables (e.g., cluster centers and dictionaries); 2) *efficiently storing and retrieving the large local state* which includes the data-points and their corresponding latent variables (e.g., cluster membership); and 3) *sequentially incorporating streaming data* (e.g., the news). We address these challenges by introducing a novel data-based aggregation system with a highly efficient communication protocol and schedule-aware out-of-core storage; and 3) approximate forward sampling to rapidly incorporate new data. We demonstrate state-of-the-art performance of our framework by easily tackling datasets two orders of magnitude larger than those addressed by the current state-of-the-art. Furthermore, we provide an optimized and easily customizable open-source implementation of the framework¹.

Categories and Subject Descriptors
G.3 [Probability And Statistics]: Statistical Computing

^{*}Visiting on internship from CMU, Department of Machine Learning, Pittsburgh PA; jegonzal@cs.cmu.edu
¹ Available at https://github.com/shravanm/Yahoo_LDA

Permission to make digital or hard copies of all or part of this work for personal use or classroom use is granted for free provided that copies are not made or distributed for profit or commercial advantage and that copies bear this notice and the full citation on the first page. To copy otherwise, to republish, to post on servers or to redistribute in lists, requires prior specific permission and/or a fee.
Copyright 2012 ACM 978-1-4503-1073-5/12/02 \$10.00.

123

General Terms

Algorithms, Experimentation, Performance

Keywords

Inference, Graphical Models, Large-scale Systems, Latent Models.

I. INTRODUCTION

In many cases, we are interested in reasoning about the underlying latent causes that give rise to the data we observe. For instance, when dealing with users we may want to elicit the underlying intents and interests that govern their activity and friendship patterns. Alternatively, we might want to discover the underlying topics of discussions on various pages across the web. More generally we may want to assign meaning to linked and interacting objects such as webpages, named entities, users, and their behavior.

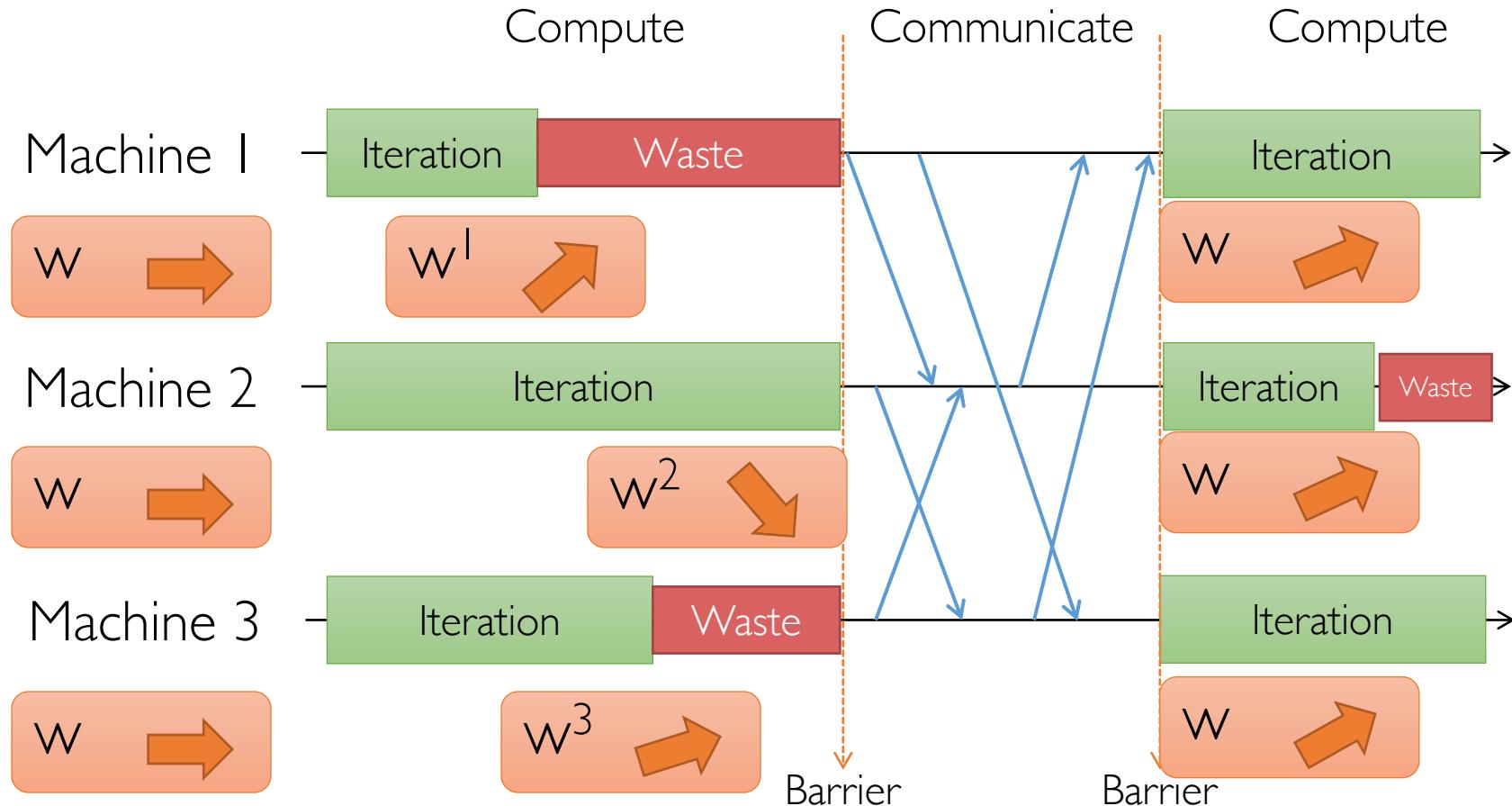
Latent variable models have become an indispensable tool for reasoning about the latent causes that give rise to data in tasks ranging from text modeling [18, 3, 6] to bioinformatics [8, 21]. The popularity of latent variable models stems from their ability to easily encode rich structured priors and then infer latent properties of the data without requiring access to costly labels or editorial feedback.

Latent variable models are constructed by introducing unobserved (latent) variables which help explain the observed data and then coupling these latent variables (often by introducing additional latent variables) to capture the underlying problem structure. For example, in the mixture of Gaussians model there are two sets of latent variables. The first denotes the cluster membership of each data point while the second describes the shape and position of the Gaussian clusters and introduces dependencies between the data-points.

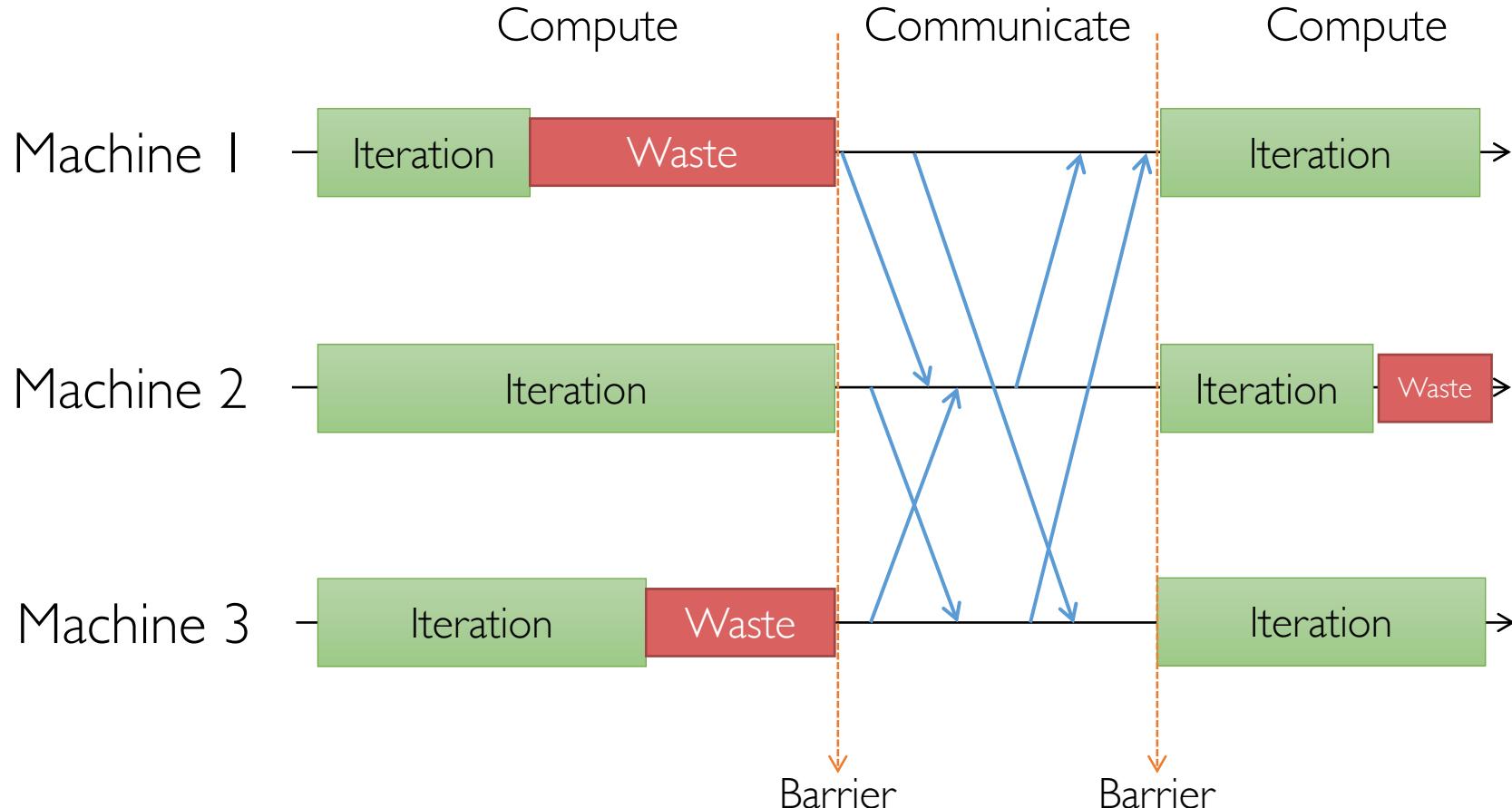
Latent variable inference is the process of estimating the most likely assignment (or posterior distribution) of all the latent variables. In the context of the mixture of Gaussians model, latent variable inference is the process of estimating the latent membership of each data-point as well as the center and shape of each cluster. Inference in latent variable models is typically computationally expensive, often requiring the solution to hard combinatorial search problems. As a consequence, approximate inference algorithms are typically employed. Unfortunately, even approximate inference algorithms can be costly, requiring iterative transformations of the latent variable assignments (i.e., transforming large amounts of program state).

In most web-scale settings, data is not collected once and then processed offline; instead, data arrives continuously and

Bulk Synchronous Parallel (BSP) Execution

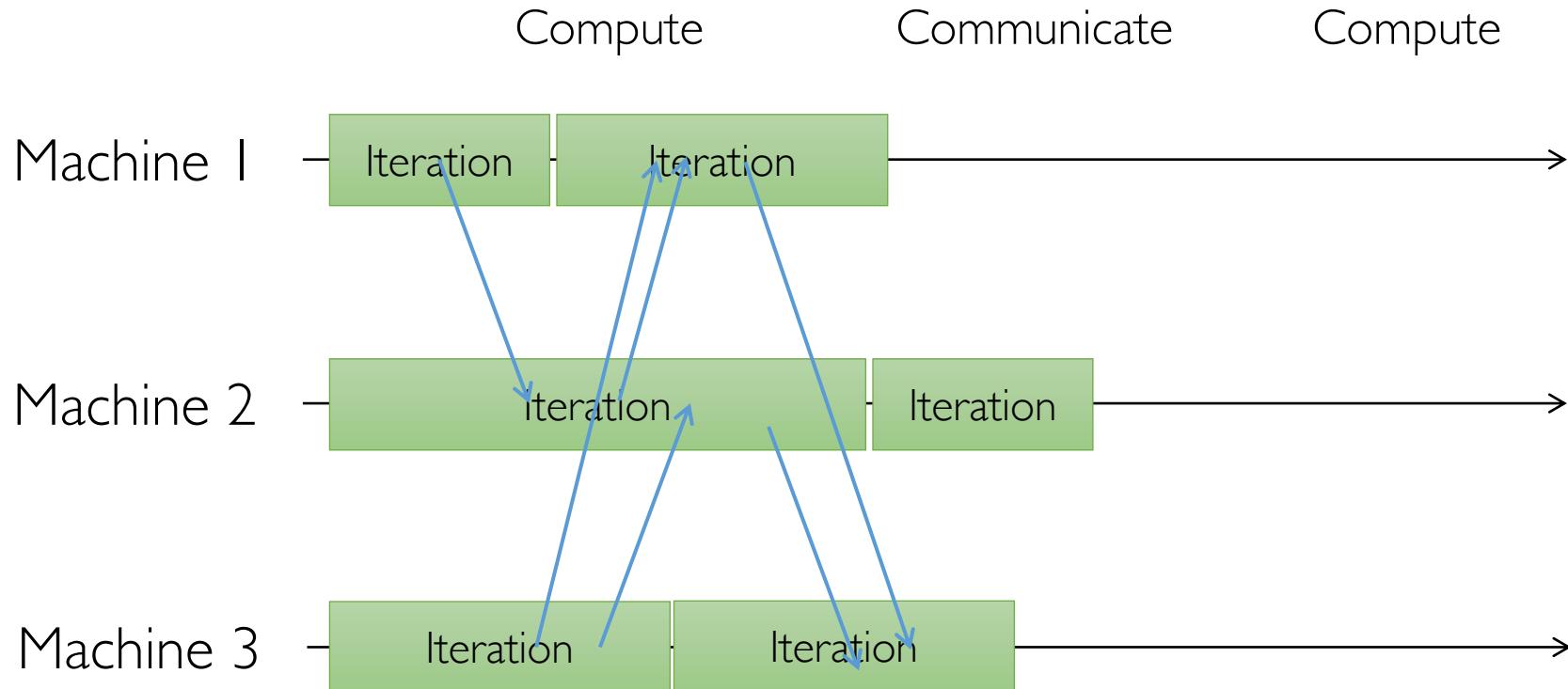


Asynchronous Execution



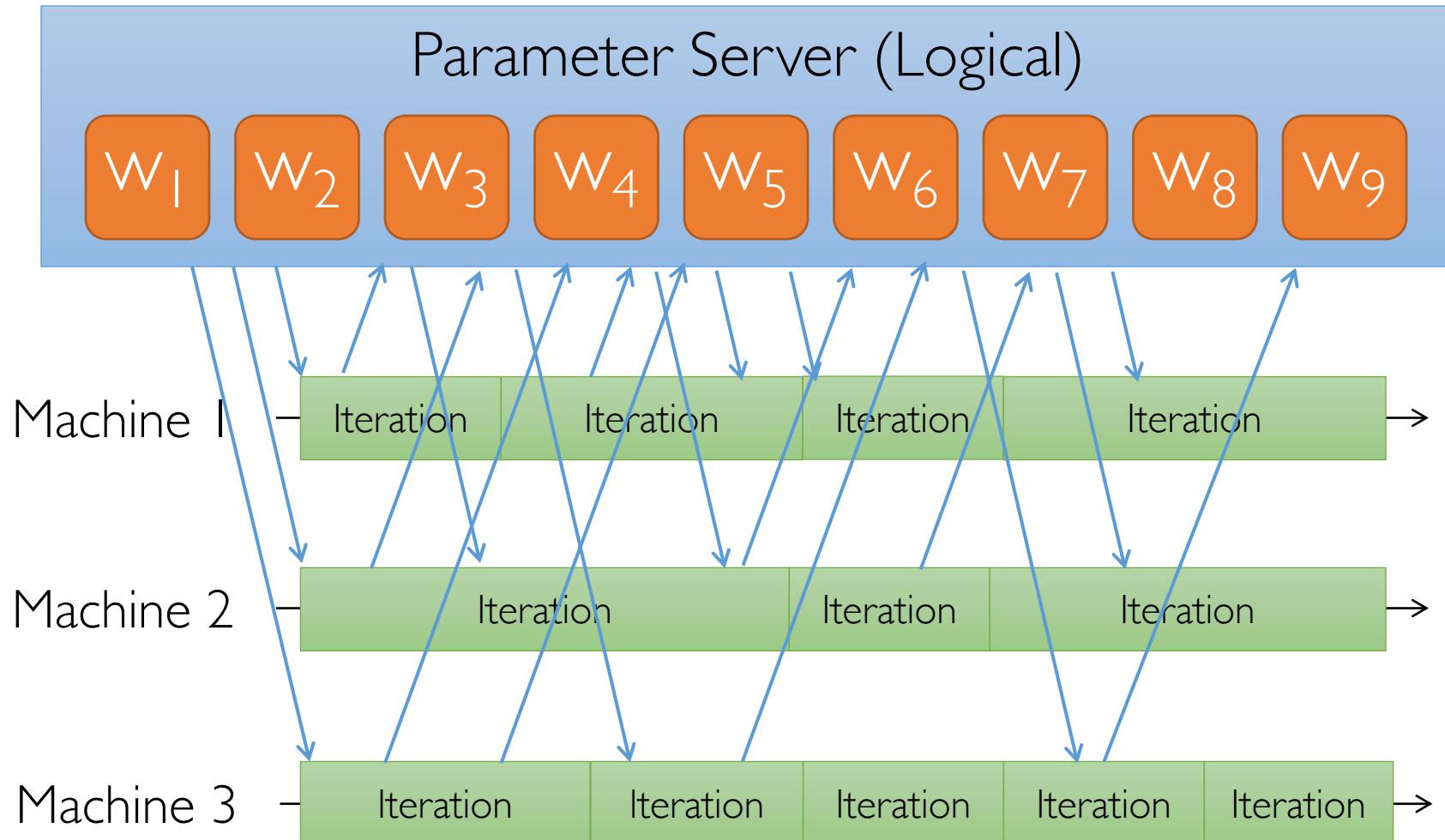
Enable more frequent coordination on parameter values

Asynchronous Execution



Enable more frequent coordination on parameter values

Asynchronous Execution



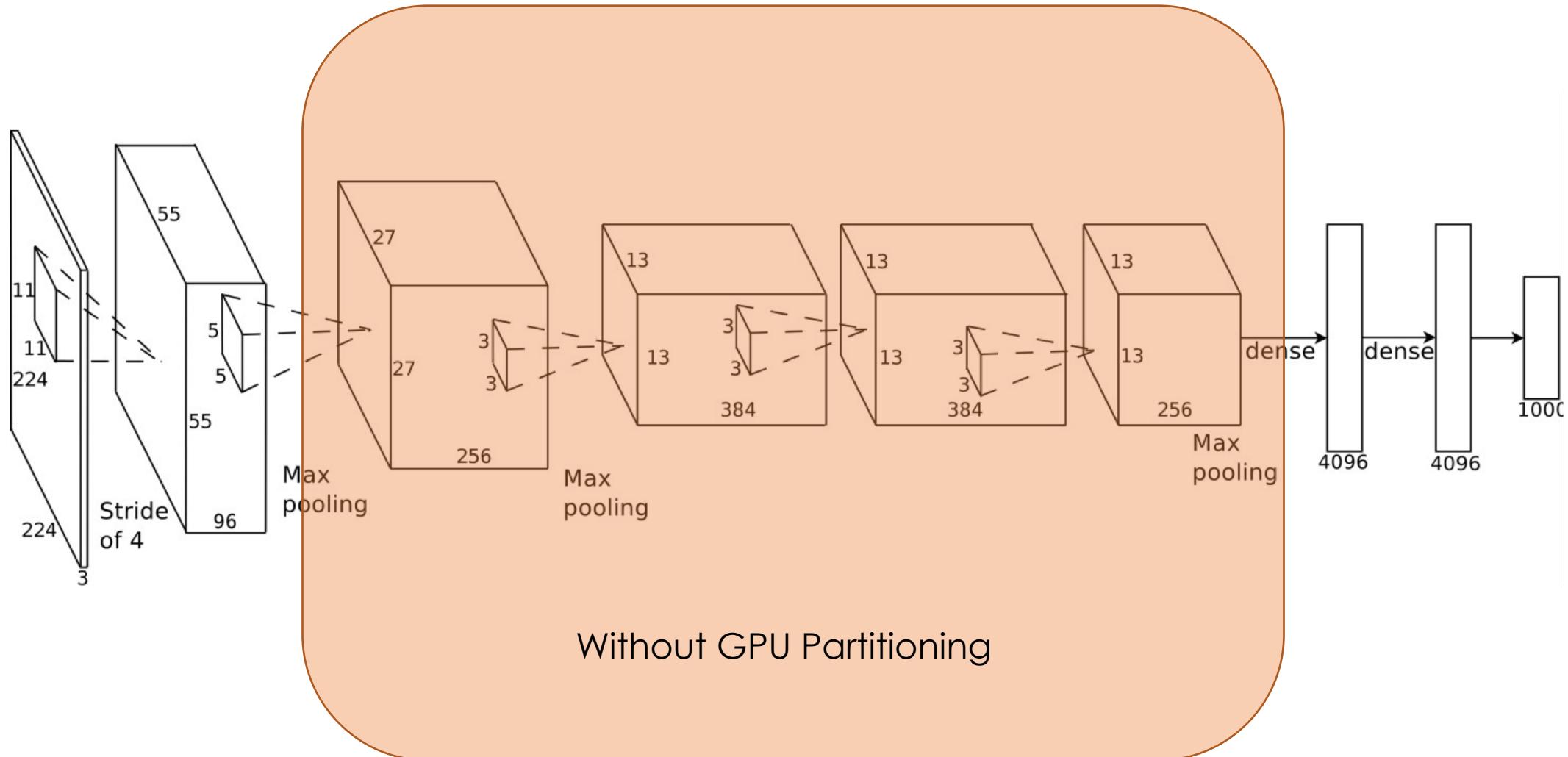
AlexNet

ImageNet Classification with Deep Convolutional Neural Networks

Alex Krizhevsky, Ilya Sutskever, Geoffrey E. Hinton

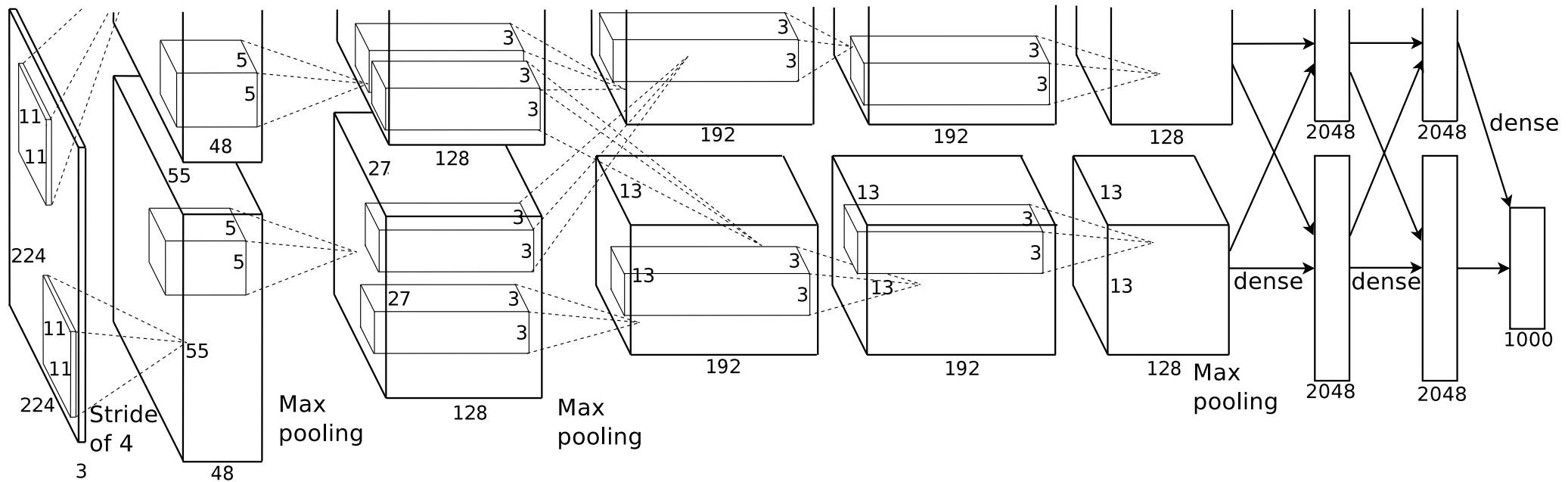
TL;DR; This paper describe the deep convolutional architecture, training techniques, and system innovations that resulted in the winning entry for the ILSVRC-2012 Benchmark. This model substantially outperformed the next best model that year.

The AlexNet* Architecture



The Actual AlexNet* Architecture

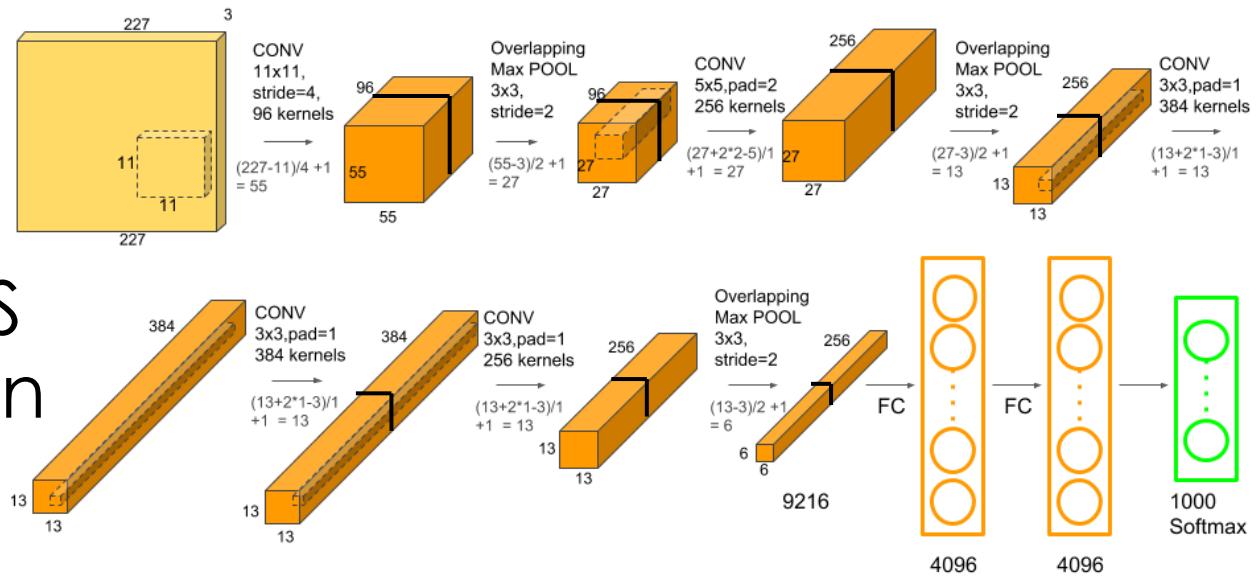
from the paper



*Posthumously Named

Training on Multiple GPUs

- Limited by GPU **memory** using Nvidia GTX 580 (3GB RAM)
 - 60M Parameters ~ **240 MB**
 - Need to cache activation maps for backpropagation
 - Batch size = 128
 - $128 * (227*227*3 + 55*55*96*2 + 96*27*27*2 + 256*27*27*2 + 256*13*13*2 + 13*13*384*2 + 256*13*13 + 6*6*256 + 4096 + 4096 + 1000) * 4 \text{ Bytes} \sim 782\text{MB Activations}$
 - That is assuming no overhead and single precision values
- Tuned splitting across GPUS to balance communication and computation



Interesting Consequence of Partitioned Training

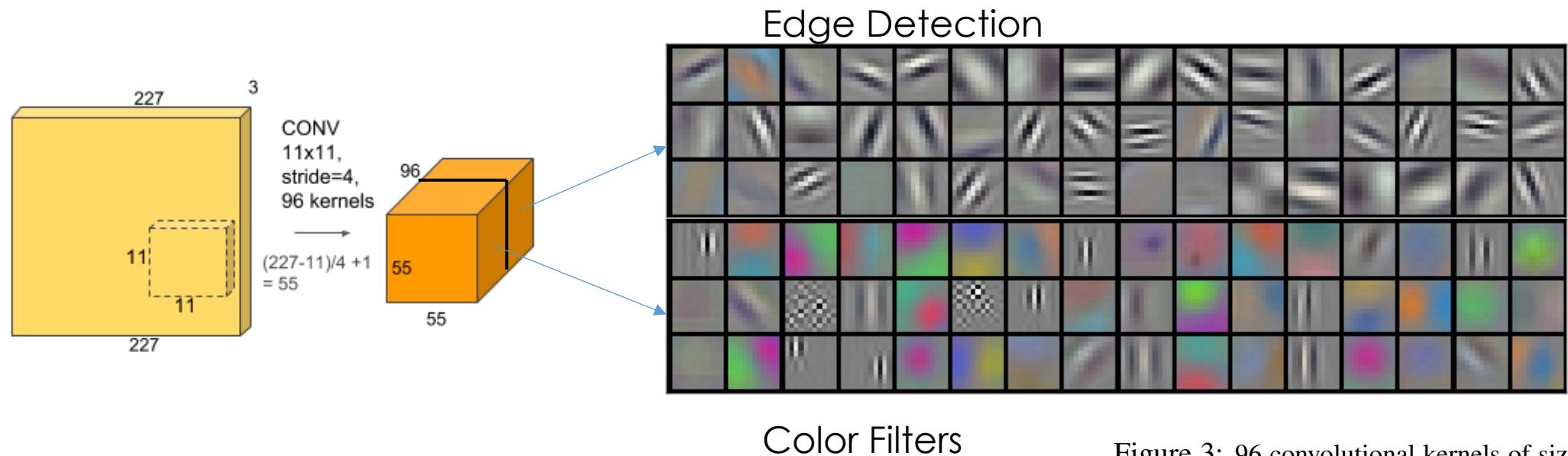
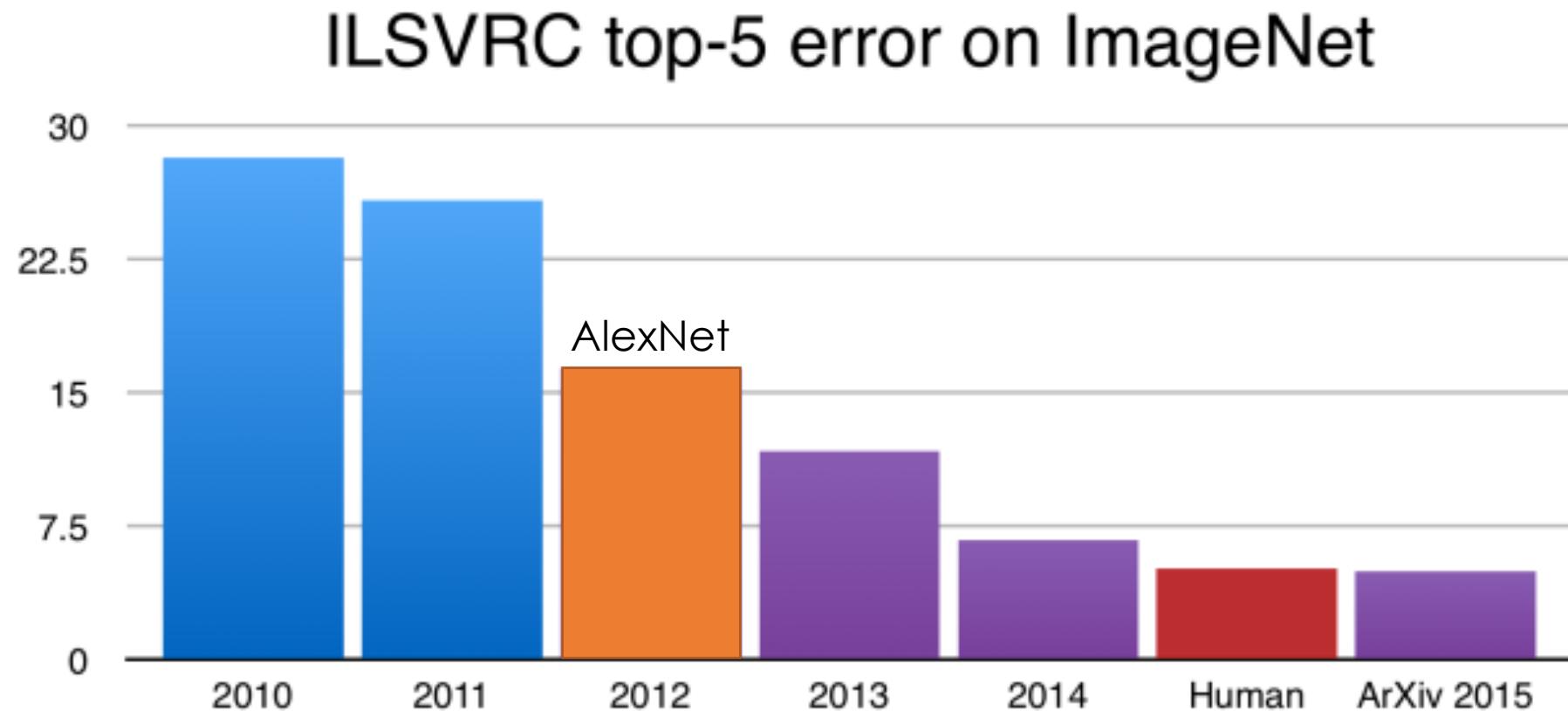


Figure 3: 96 convolutional kernels of size $11 \times 11 \times 3$ learned by the first convolutional layer on the $224 \times 224 \times 3$ input images. The top 48 kernels were learned on GPU 1 while the bottom 48 kernels were learned on GPU 2. See Section 6.1 for details.

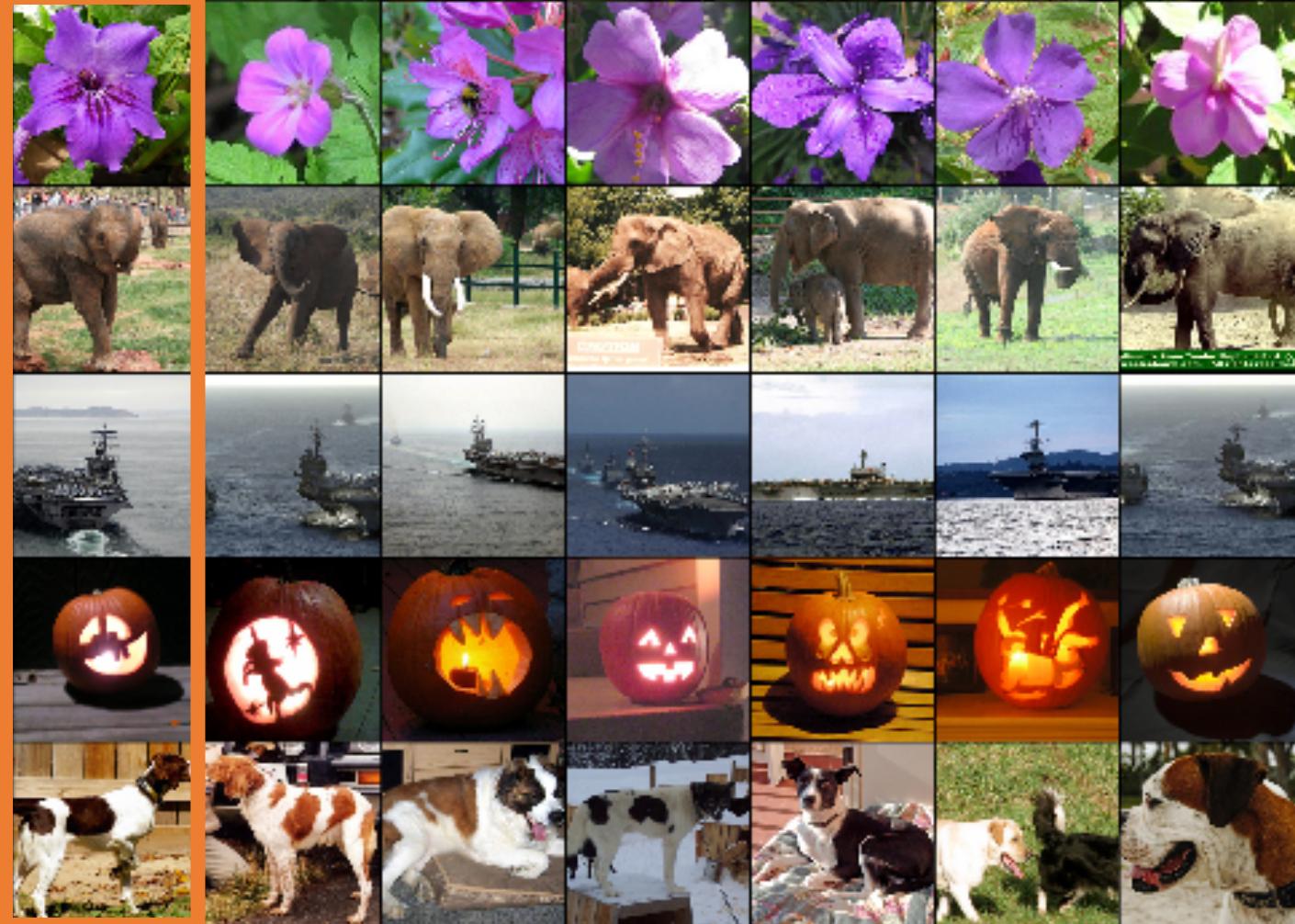
Put into historical context



Good Embeddings ...

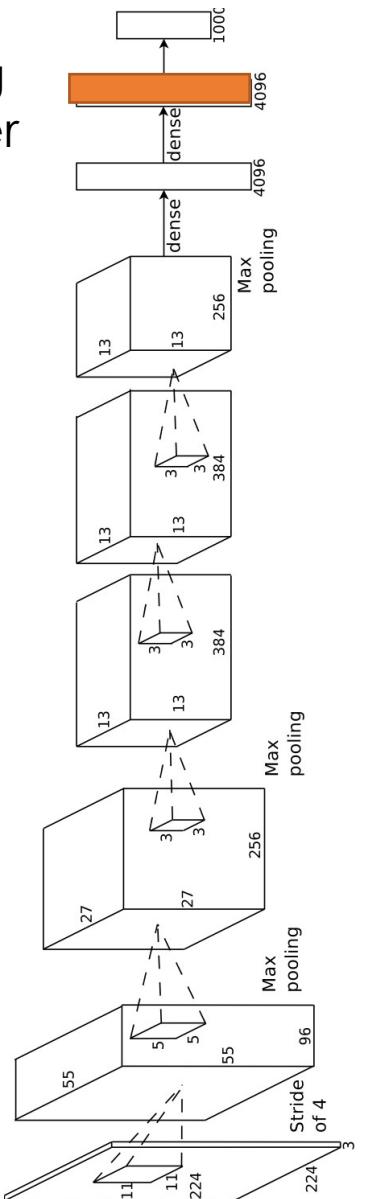
This will later be the foundation of **many** papers

Query



Embedding Layer

Images with
largest dot
product
with query



DistBelief

Large Scale Distributed Deep Networks

Described the system for the 2012 ICML Paper

Building High-level Features Using Large Scale Unsupervised Learning

Quoc V. Le
Marc'Aurelio Ranzato
Rajat Monga
Matthieu Devin
Kai Chen
Greg S. Corrado
Jeff Dean
Andrew Y. Ng

Abstract

We consider the problem of learning high-level, class-specific features from unlabeled data, without any labeled data. For example, it is possible to learn a face detector from a large collection of unlabeled images using unsupervised learning. To answer this, we train a deep neural network with a sparse autoencoder architecture. It consists of two layers of connected sparse autoencoders, with local contrast normalization. The first layer takes a dataset of images (the ImageNet dataset has 15 million connections, the dataset has 10 million



but current experimental evidence suggests the possibility that some neurons in the temporal cortex are

NIPS 2012 (Same Year as AlexNet)

Large Scale Distributed Deep Networks

Jeffrey Dean, Greg S. Corrado, Rajat Monga, Kai Chen,
Matthieu Devin, Quoc V. Le, Mark Z. Mao, Marc'Aurelio Ranzato,
Andrew Senior, Paul Tucker, Ke Yang, Andrew Y. Ng
{jeff, gcorrado}@google.com
Google Inc., Mountain View, CA

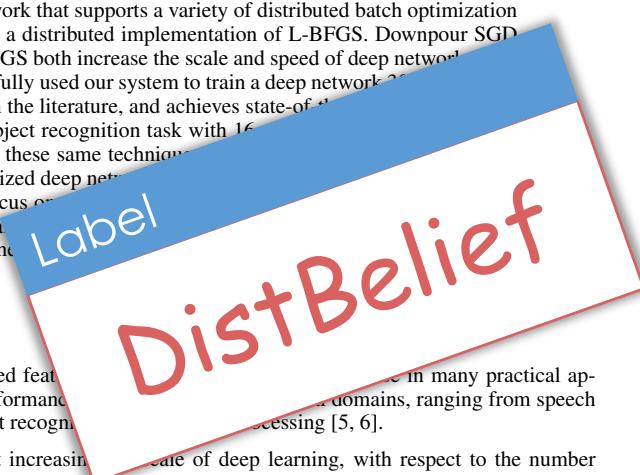
Abstract

Recent work in unsupervised feature learning and deep learning has shown that being able to train large models can dramatically improve performance. In this paper, we consider the problem of training a deep network with billions of parameters using tens of thousands of CPU cores. We have developed a software framework called *DistBelief* that can utilize computing clusters with thousands of machines to train large models. Within this framework, we have developed two algorithms for large-scale distributed training: (i) Downpour SGD, an asynchronous stochastic gradient descent procedure supporting a large number of model replicas, and (ii) Sandblaster, a framework that supports a variety of distributed batch optimization procedures, including a distributed implementation of L-BFGS. Downpour SGD and Sandblaster L-BFGS both increase the scale and speed of deep network training. We have successfully used our system to train a deep network 20 times larger than previously reported in the literature, and achieves state-of-the-art performance on ImageNet, a visual object recognition task with 16 million images and 1000 categories. We show that these same techniques can be applied to training a much smaller, but more complex, model of a more modestly-sized deep network. Although we focus on training large neural networks, the techniques we present are applicable to training large neural networks using gradient-based machine learning.

1 Introduction

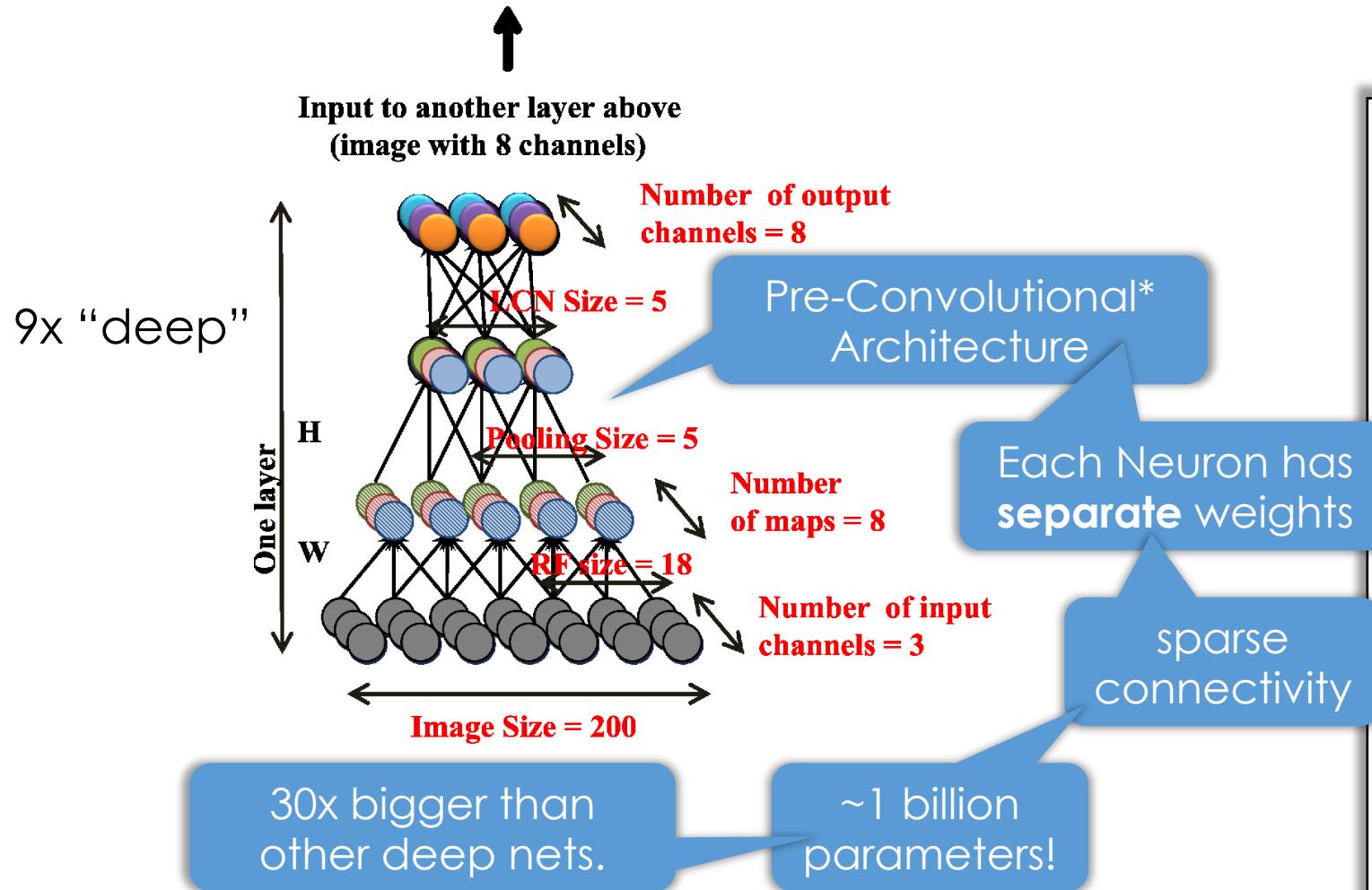
Deep learning and unsupervised feature learning have had significant success in many practical applications. State-of-the-art performance has been achieved in domains, ranging from speech recognition [1, 2], visual object recognition [3, 4] to natural language processing [5, 6].

It has also been observed that increasing the scale of deep learning, with respect to the number of training examples, the number of model parameters, or both, can drastically improve ultimate classification accuracy [3, 4, 7]. These results have led to a surge of interest in scaling up the training and inference algorithms used for these models [8] and in improving applicable optimization procedures [7, 9]. The use of GPUs [1, 2, 3, 8] is a significant advance in recent years that makes the training of such large models tractable. A large limitation of the GPU approach is



Building High-Level Features Using Large Scale Unsupervised Learning

ICML 2012



*This pre-dates AlexNet but is two decades after LeNet.

Building High-level Features
Using Large Scale Unsupervised Learning

Quoc V. Le
Marc'Aurelio Ranzato
Rajat Monga
Mathieu Devin
Kai Chen
Greg S. Corrado
Jeff Dean
Andrew Y. Ng

QUOCLE@CS.STANFORD.EDU
RANZATO@GOOGLE.COM
RAJATMONGA@GOOGLE.COM
MDEVIN@GOOGLE.COM
KAICHEN@GOOGLE.COM
GCORRADO@GOOGLE.COM
JEFF@GOOGLE.COM
ANG@CS.STANFORD.EDU

Abstract

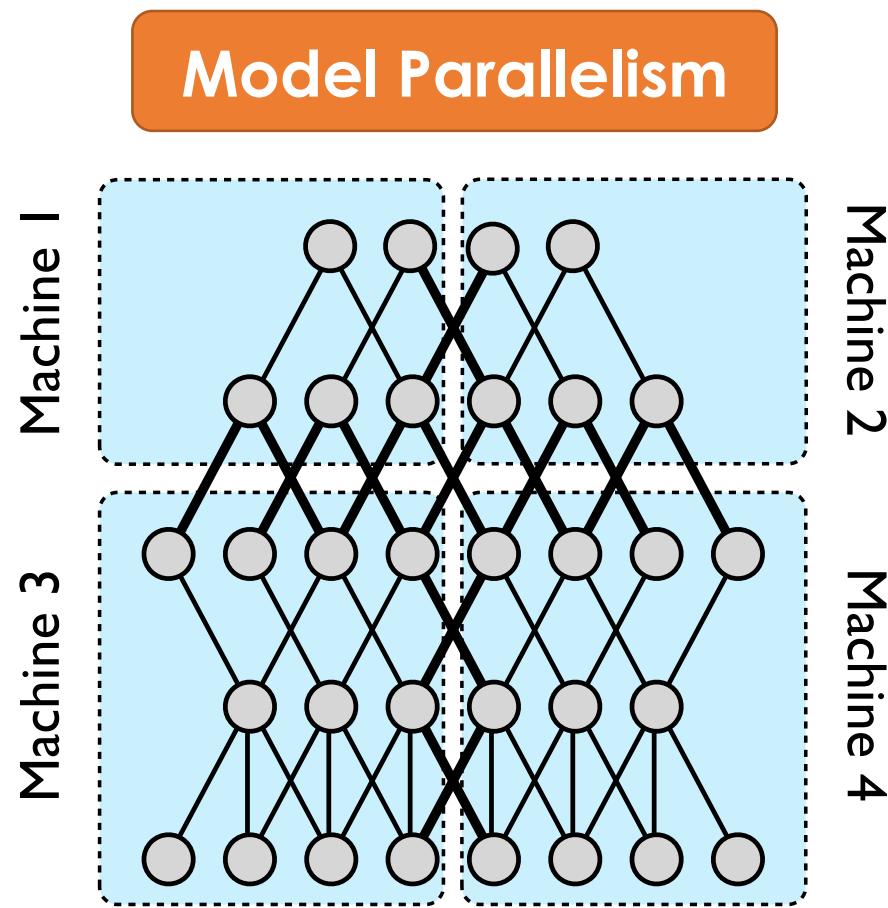
We consider the problem of building high-level, class-specific feature detectors from *unlabeled* images. For instance, we would like to understand if it is possible to build a face detector from only unlabeled images. This approach is inspired by the neuroscientific conjecture that there exist highly class-specific neurons in the human brain, generally and informally known as “grandmother neurons.” The extent of class-specificity of neurons in the brain is an area of active investigation, but current experimental evidence suggests the possibility that some neurons in the temporal cortex are highly selective for object categories such as faces or hands (Desimone et al., 1984), and perhaps even specific people (Quiroga et al., 2005).

1. Introduction

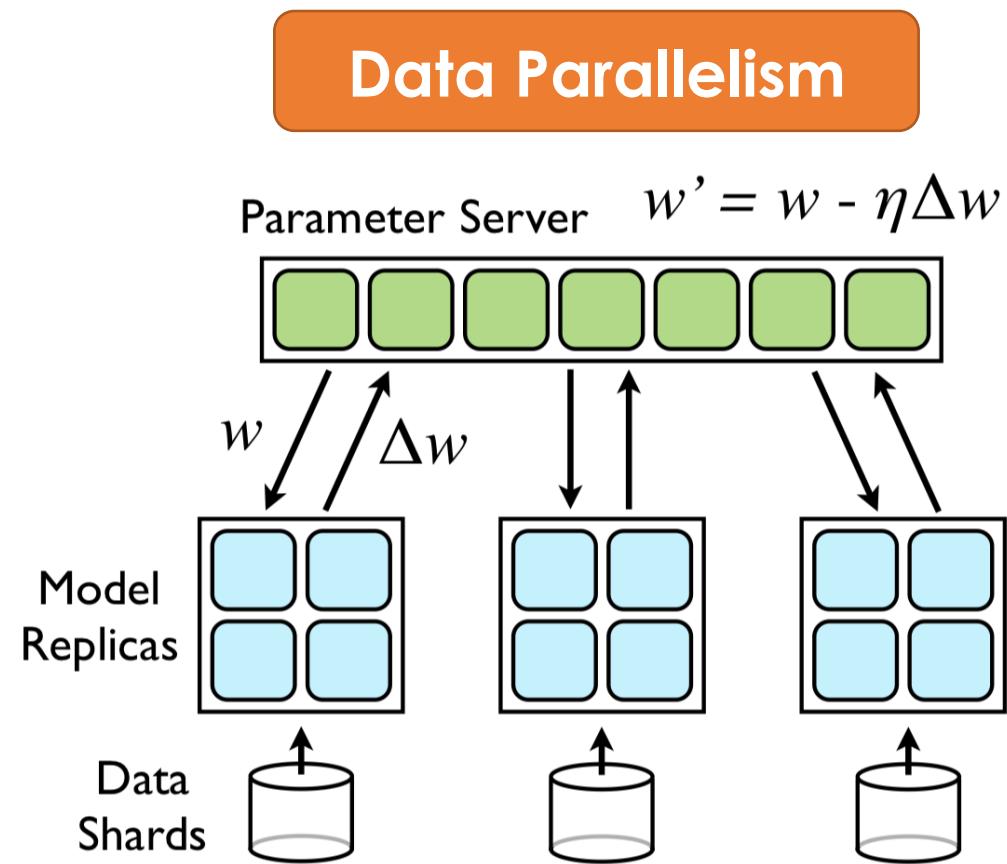
Contemporary computer vision methodology typically emphasizes the role of *labeled* data to obtain these class-specific feature detectors. For example, to build a face detector, one needs a large collection of images labeled as containing faces, often with a bounding box around the face. The need for large labeled sets poses a significant challenge for problems where labeled data are rare. Although approaches that make use of inexpensive unlabeled data are often preferred, they have not been shown to work well for building high-level features.

This work investigates the feasibility of building high-level features from only *unlabeled* data. A positive result in this setting will indicate a significant

Combine Model and Data Parallelism



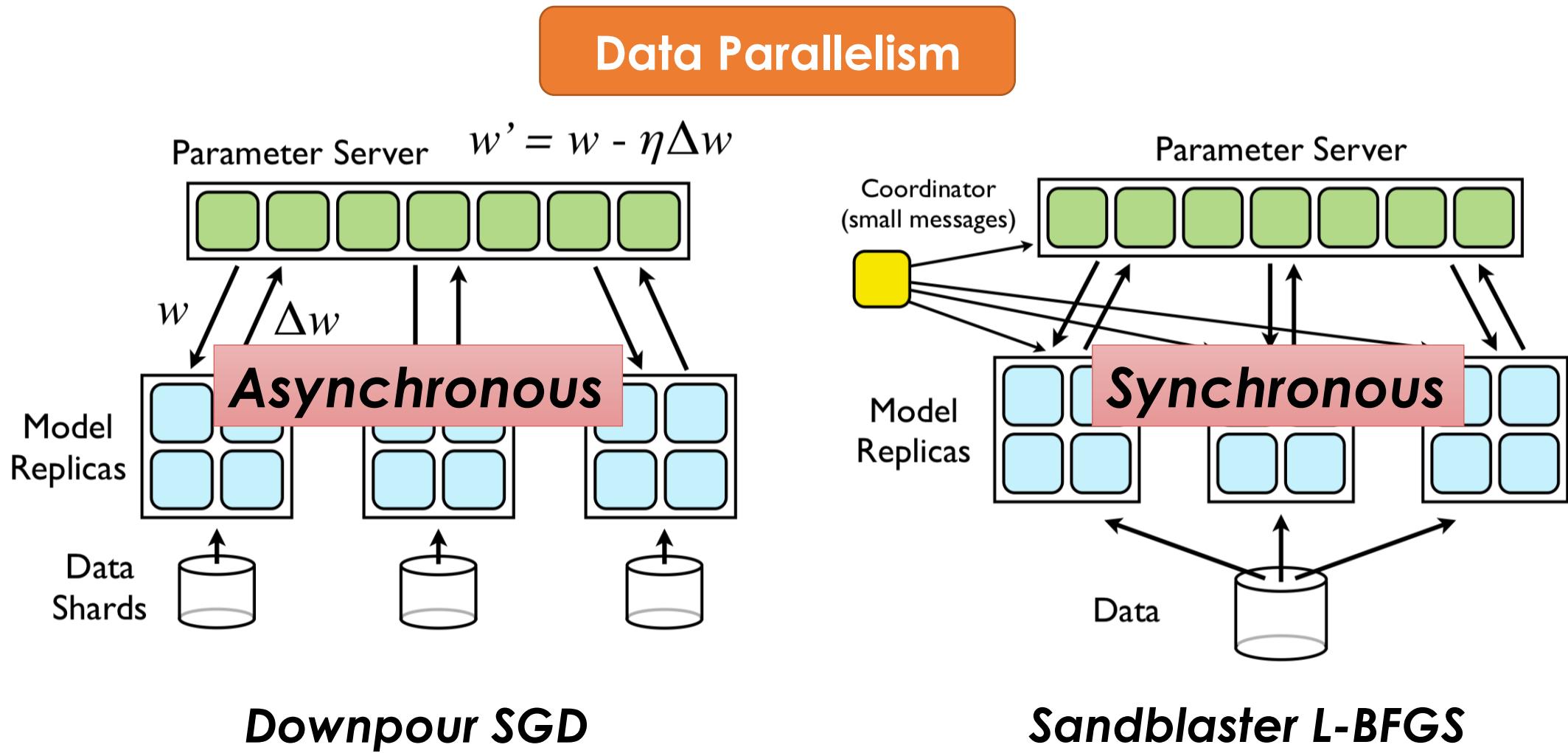
This appears in earlier work on graph systems ...



Downpour SGD

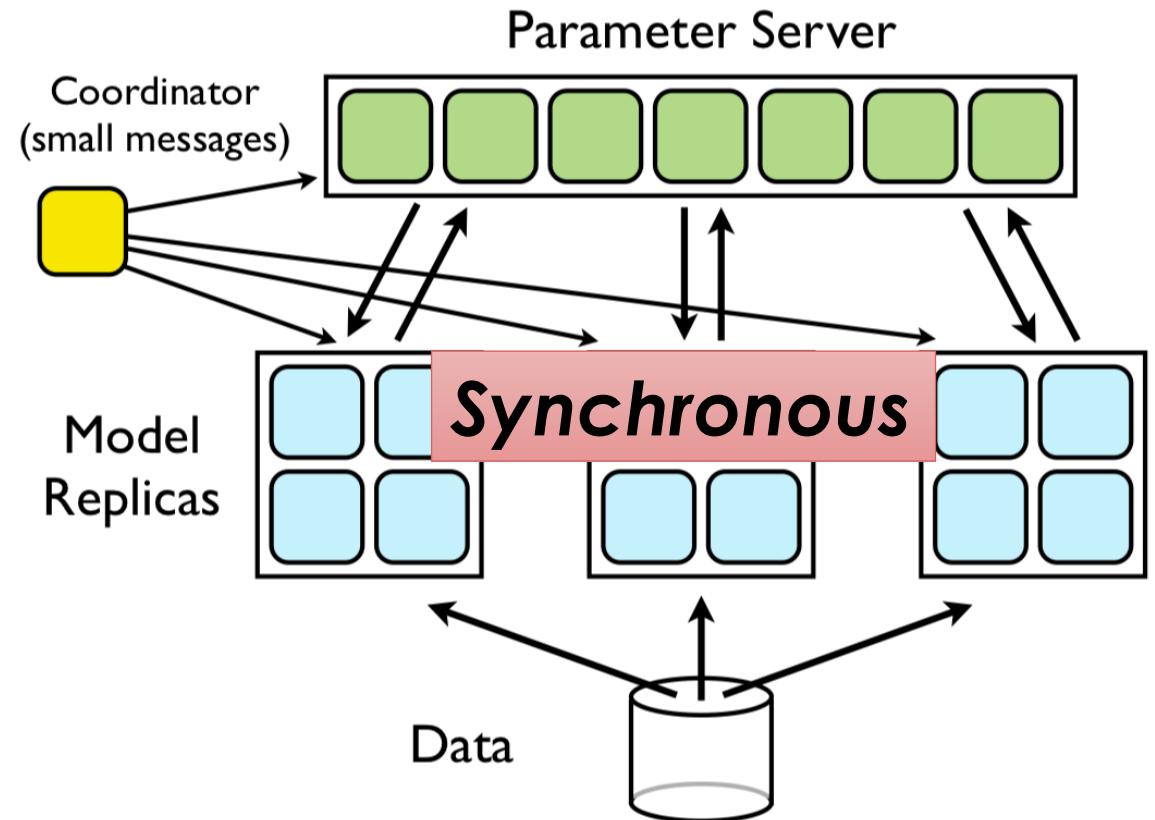
Combine Model and Data Parallelism

Machine 2
Machine 4



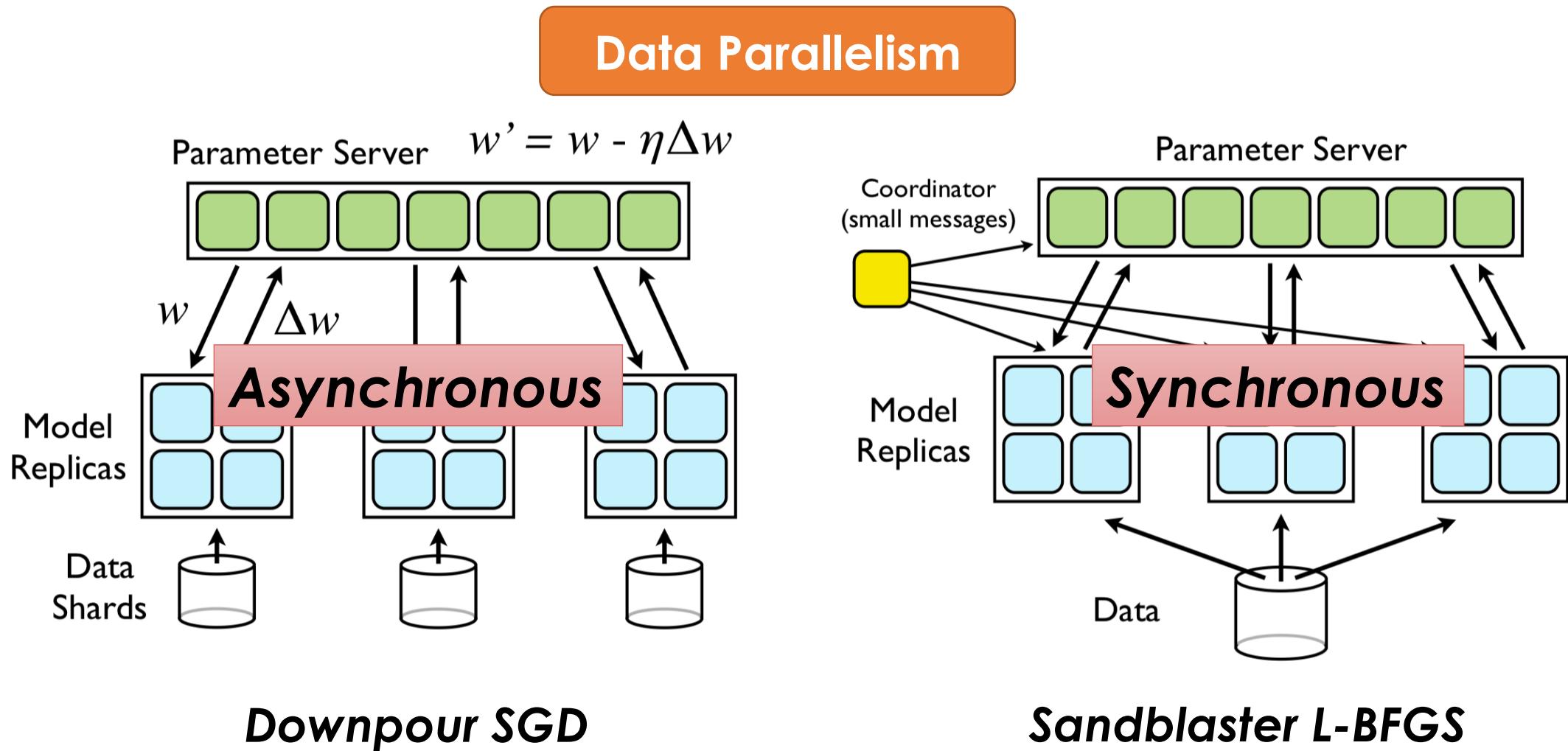
Sandblaster L-BFGS

- L-BFGS
 - Commonly used for convex opt. problems
 - Requires repeated scans of all data
 - Robust, minimal tuning
- Naturally fits map-reduce pattern
- **Innovations:**
 - accumulate gradients and store outputs in a sharded key value store (parameter server)
 - Tiny tasks + backup tasks to mitigate stragglers



Combine Model and Data Parallelism

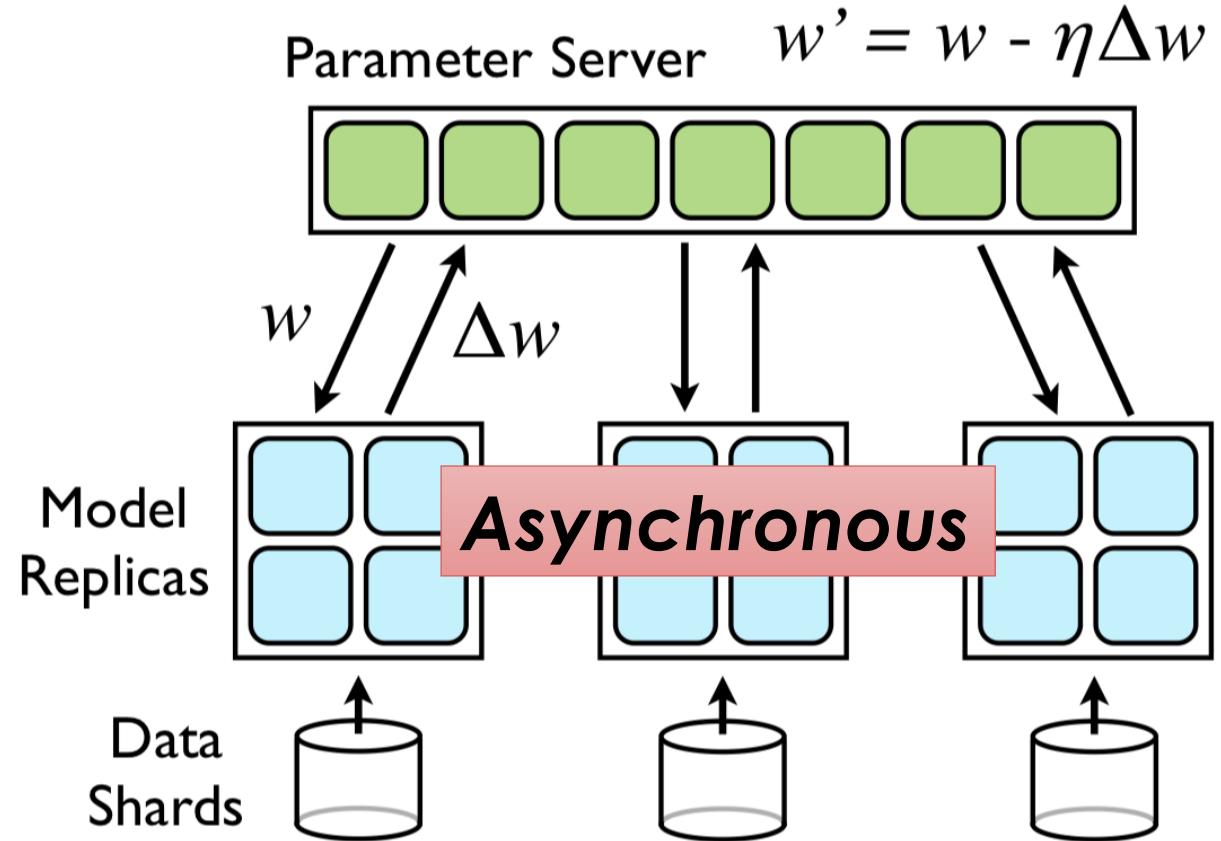
Machine 2
Machine 4



Downpour SGD

Claimed Innovations

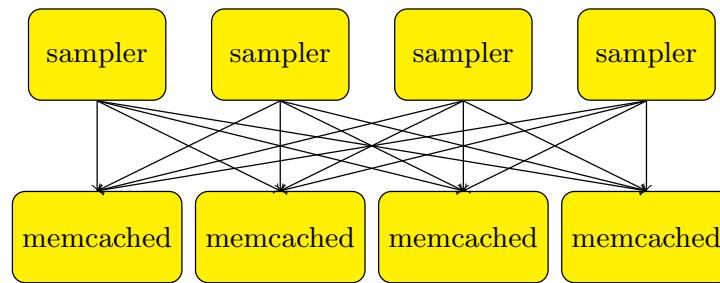
- Parameter Server
- Combine model and data parallelism in an async. execution.
- Adagrad stabilization
- Warmstarting



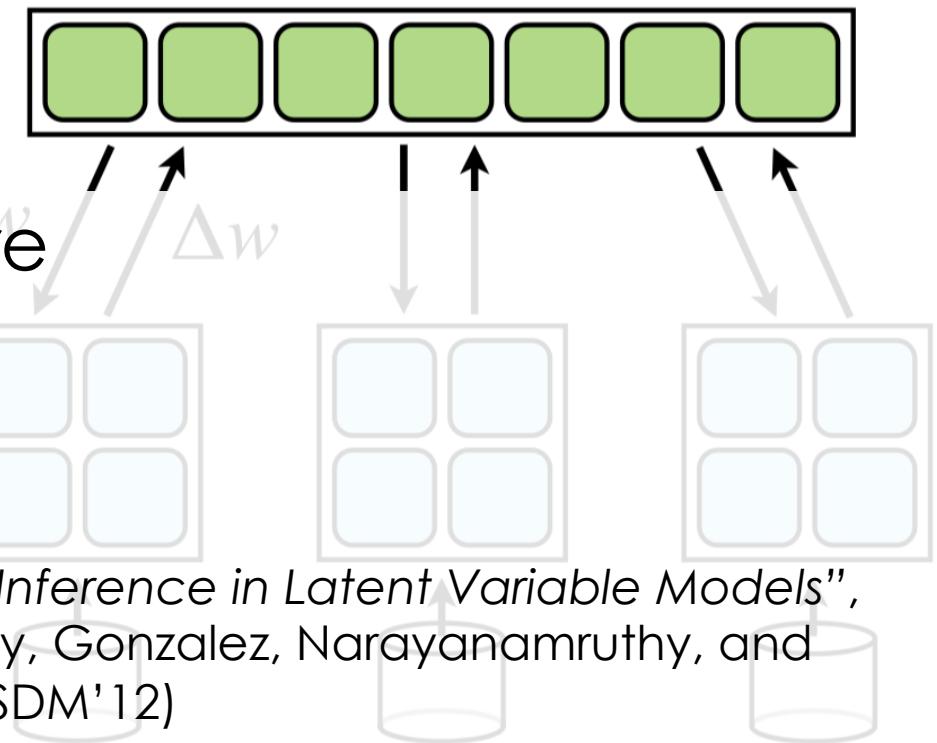
Parameter Servers

- Essentially a **sharded** key-value store
 - support for put, get, **add**
- Idea appears in earlier papers:

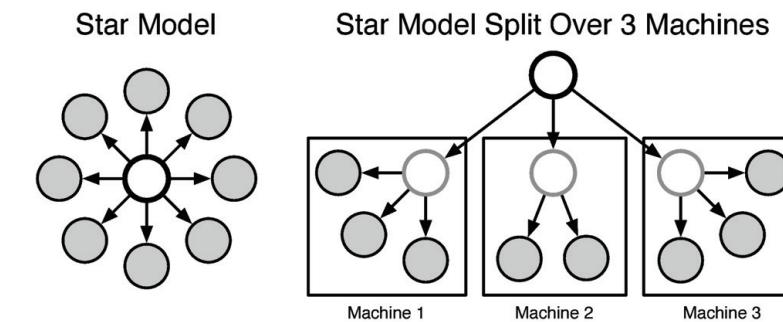
"An Architecture for Parallel Topic Models", Smola and Narayananamruthy.
(VLDB'10)



$$\text{Parameter Server} \quad w' = w - \eta \Delta w$$



"Scalable Inference in Latent Variable Models",
Ahmed, Aly, Gonzalez, Narayananamruthy, and
Smola. (WSDM'12)

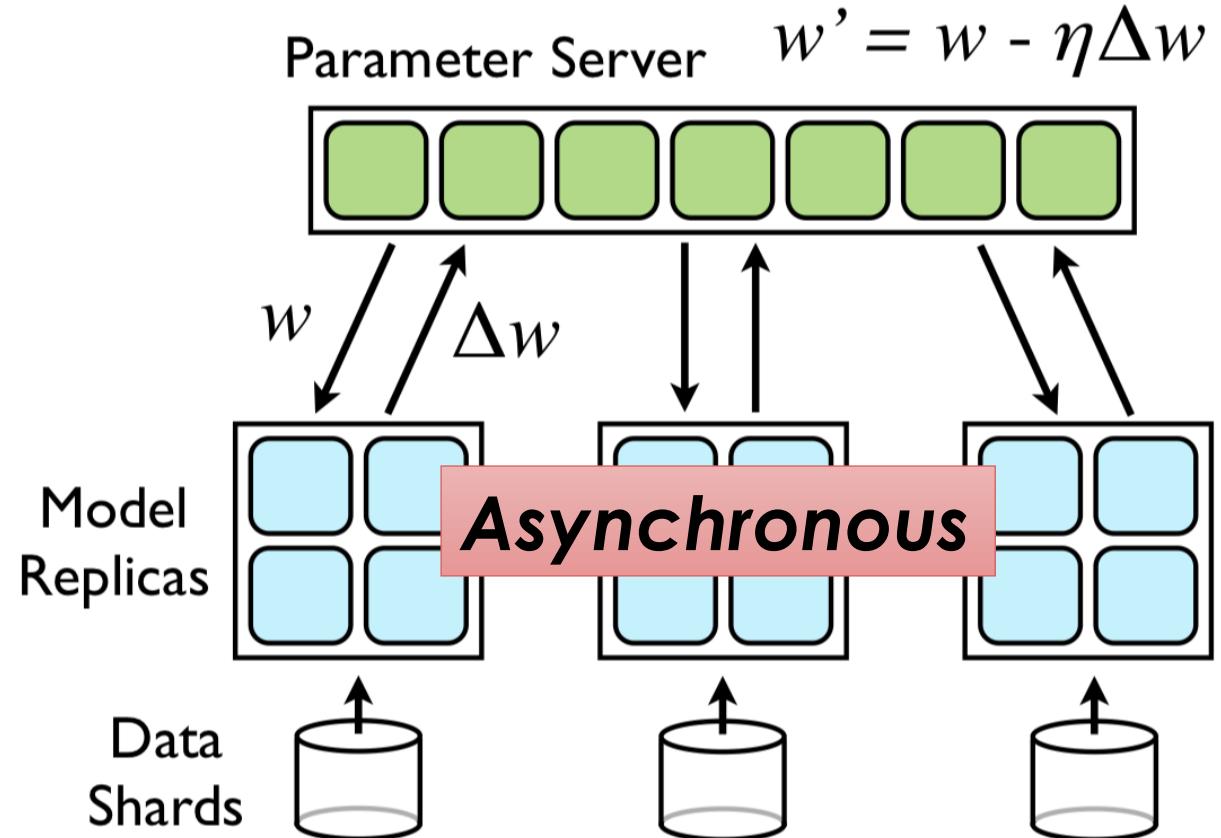


DistBelief was probably the first paper to call a sharded key-value store a Parameter Server.

Downpour SGD

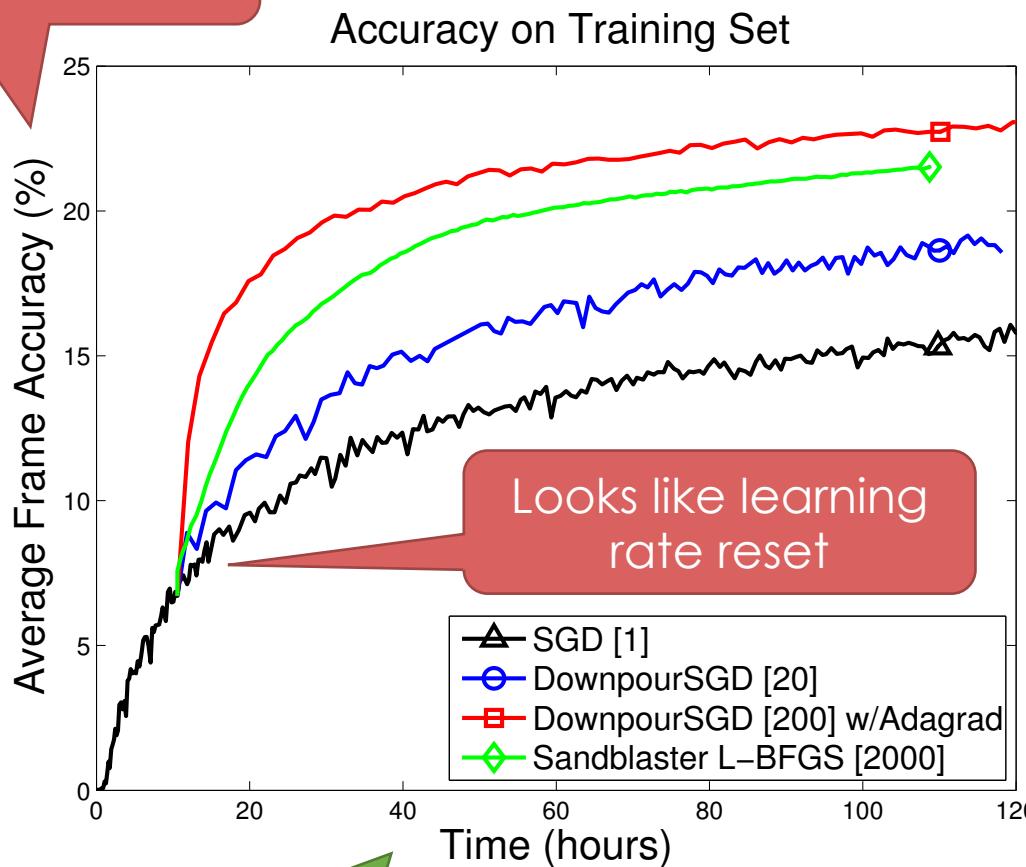
Claimed Innovations

- Parameter Server
- Combine model and data parallelism in an **async. execution.**
- Adagrad stabilization
- Warmstarting

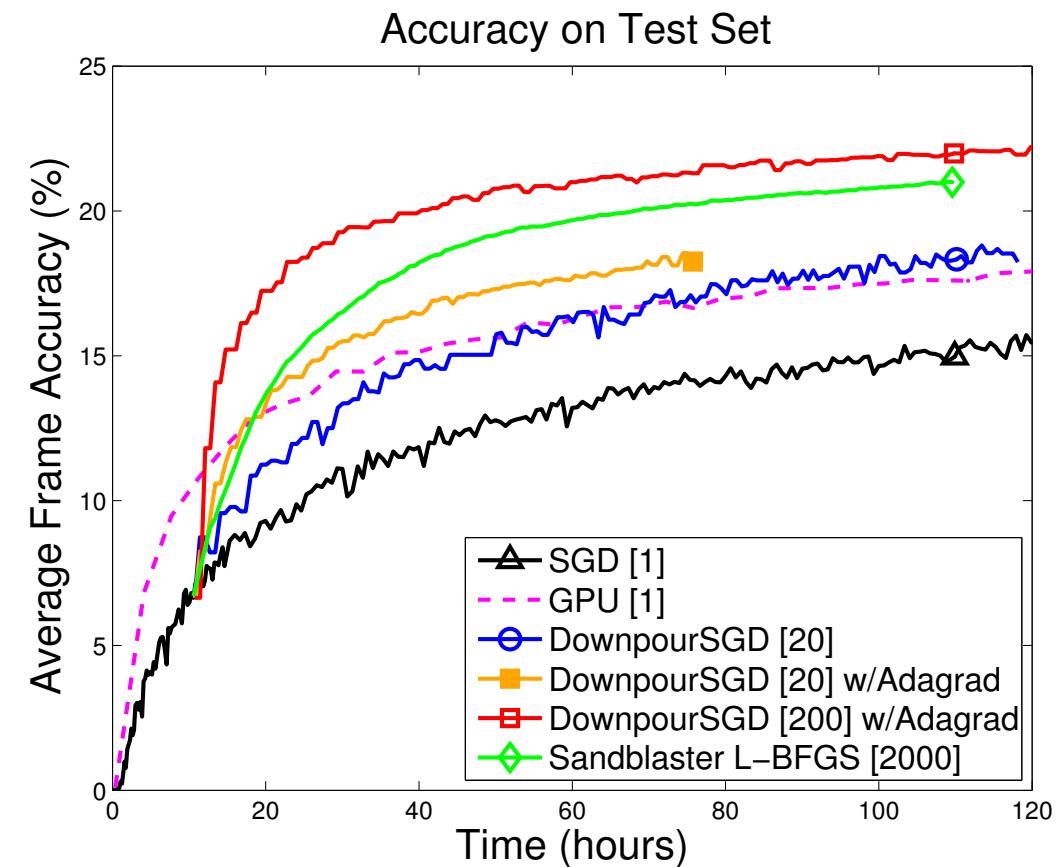


Key Results: Training and Test Error

Weird 20K
Error Metric



Wall clock
time is good.



Why are they in the NY Times

- Trained a 1.7 billion parameter model (30x larger than state-of-the-art) (was it necessary?)
- Using 16,000 cores (efficiently?)
- Achieves 15.8 accuracy on ImageNet 20K (70% improvement over state of the art).
 - Non-standard benchmark
- Qualitatively interesting results



Figure 6. Visualization of the cat face neuron (left) and human body neuron (right).

Long-term Impact

- The **parameter server** appears in many later machine learning systems
- Downpour (**Asynchronous**) SGD has been largely **replaced by synchronous systems** for supervised training
 - Asynchrony is still popular in RL research
 - Why?
- Model parallelism is still used for large language models
 - Predated this work
- The neural network architectures studied here have been largely replaced by convolutional networks

More recent large-scale training

- Generated a lot of press
 - Surpassed by
Fast.ai: “Now anyone can train ImageNet in 18 minutes for \$40.” blog post
 - Popularized linear learning rate scaling

2018 (Unpublished on Arxiv)

**Accurate, Large Minibatch SGD:
Training ImageNet in 1 Hour**

Priya Goyal Piotr Dollár Ross Girshick Pieter Noordhuis
Lukasz Wesolowski Aapo Kyrola Andrew Tulloch Yangqing Jia Kaiming He
Facebook

arXiv:1706.02677v2 [cs.CV] 30 Apr 2018

Abstract

Deep learning thrives with large neural networks and large datasets. However, larger networks and larger datasets result in longer training times that impede research and development progress. Distributed synchronous SGD offers a potential solution to this problem by dividing SGD minibatches over a pool of parallel workers. Yet to make this scheme efficient, the per-worker workload must be large, which implies nontrivial growth in the SGD minibatch size. In this paper, we empirically show that on the ImageNet dataset large minibatches cause optimization difficulties, but when these are addressed the trained networks exhibit good generalization. Specifically, we show no loss of accuracy when training with large minibatch sizes up to 8192 images. To achieve this result, we adopt a hyper-parameter-free linear scaling rule for adjusting learning rates as a function of minibatch size and develop a new warmup scheme that overcomes optimization challenges early in training. With these simple techniques, our Caffe2-based system trains ResNet-50 with a minibatch size of 8192 on 256 GPUs in one hour, while matching small minibatch accuracy. Using commodity hardware, our implementation achieves ~90% scaling efficiency when moving from 8 to 256 GPUs. Our findings enable training visual recognition models on internet-scale data with high efficiency.

1. Introduction

Scale matters. We are in an unprecedented era in AI research history in which the increasing data and model scale is rapidly improving accuracy in computer vision [22, 41, 34, 35, 36, 16], speech [17, 40], and natural language processing [7, 38]. Take the profound impact in computer vision as an example: visual representations learned by deep convolutional neural networks [23, 22] show excellent performance on previously challenging tasks like ImageNet classification [33] and can be transferred to difficult perception problems such as object detection and segmen-

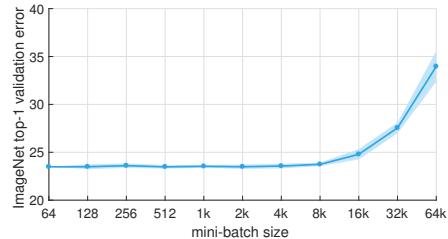


Figure 1. **ImageNet top-1 validation error vs. minibatch size.** Error range of plus/minus two standard deviations is shown. We present a simple and general technique for scaling distributed synchronous SGD to minibatches of up to 8k images while maintaining the top-1 error of small minibatch training. For all minibatch sizes we set the learning rate as a linear function of the minibatch size and apply a simple warmup phase for the first few epochs of training. All other hyper-parameters are kept fixed. Using this simple approach, accuracy of our models is invariant to minibatch size (up to an 8k minibatch size). Our techniques enable a linear reduction in training time with ~90% efficiency as we scale to large minibatch sizes, allowing us to train an accurate 8k minibatch ResNet-50 model in 1 hour on 256 GPUs.

tion [8, 10, 28]. Moreover, this pattern generalizes: larger datasets and neural network architectures consistently yield improved accuracy across all tasks that benefit from pre-training [22, 41, 34, 35, 36, 16]. But as model and data scale grow, so does training time; discovering the potential and limits of large-scale deep learning requires developing novel techniques to keep training time manageable.

The goal of this report is to demonstrate the feasibility of, and to communicate a practical guide to, large-scale training with distributed *synchronous* stochastic gradient descent (SGD). As an example, we scale ResNet-50 [16] training, originally performed with a minibatch size of 256 images (using 8 Tesla P100 GPUs, training time is 29 hours), to larger minibatches (see Figure 1). In particular, we show that *with a large minibatch size of 8192, we can train ResNet-50 in 1 hour using 256 GPUs while maintaining*

Contrasting to the first paper

- **Synchronous SGD**
 - Much of the recent work has focused on synchronous setting
 - Easier to reason about
- Focus exclusively on data parallelism: **batch-size scaling**
- Focuses on the **generalization gap problem**

How do you distribute SGD?

Stochastic Gradient Descent

$\theta^{(0)} \leftarrow$ initial vector (random, zeros ...)

For t from 0 to convergence:

$\mathcal{B} \sim$ Random subset of indices

$$\theta^{(t+1)} \leftarrow \theta^{(t)} - \eta_t \left(\frac{1}{|\mathcal{B}|} \sum_{i \in \mathcal{B}} \nabla_{\theta} \mathbf{L}(y_i, f(x_i; \theta)) \Big|_{\theta=\theta^{(t)}} \right)$$

Data
Parallelism

Slow? (~150ms)
Depending on size of B

Batch Size Scaling

- Increase the batch size by adding machines

$$\theta^{(t+1)} \leftarrow \theta^{(t)} - \hat{\eta} \left(\frac{1}{k} \sum_{j=1}^k \frac{1}{|\mathcal{B}_j|} \sum_{i \in \mathcal{B}_j} \nabla_{\theta} \mathbf{L}(y_i, f(x_i; \theta)) \Big|_{\theta=\theta^{(t)}} \right)$$

- Each server processes a fixed batch size (e.g., n=32)
- As more servers are added (k) the effective overall batch size increases linearly
- Why do these additional servers help?

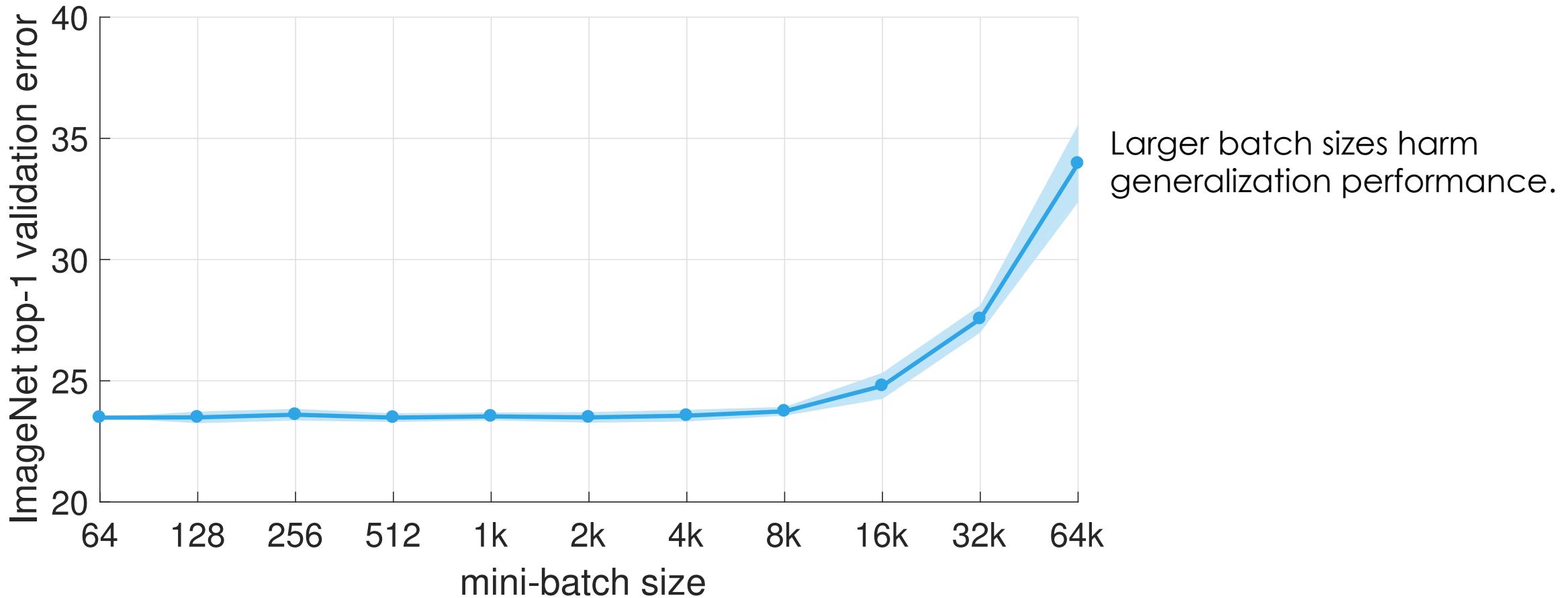
Bigger isn't Always Better

- Motivation for larger batch sizes
 - More opportunities for parallelism → but is it useful?
 - Recall ($1/n$ variance reduction):

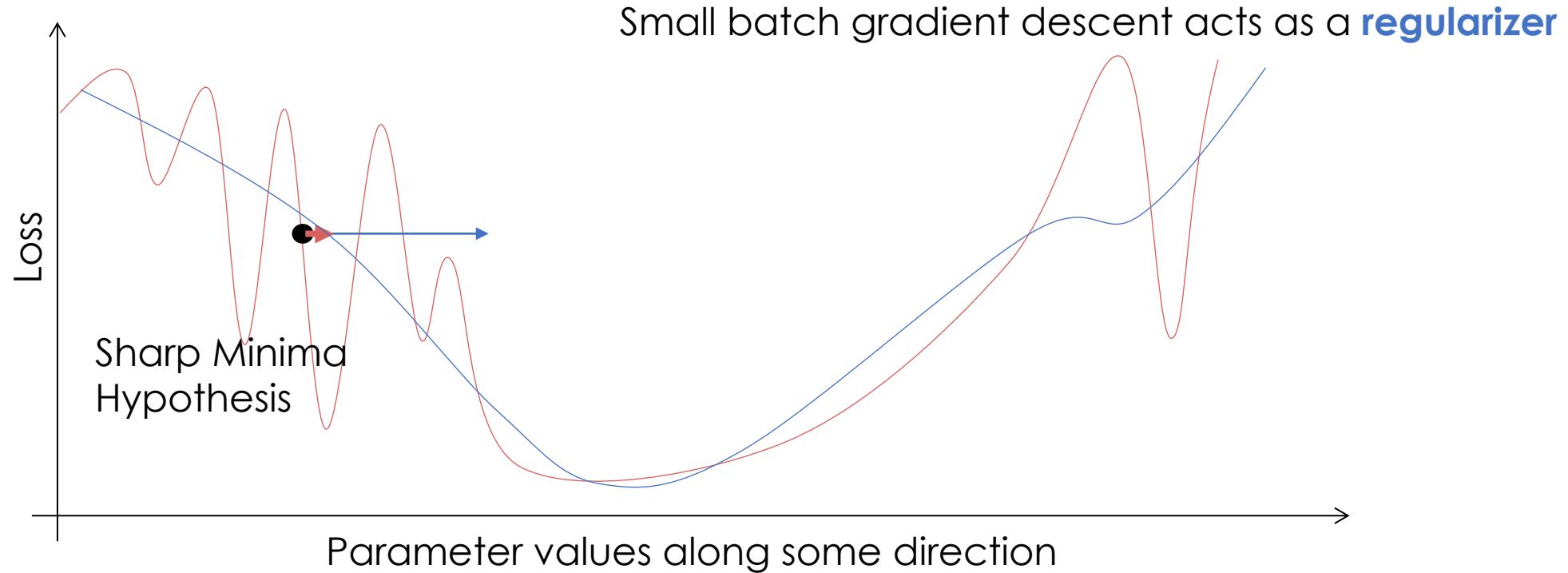
$$\frac{1}{n} \sum_{i=1} \nabla_{\theta} \mathbf{L}(y_i, f(x_i; \theta)) \approx \frac{1}{|\mathcal{B}|} \sum_{i \in \mathcal{B}} \nabla_{\theta} \mathbf{L}(y_i, f(x_i; \theta))$$

- Is a variance reduction helpful?
 - Only if it lets you take bigger steps (move faster)
 - Does it affect the final prediction accuracy?

Generalization Gap Problem



Rough “Intuition”



Key problem: Addressing the generalization gap for large batch sizes.

Solution: Linear Scaling Rule

- Scale the learning rate linearly with the batch size

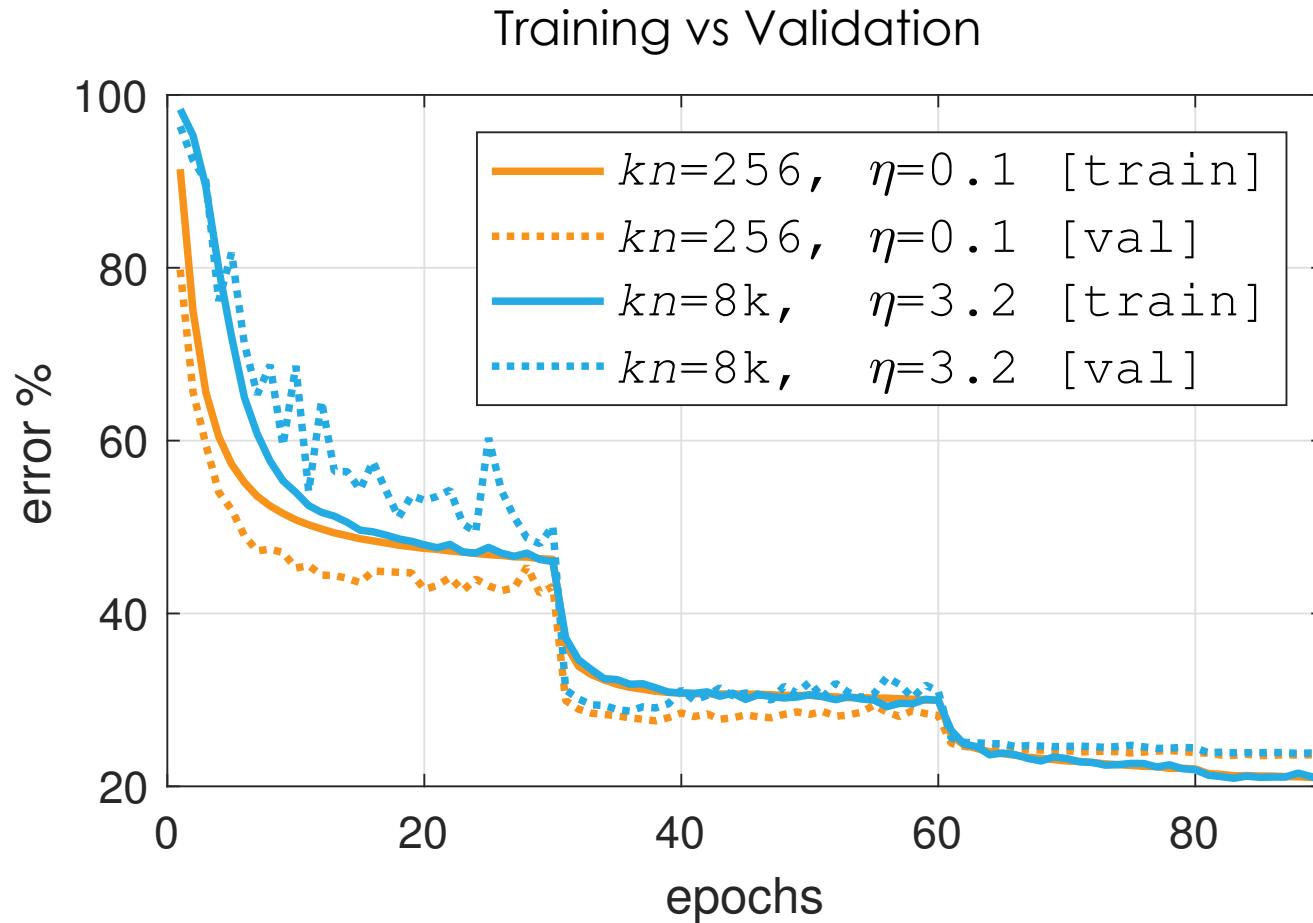
$$\theta^{(t+1)} \leftarrow \theta^{(t)} - \hat{\eta} \left(\frac{1}{k} \sum_{j=1}^k \frac{1}{|\mathcal{B}_j|} \sum_{i \in \mathcal{B}_j} \nabla_{\theta} \mathbf{L}(y_i, f(x_i; \theta)) \Big|_{\theta=\theta^{(t)}} \right)$$

- Addresses generalization performance by **taking larger steps** (also improves training convergence)
- **Sub-problem:** Large learning rates can be destabilizing in the beginning. Why?
 - **Gradual warmup solution:** increase learning rate scaling from constant to linear in first few epochs
 - Doesn't help for very large k...

Other Details

- **Independent Batch Norm:** Batch norm calculation applies only to local batch size (n).
- **All-Reduce:** Recursive halving and doubling algorithm
 - Used instead of popular ring reduction (fewer rounds)
- **Gloo** a library for efficient collective communications
- **Big Basin GPU Servers:** Designed for deep learning workloads
 - Analysis of communication requirements → latency bound
- **No discussion on straggler or fault-tolerance**
 - Why?!

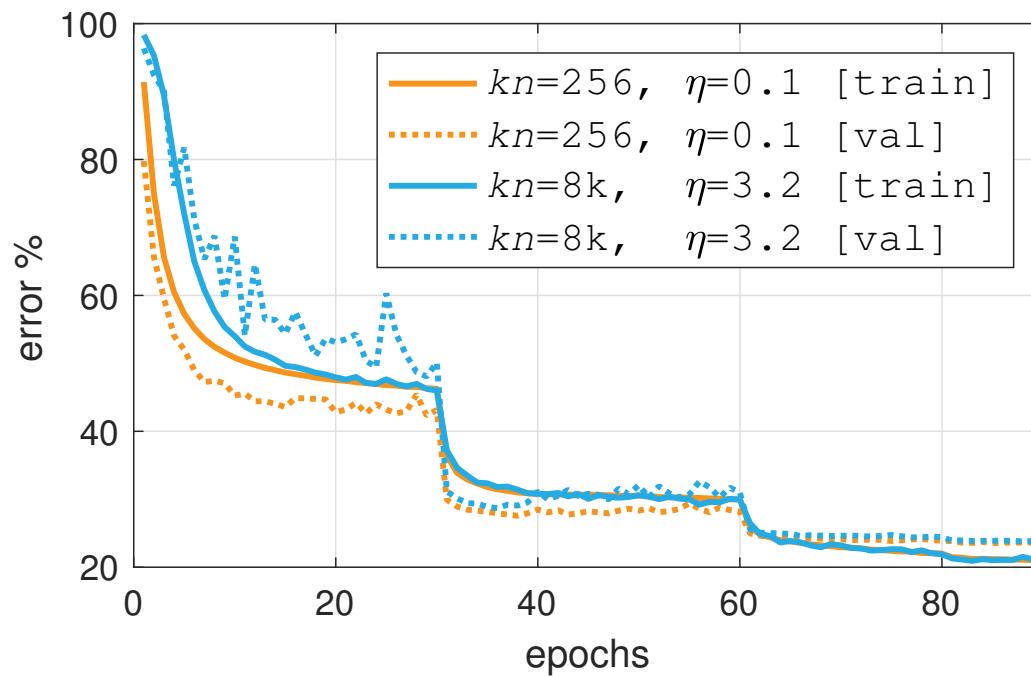
Key Results



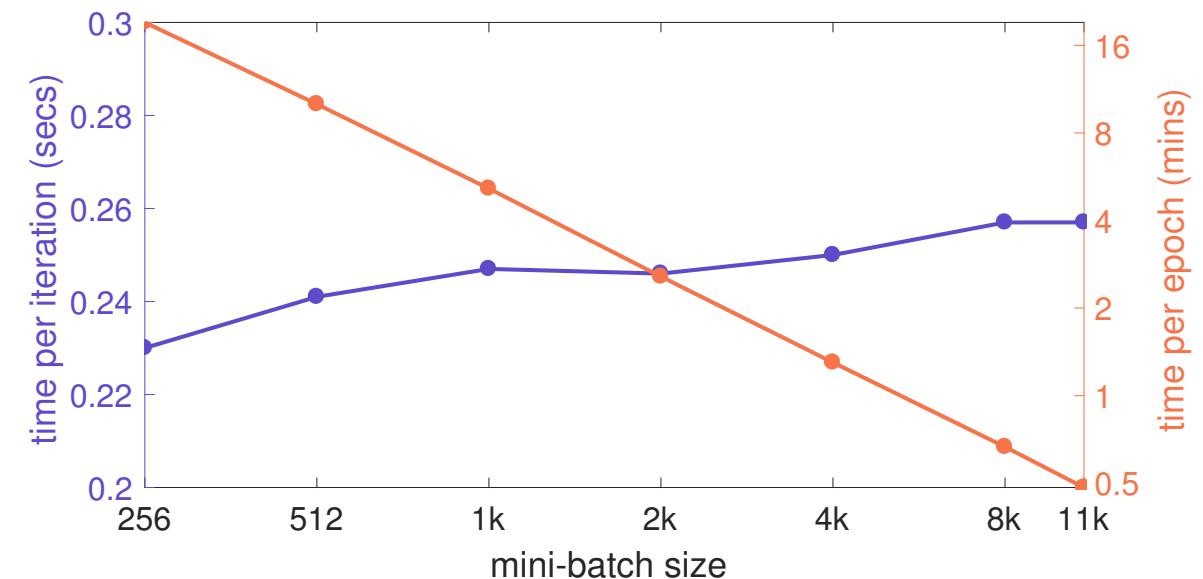
All curves closely match using the linear scaling rule.

Note learning rate schedule drops.

Key Results



“Learning”
Epoch
Machine Learning



Epoch
Second
System

Key Results

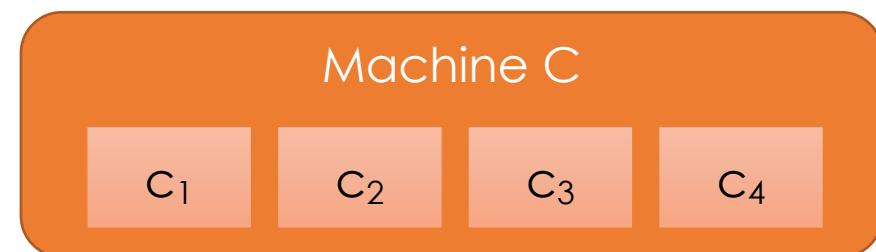
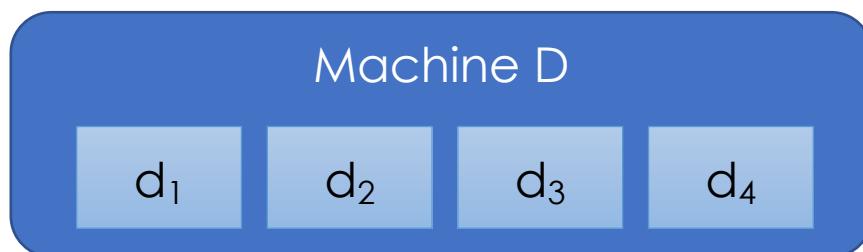
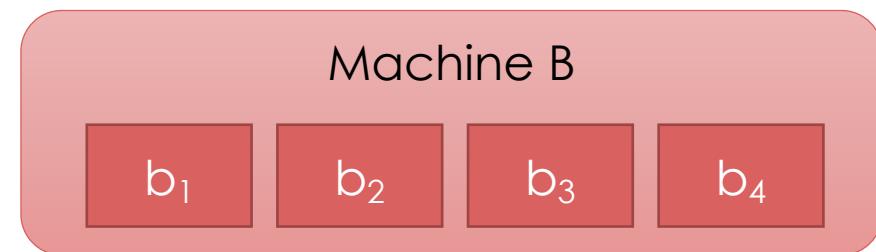
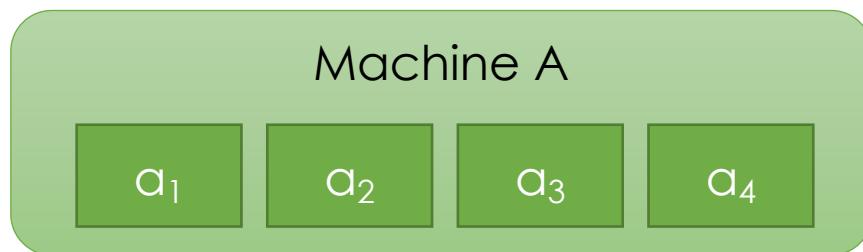
- Train ResNet-50 to state-of-the-art on 256 GPUs in 1 hour
 - 90% scaling efficiency
- Fairly careful study of the linear scaling rule
 - Observed limits to linear scaling do not depend on dataset size
 - Cannot scale parallelism with dataset size

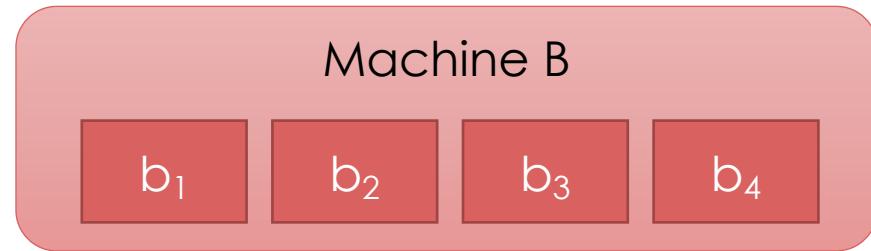
All-Reduce

All Reduce

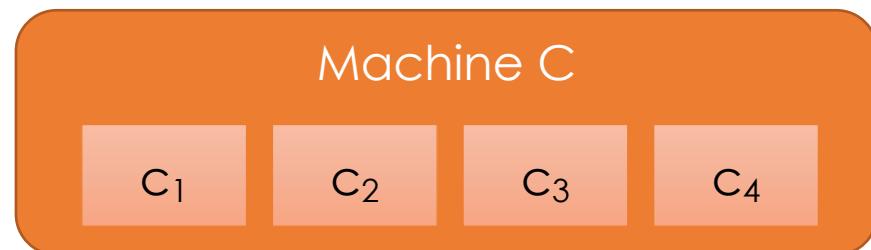
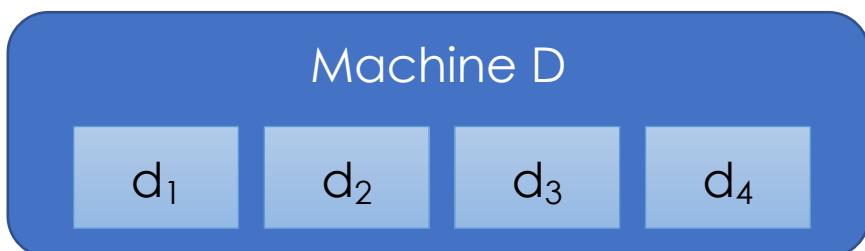
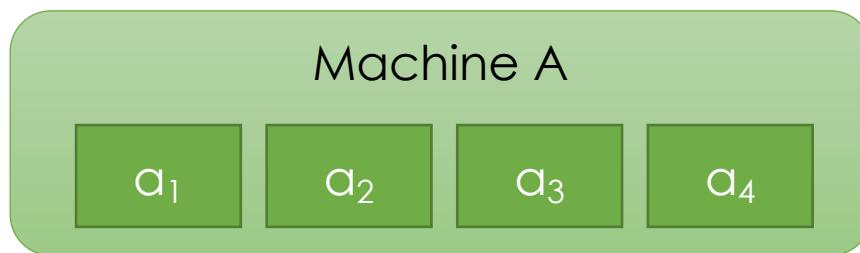
Mechanism to sum and distribute data across machines.

- Used to sum and distribute the gradient



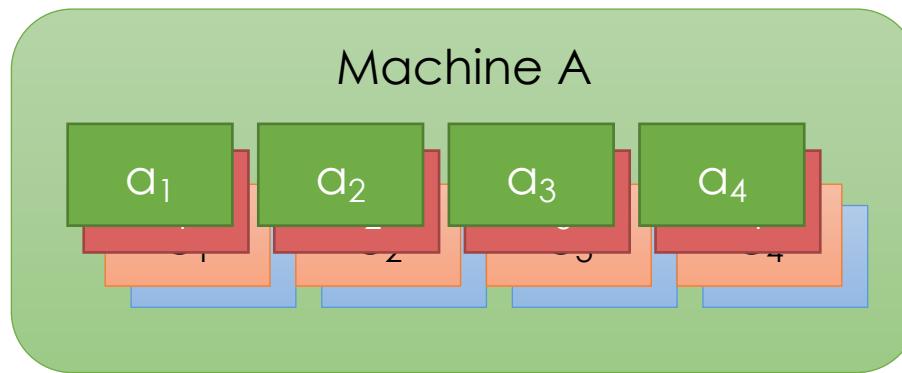


Single Master All-Reduce



Single Master All-Reduce

Machine B



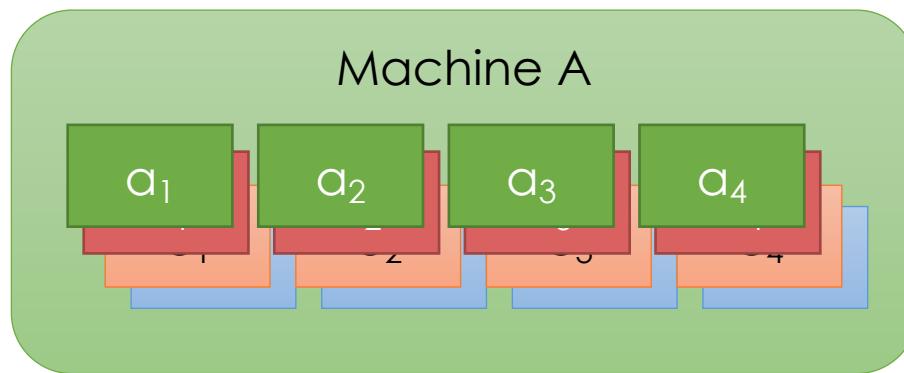
Sends **(P-1) * N** Data
➤ **P** Machines
➤ **N** Parameters

Machine D

Machine C

Single Master All-Reduce

Machine B



Sends **(P-1) * N** Data
➤ **P** Machines
➤ **N** Parameters

$$s_i = a_i + b_i + c_i + d_i$$

Machine D

Machine C

Single Master All-Reduce

Machine B

Machine A

s_1 s_2 s_3 s_4

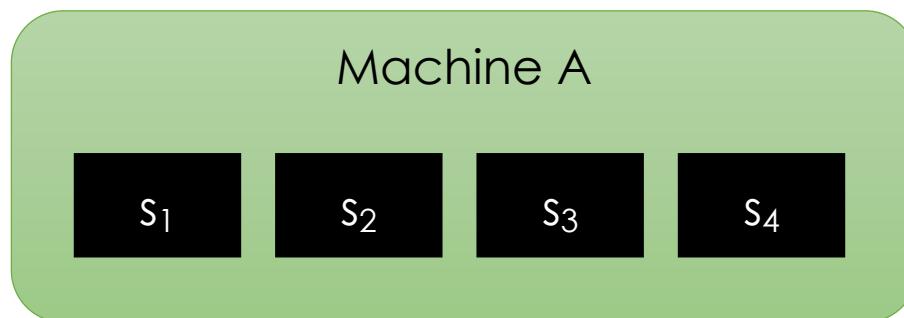
Sends **(P-1) * N** Data
➤ **P** Machines
➤ **N** Parameters

$$s_i = a_i + b_i + c_i + d_i$$

Machine D

Machine C

Single Master All-Reduce



Sends **(P-1) * N** Data
➤ **P** Machines
➤ **N** Parameters

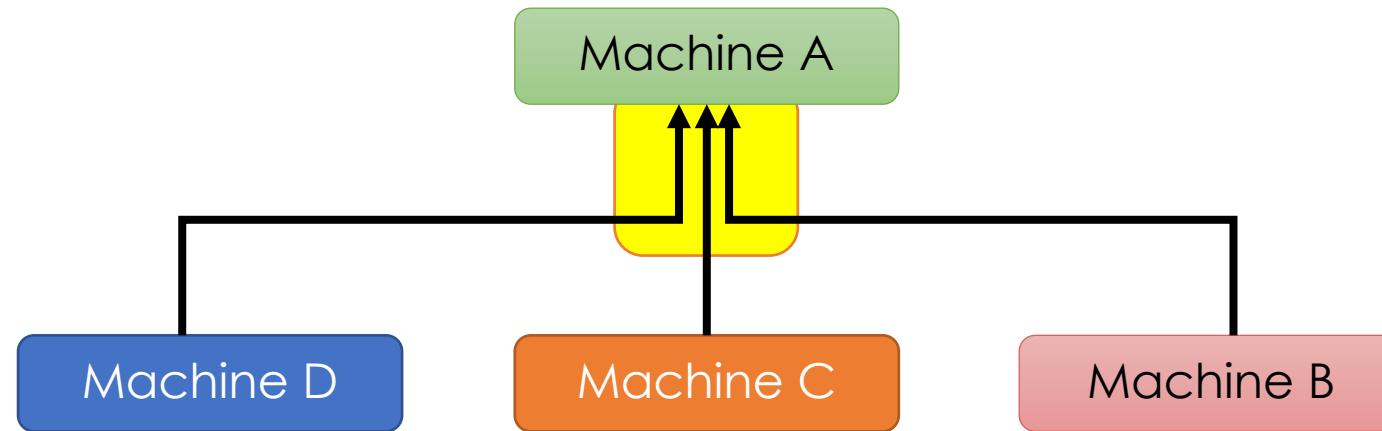
^{*2}

$$s_i = a_i + b_i + c_i + d_i$$



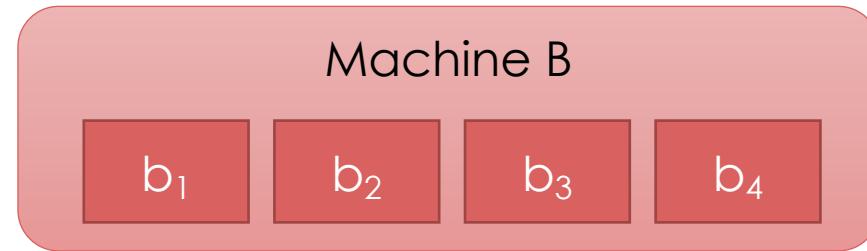
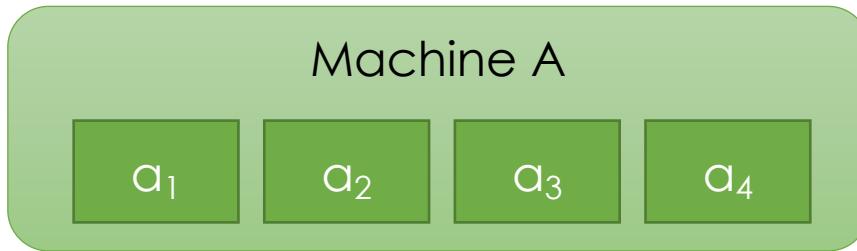
Single Master All-Reduce

Sends $(P-1) * N$ Data
➤ P Machines
➤ N Parameters

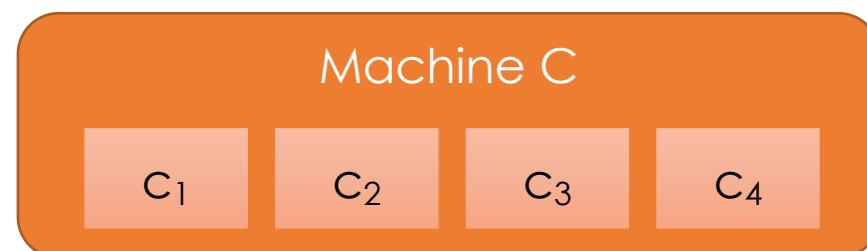
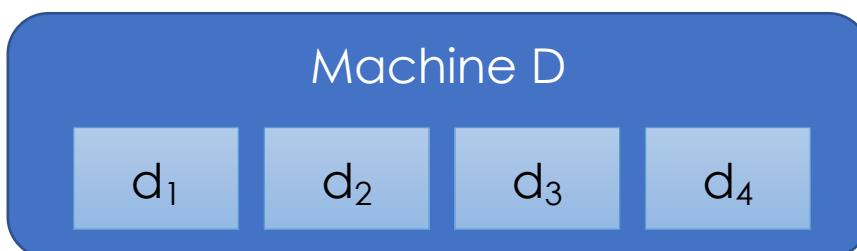


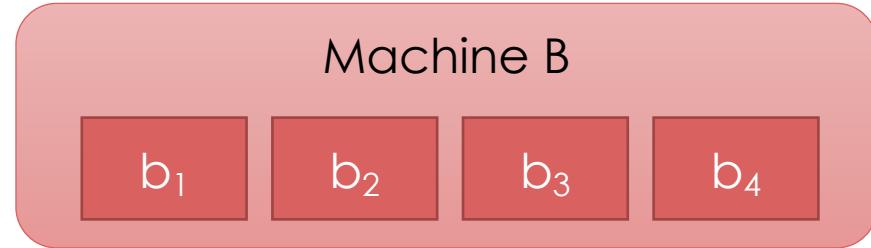
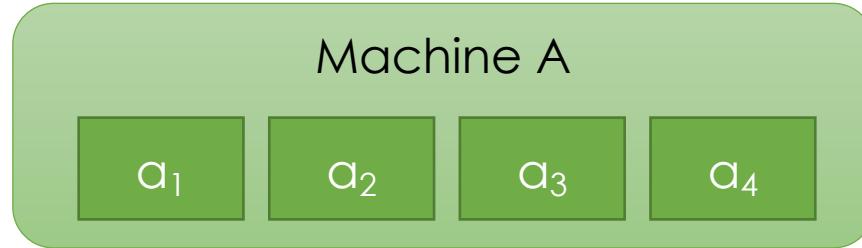
Issues?

- High **fan-in** on Machine A
- **$(P-1) * N$ Bandwidth** for Machine A



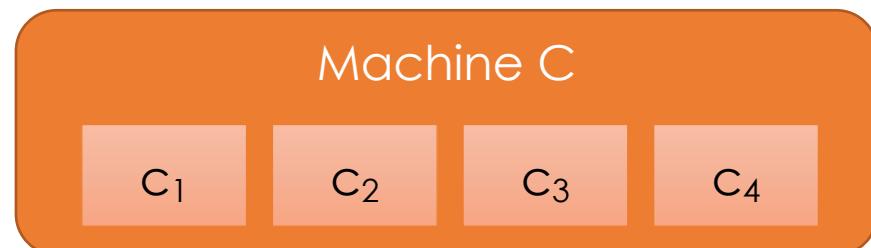
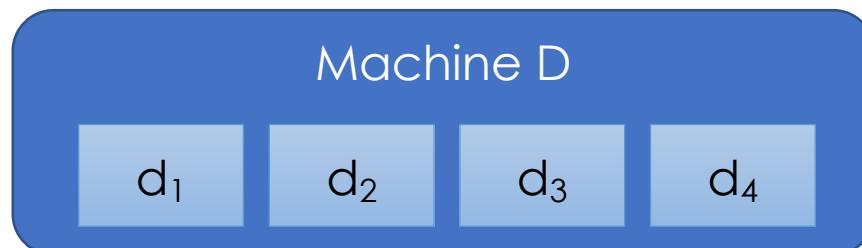
Parameter Server All Reduce

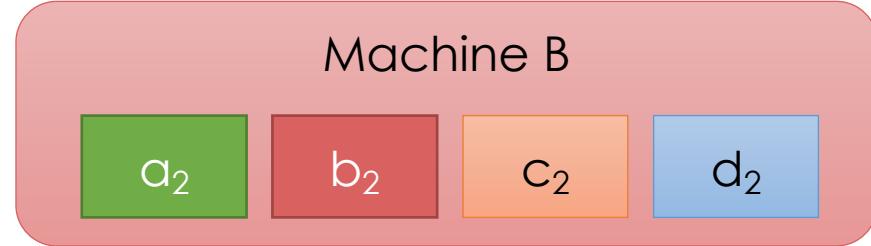
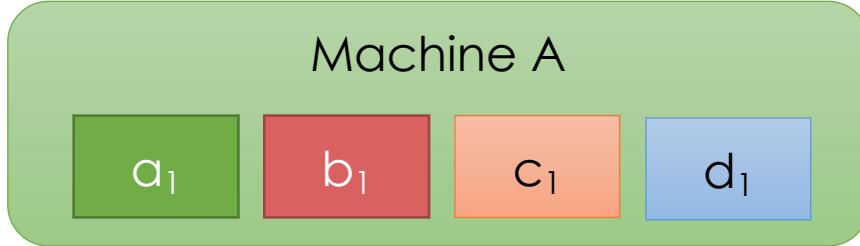




Send each entry to parameter server for that entry.

- Key 1 → A
- Key 2 → B
- Key 3 → C
- Key 4 → D

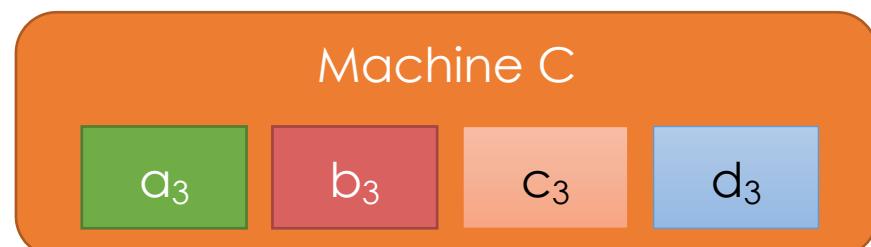




Each machine sends N/P data to all other machines.

$$P * (P-1) * N/P = (P-1) * N$$

- **P** Machines
- **N** Parameters



Machine A

s_1

Machine B

s_2

Compute local sum on each machine

$$s_i = a_i + b_i + c_i + d_i$$

Machine D

s_4

Machine C

s_3

Machine A

s_1

Machine B

s_2

Broadcast sum to each machine

Machine D

s_4

Machine C

s_3

Machine A

s_1

s_2

s_3

s_4

Machine B

s_1

s_2

s_3

s_4

Broadcast sum to each machine

Machine D

s_1

s_2

s_3

s_4

Machine C

s_1

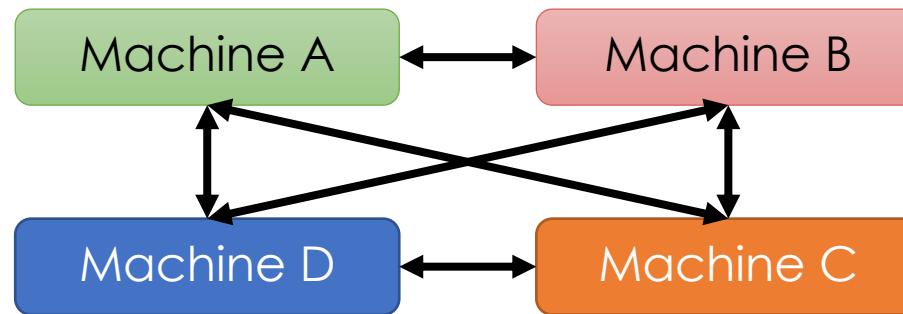
s_2

s_3

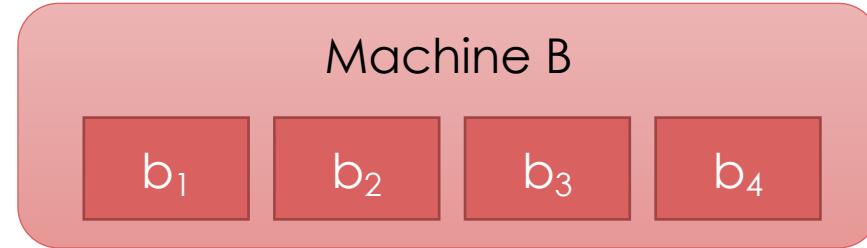
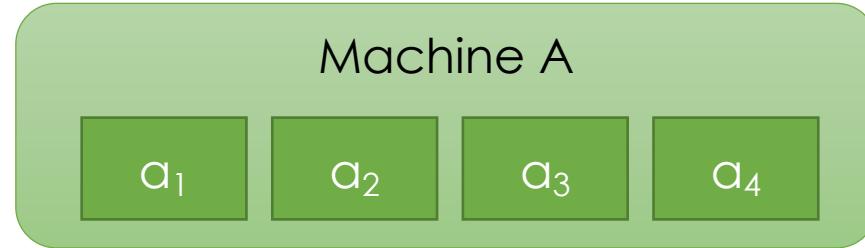
s_4

Parameter Server All-Reduce

- Same amount of data transmitted as before

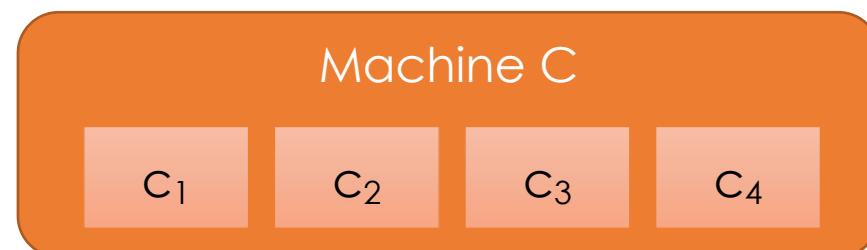
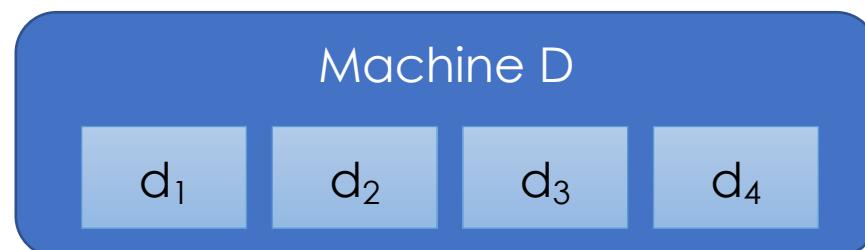


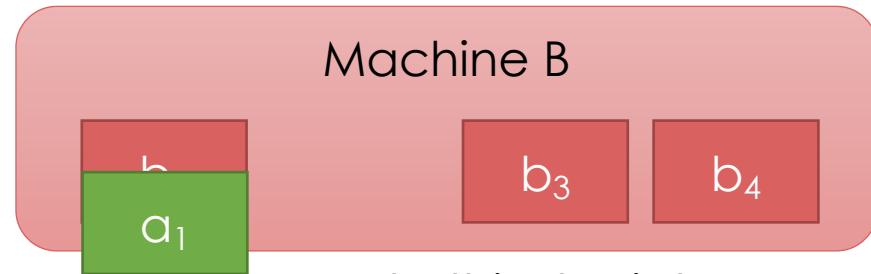
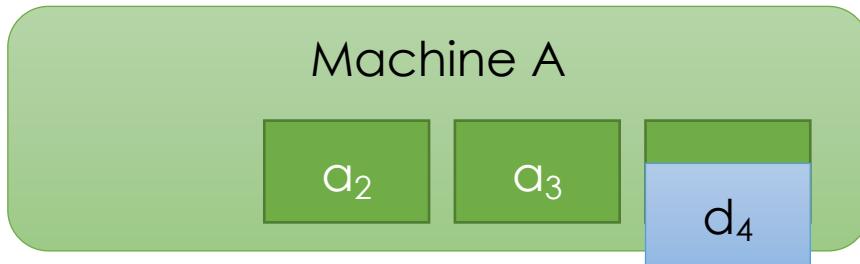
- Same **high fan-in** ($P-1$)
- **Reduced** Inbound Bandwidth = $(P-1)N/P$
 - Previously $(P-1)*N$



Ring All Reduce

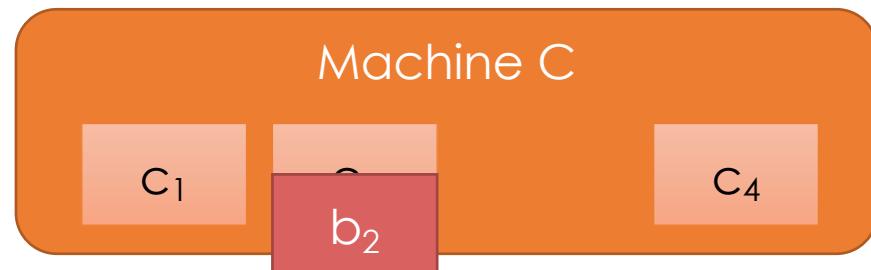
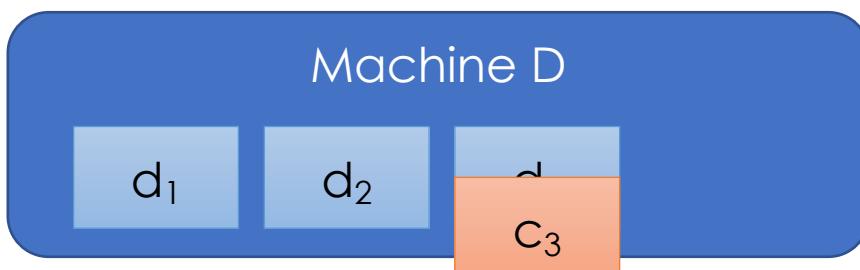
Send messages in a ring using to reduce fan-in.



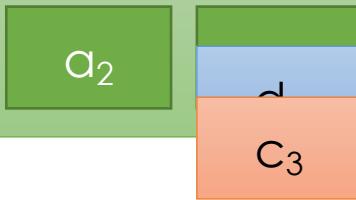


← Note this depicts a partial sum and not a bigger message.

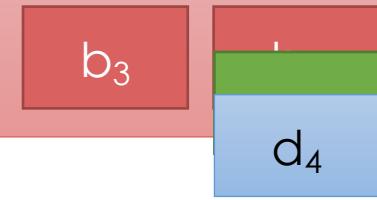
Ring All Reduce



Machine A



Machine B



Ring All Reduce

Machine D



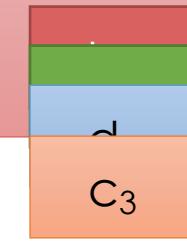
Machine C



Machine A



Machine B

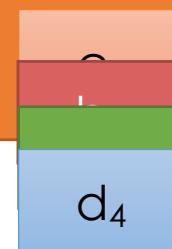


Ring All Reduce

Machine D



Machine C



Machine A

s_2

Machine B

s_3

Ring All Reduce

Each machine sends N/P data to next machine each of $(P-1)$ rounds:

$$(P-1) * P * N/P = (P-1) * N$$

➤ **Bandwidth** per round:

➤ $P(N/P) = N$ (doesn't depend on P)

➤ **Fan-in Per Round:**

➤ 1 (doesn't depend on P)

Machine D

s_1

Machine C

s_4

Machine A

s_2

Machine B

s_3

Ring All Reduce

Broadcast stage repeats process sending messages forwarding sums (same communication costs).

Machine D

s_1

Machine C

s_4

Machine A

s_1

s_2

Machine B

s_2

s_3

Ring All Reduce

Machine D

s_1

s_4

Machine C

s_3

s_4

Machine A

s_1

s_2

s_4

Machine B

s_1

s_2

s_3

Ring All Reduce

Machine D

s_1

s_3

s_4

Machine C

s_2

s_3

s_4

Machine A

s_1

s_2

s_3

s_4

Machine B

s_1

s_2

s_3

s_4

Ring All Reduce

Machine D

s_1

s_2

s_3

s_4

Machine C

s_1

s_2

s_3

s_4

Machine A

s_1

s_2

s_3

s_4

Machine B

s_1

s_2

s_3

s_4

Ring All Reduce

Machine D

s_1

s_2

s_3

s_4

Machine C

s_1

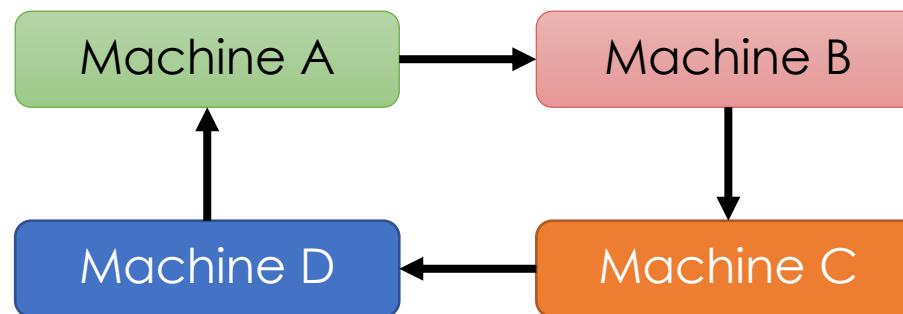
s_2

s_3

s_4

Ring All-Reduce

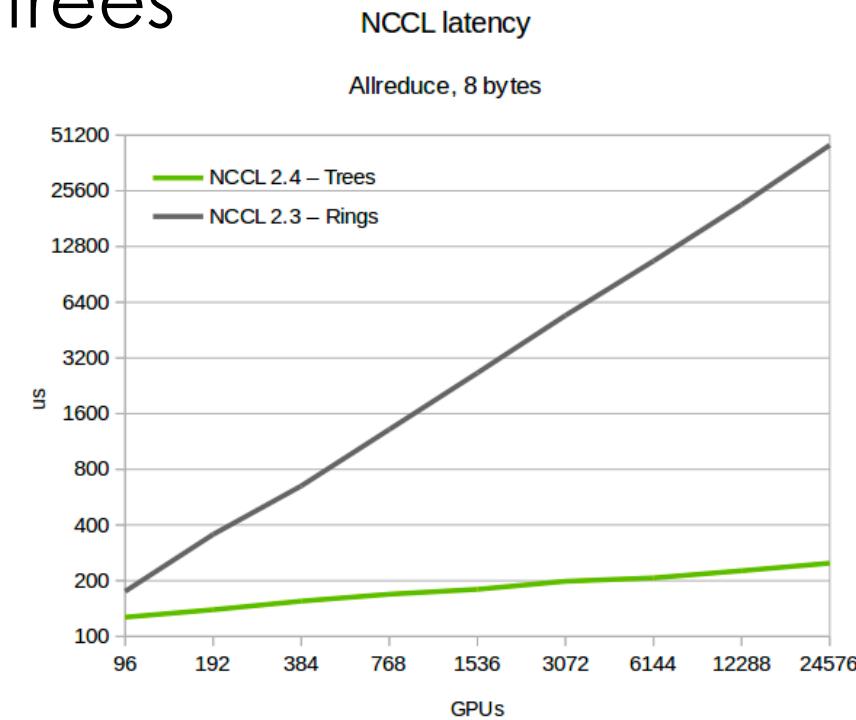
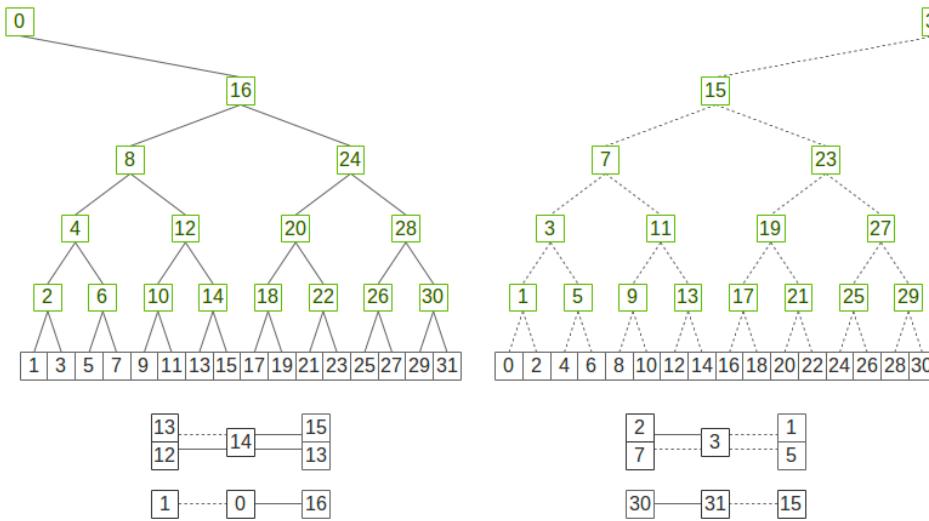
- Simplified communication topology with low fan-in



- Overall communication
 - Same total communication: $2*(P-1)*N$
 - **Bandwidth** per round (N) doesn't depend on P
 - **Fan-in** is constant (doesn't depend on P)
- **Issue:** Number of communication rounds ($P-1$)

Double Binary Tree All-Reduce

- Two overlaid binary reduction trees



- Double the fan-in → $\log(p)$ rounds of communication
 - Currently used on Summit super-computer and latest NCCL

Review:

Dimensions of Parallelism

Data Parallelism

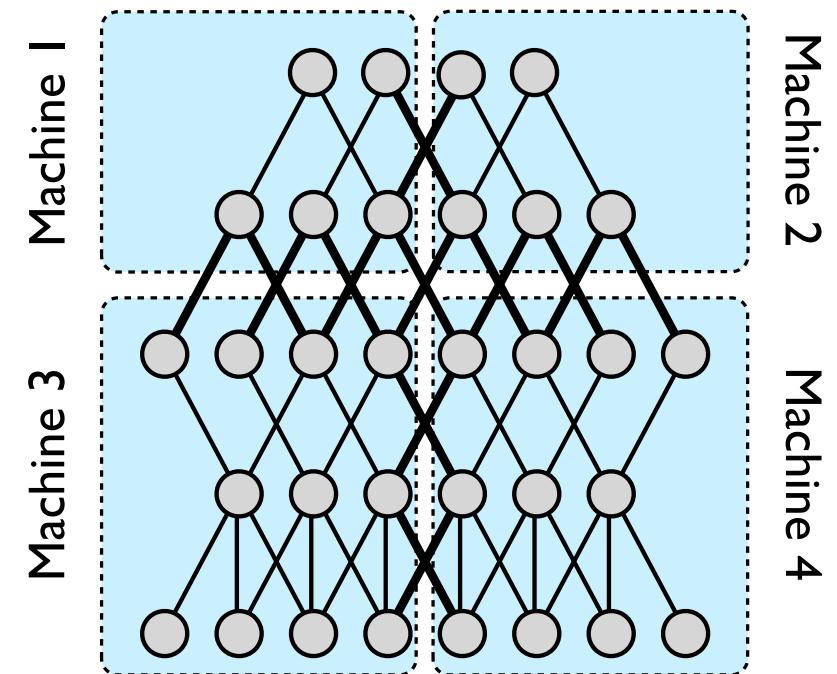
Parallelizing mini-batch gradient calculation with model replicated to all machines.

- Synchronous Execution (Most Common)
 - **Strengths:** deterministic, parallelism does not effect result
 - **Weaknesses:** need large batch sizes, frequent blocking comm., learning rate scaling, doesn't work with batch normalization
- Asynchronous Execution (Popular in Research)
 - **Strengths:** eliminate blocking and use background comm., batches don't need to span machines
 - **Weaknesses:** affects convergence (stability)
- Issues:
 - Model and activations must fit in each machine

Model Parallelism

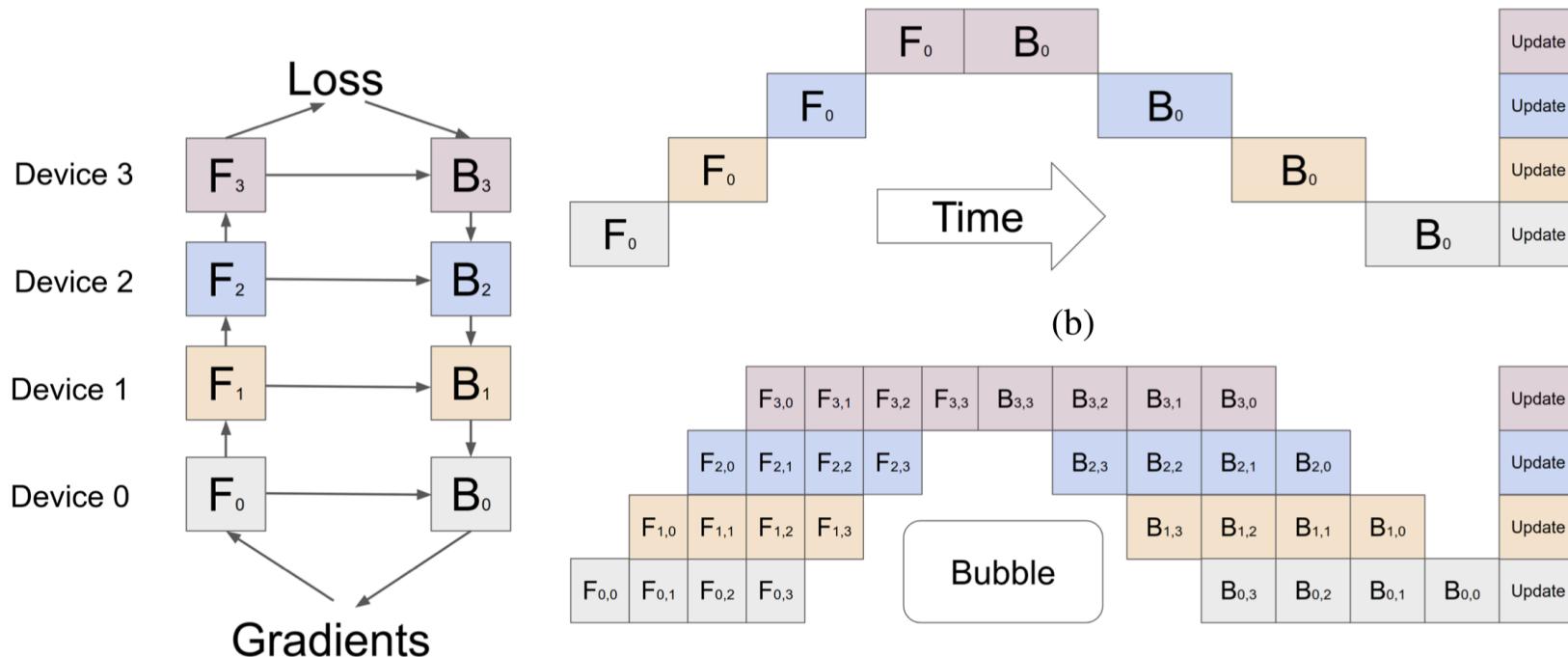
Divide the model across machines and replicate the data.

- Supports large models and activations
- Requires communication within single evaluation
- How to best divide a model?
 - Split individual layers
 - which dimension?
 - Batch or Spatial → depends on operation
 - Split across layers
 - Only one set of layers active a time → poor work balance
 - Soln: Pipelining Parallelism



Pipeline Parallelism

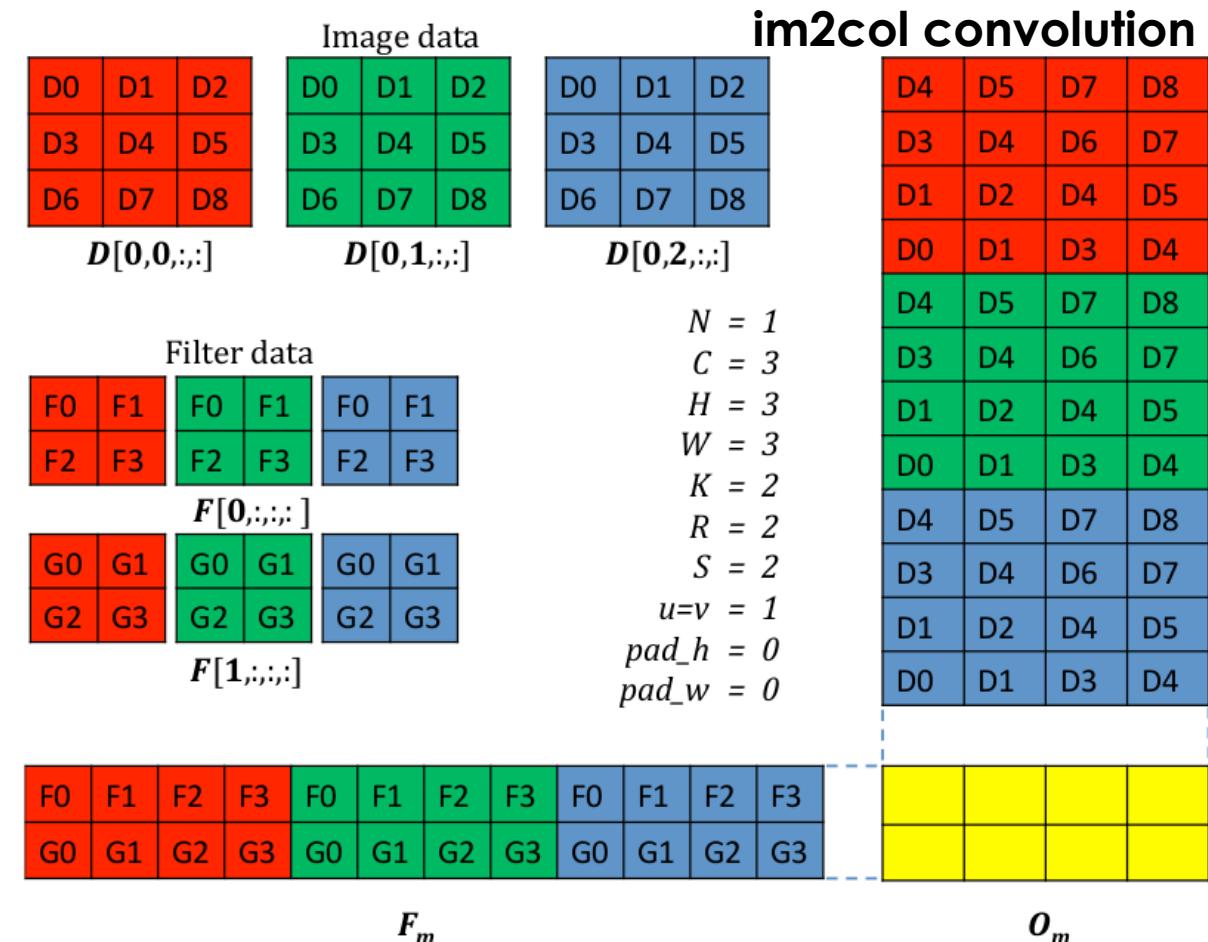
- Combine model and data parallelism to concurrently process multiple layers and batches.
- Originally described in GPipe*



*[GPipe: Easy Scaling with Micro-Batch Pipeline Parallelism](#)

Operator Level Parallelism

- Exploiting the parallelism within linear algebra and convolution operations (a form of model parallelism)
- Multiple dimensions
 - Batch, spatial, time, ...
- Typically cast operators as linear alg. routines and leverage optimizes BLAS libraries

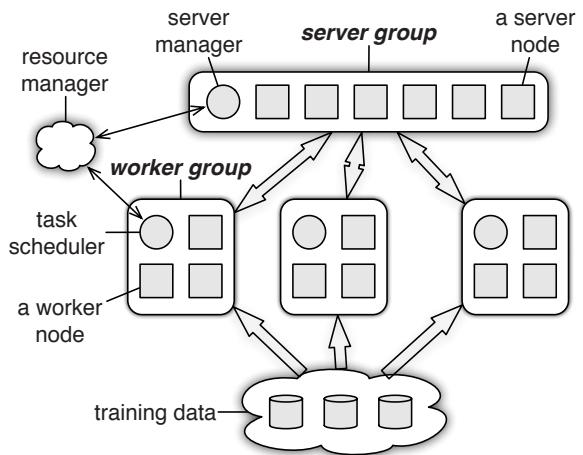


This weeks readings

Reading for the Week

- [Scaling Distributed Machine Learning with the Parameter Server](#) (OSDI'14)
 - Paper describing the parameter server system
- [PipeDream: Generalized Pipeline Parallelism for DNN Training](#) (SOSP'19)
 - Latest paper exploring pipeline parallel training
- [Adaptive Communication Strategies to Achieve the Best Error-Runtime Trade-off in Local-Update SGD](#) (SysML'19)
 - Dynamic averaging approach to distributed training

Scaling Distributed Machine Learning with the Parameter Server (OSDI'14)

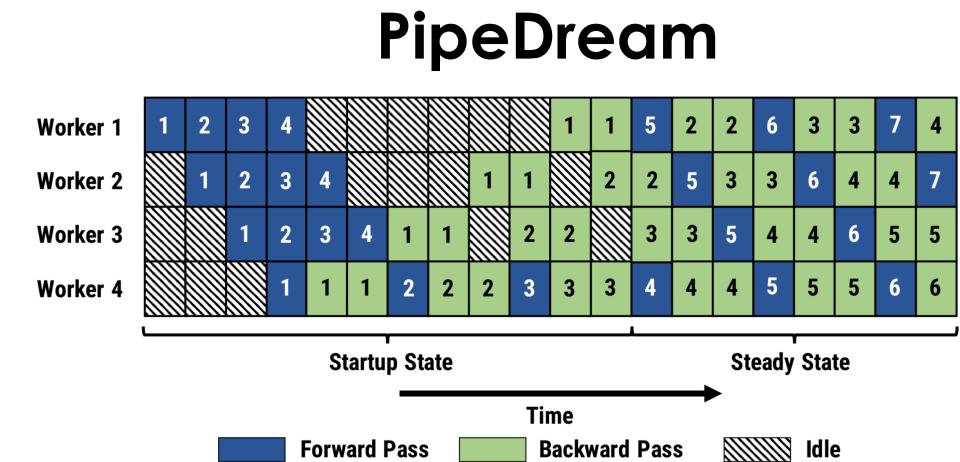
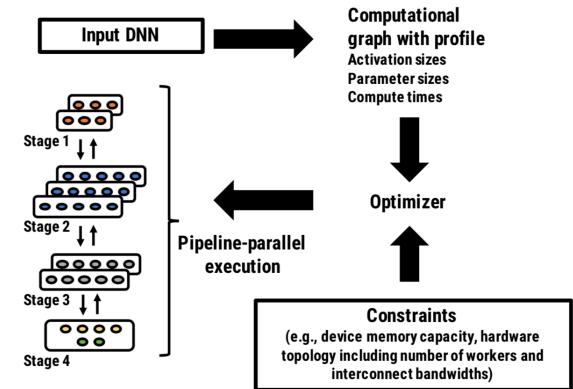
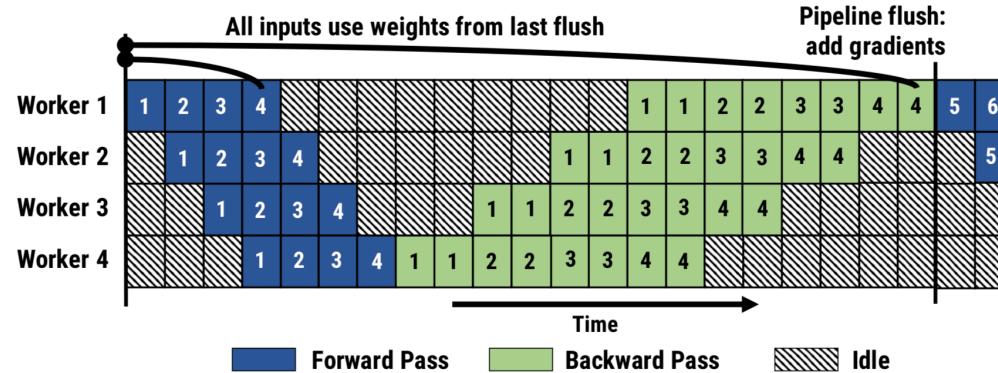


- Describes the key-value store customized for machine learning
 - Builds on earlier work in parameter servers
- **Additional Context:** focused on topic modeling and sparse regression
- **Key Ideas:** There are many ideas ...
 - Keys – Value pairs with **linear algebra** semantics (e.g., get by range)
 - User defined **event handlers** on parameter servers and workers
 - Several different **consistency models**
 - **User defined filters** to determine when updates are communicated

PipeDream: Generalized Pipeline Parallelism for DNN Training (SOSP'19)

- Contemporaneously published with:
 - [GPipe: Efficient Training of Giant Neural Networks using Pipeline Parallelism](#) (arXiv'18)
- **Key idea:** Leverage pipeline parallelism during training
 - **Automatically** constructs pipeline partition + schedule
 - Leverage **bounded staleness** + **versioned activations** to eliminate **bubbles**

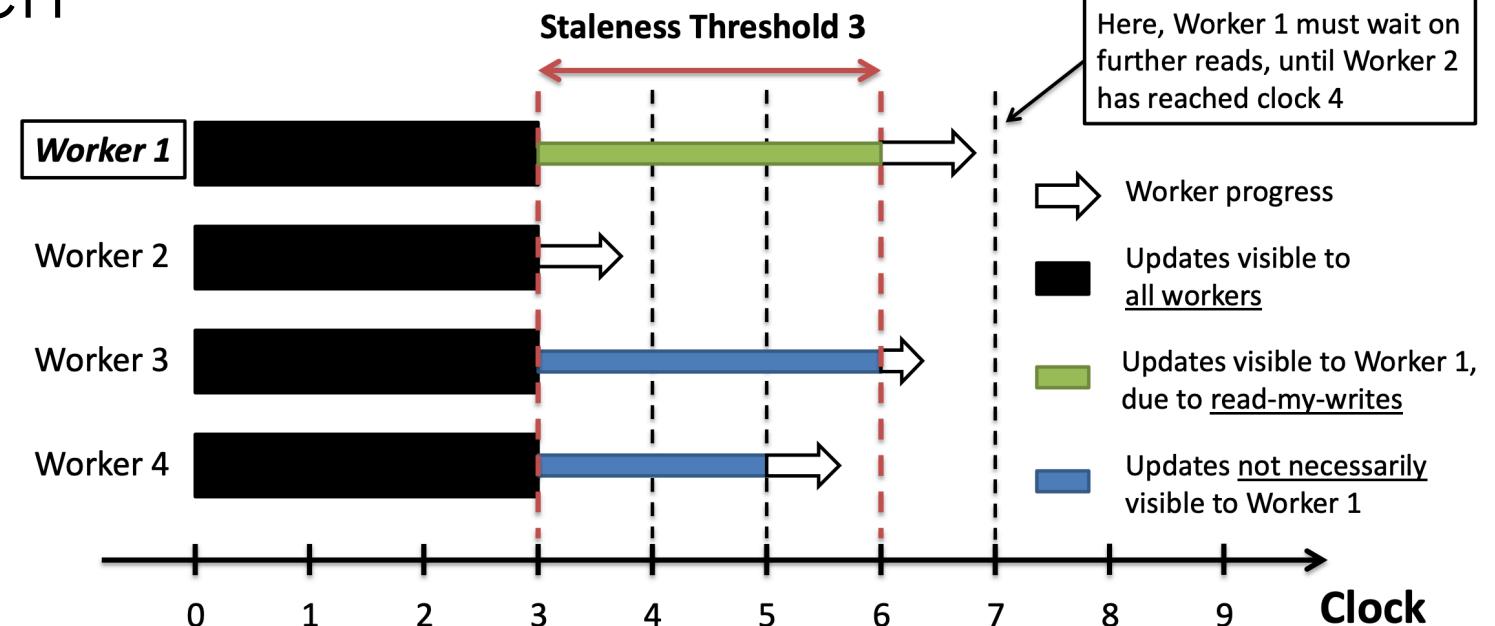
GPipe



Bounded Staleness

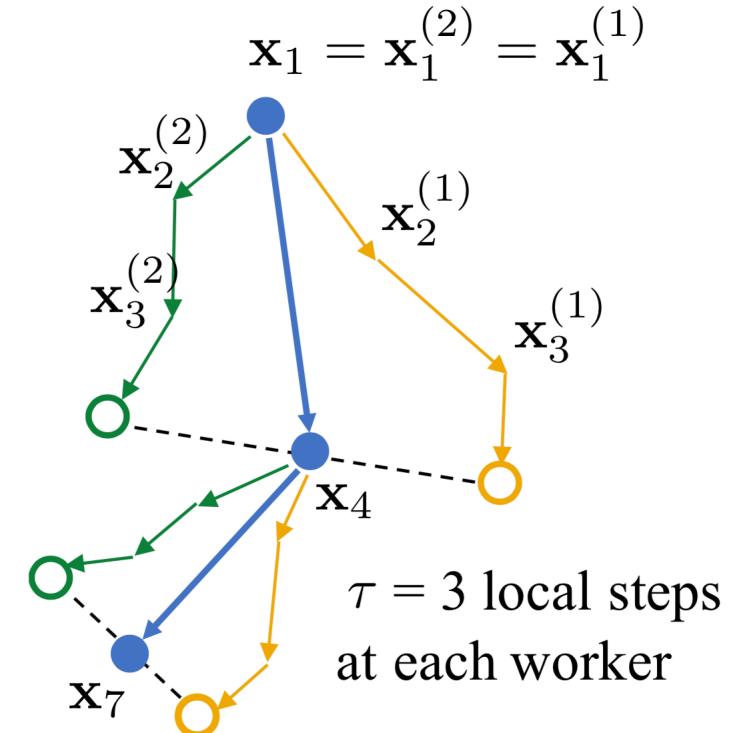
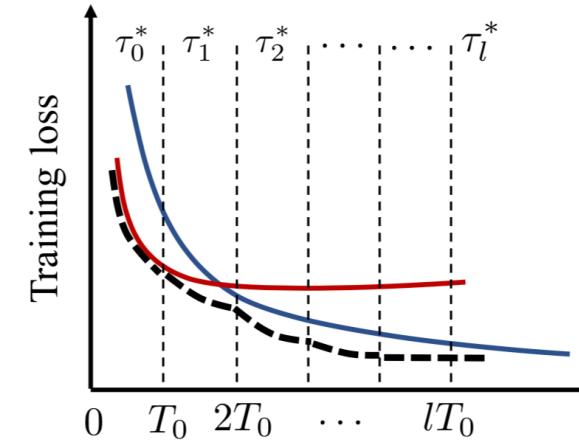
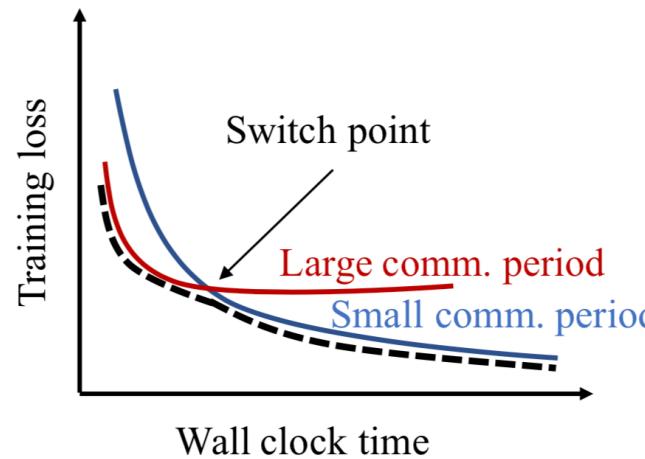
- Developed as part of the parameter server work at CMU
 - [More Effective Distributed ML via a Stale Synchronous Parallel Parameter Server \(NIPS'13\)](#)
- Compromise between Hogwild and BSP
- Unclear implications for deep learning
 - Non-convex loss

SSP: Bounded Staleness and Clocks



Adaptive Communication Strategies to Achieve the Best Error-Runtime Trade-off in Local-Update SGD (SysML'19)

- Studies Periodic Averaging SGD (PASGD)
- **Key Idea:** Change τ as algorithm converges



- More theoretical than previous reading
- Theoretical results do not make convex assumptions!

Old Stuff