# PipeDream: Generalized Pipeline Parallelism for DNN Training &

A Unified Architecture for Accelerating Distributed DNN Training in Heterogeneous GPU/CPU Clusters

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# PipeDream: Generalized Pipeline Parallelism for DNN Training

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SOSP 2019 - Symposium on Operating Systems Principles

EECS 598 – W21

#### Overview

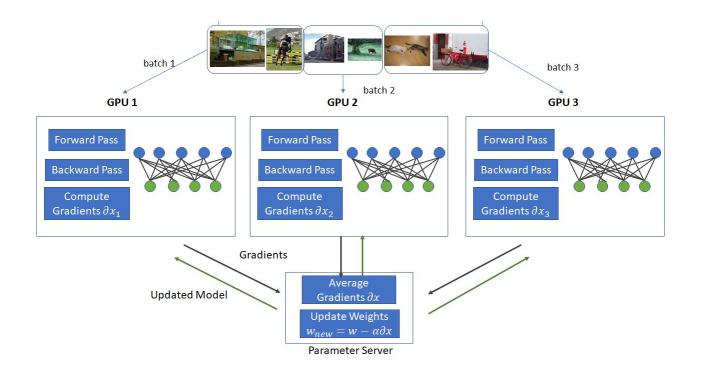
- Introduction
  - Data parallelism
  - Model parallelism
  - Pipelining (GPipe)
- PipeDream
  - Principles
  - Challenges
- Implementation and Evaluation

#### Distributed Training Approaches

- Intra-batch
  - Data parallelism (Data Partition)
  - Model parallelism (Model Partition)
  - Hybrid (e.g., For AlexNet, data partition in the convolutional layers)

- Inter-batch
  - Pipelining (GPipe)
- PipeDream: integrated intra-batch and inter-batch

#### Data parallelism



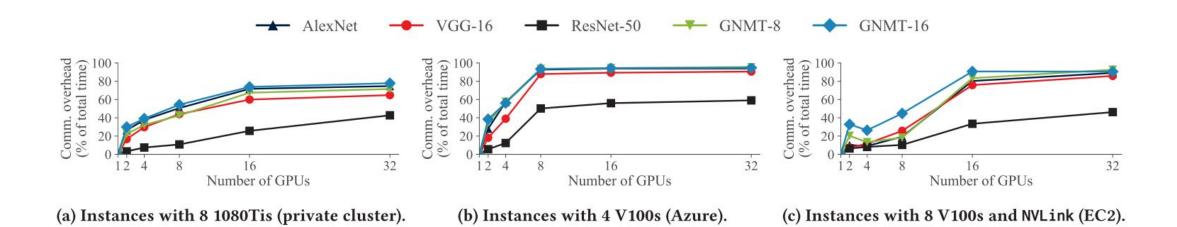
 Each worker will have the whole model but part of data

- Client-worker architecture (Parameter server)
- peer-to-peer architecture (ring all reduce algorithm)

#### Data parallelism

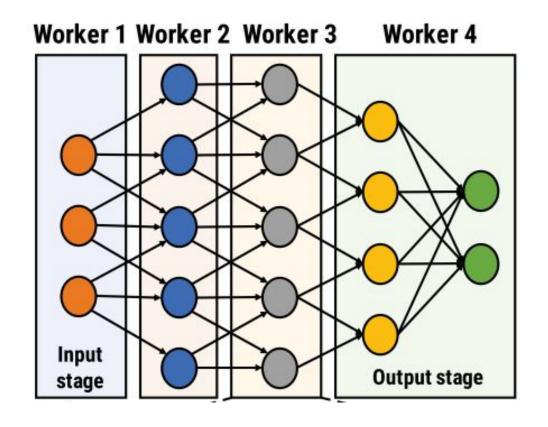
Significant communication overhead

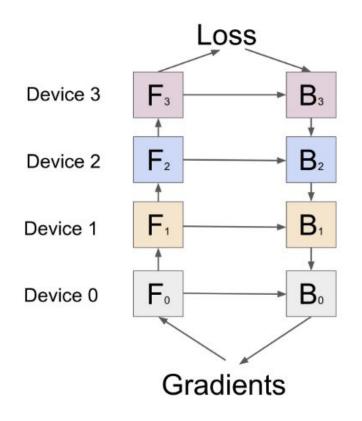
- Scales well for ResNet-50 (convolutional layers)
- · Does not scale well for dense connected model



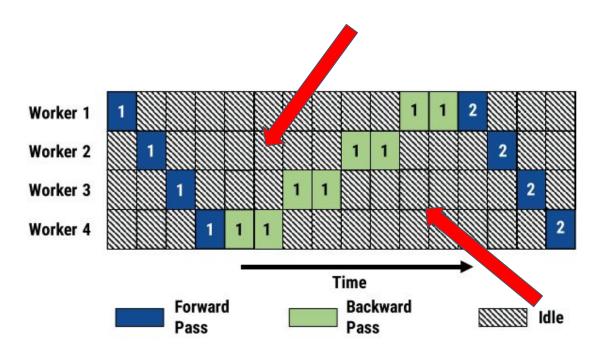
#### Model Parallelism

· Each worker will get the same data but have part of the model





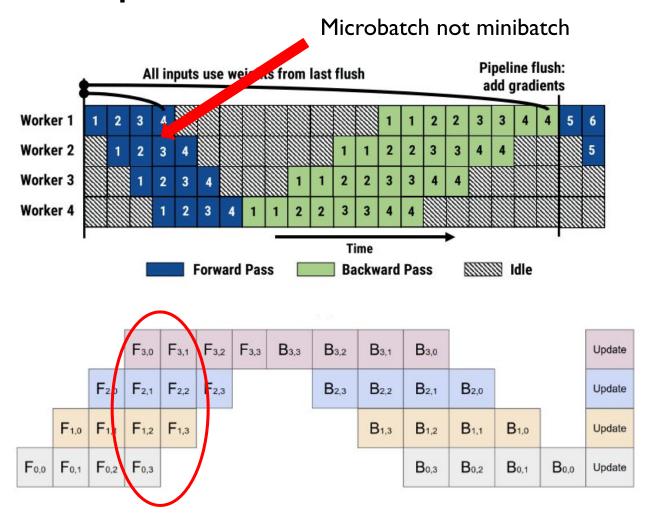
#### Model Parallelism



- The amount of data communicated is the size of intermediate outputs and and corresponding gradients
- Will not accelerate training
  - under-utilization of computing resources

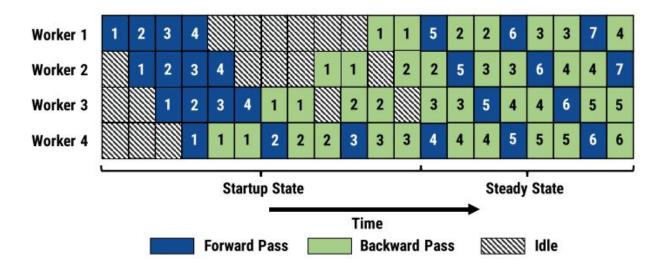
manually partition the mode

#### **GPipe**



- Divide minibatch into microbatch.
   Pipeline microbatch across workers
- Works can be used concurrently (improves the model parallelism)
- Frequent pipeline flush reduces the hardware efficiency
- Still has relatively large idle in the system
- Manually partition model

#### **PipeDream**



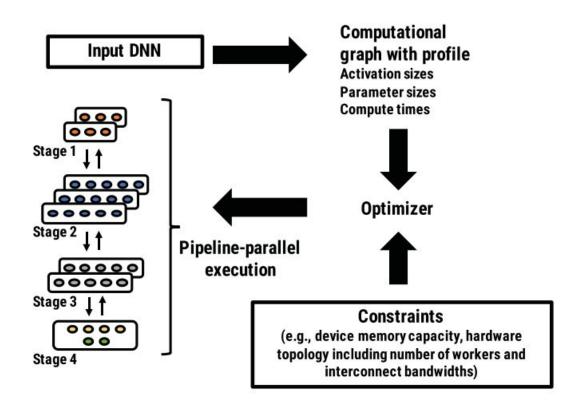
- In the startup phase: Inject multiple minibatchs to keep pipeline full (Similar to GPipe)
- In the steady state: each worker alters between forward and backwarpasses.
- Each worker once complete the forward pass for a minibatch, each stage will asynchronously sends the output to the next stage while starting to process another minibatch

#### **PipeDream**

No pipeline stalls in the steady state

- Pipelining communicates less. Each worker in PipeDream only need to communicate a subset of gradients and output activations, to only a single other worker. While in data parallelism, each worker need to send and receive the all parameters on the model.
- Pipelining overlaps computation and communication: Asynchronous communication of forward activations and backwards gradients across stages lead to the overlap of communication and computation

# PipeDream Challenge I:Work Partitioning



- PipeDream uses partitioning algorithm to automatically split the model
- Profiler:
  - Using a short (few minutes)
     profiling run 1000 minibatches on a single GPU
  - Record total computation time, the size of output activation (and input gradients) in bytes, and the size of weight parameters in bytes for each layer
  - Communication time is related to number of workers and parameter size in each layer

# PipeDream Challenge I:Work Partitioning

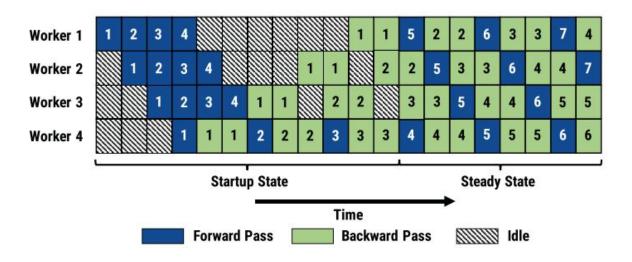


- Partitioning algorithm: (1) partition layers into stages, (2) the number of workers in each stage, and (3) optimal number of in-flight minibatches to full the pipeline
- The high level intuition is that the PipeDream's optimizer solves the dynamic programing problem and finds the optimal partitioning within a server and then split a model across servers
- $T^k(i \to j, m)$  is the total time taken by as single stage (layer i to j) replicated over m workers;  $A^k(i \to j, m)$  is the slowest stage (layer i to j) in the optimal pipeline

$$T^{k}(i \to j, m) = \frac{1}{m} \max \begin{cases} A^{k-1}(i \to j, m_{k-1}) \\ \frac{2(m-1)\sum_{l=i}^{j} |w_{l}|}{B_{k}} \end{cases} \qquad A^{k}(i \to j, m) = \min_{i \le s < j} \min_{1 \le m' < m} \max \begin{cases} A^{k}(i \to s, m - m') \\ \frac{2a_{s}/B_{k}}{T^{k}(s+1 \to j, m')} \end{cases}$$

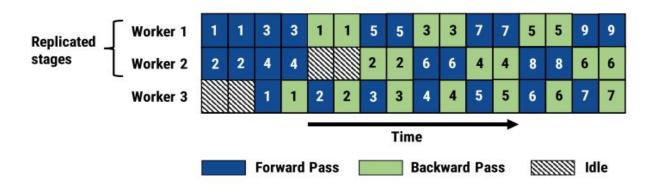
# PipeDream Challenge 2:Work Scheduling

- To determine whether the worker should compute the forward or backward.
- In the startup phase: need to decide how many minibatch to inject to keep the pipeline full in the steady stage. [#workers / #replicas in the input stage]
- In the steady phase: one-forward-one-backward (IFIB) schedule. To ensure each stage to produce the output at roughly the same rate



# PipeDream Challenge 2:Work Scheduling

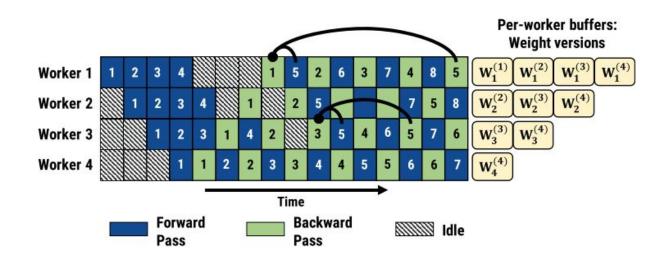
- When there is data parallelism in the stage, use the one-forward-one-backward round-robin (IFIB-RR).
- The figure below shows a 2-1 configuration. Worker 1 and worker 2 process different minibatch, while worker 3 process all minibatches.
- Worker I and 2 process replicated minibatches, worker 3 unreplicate.



# PipeDream Challenge 3: Effective Learning

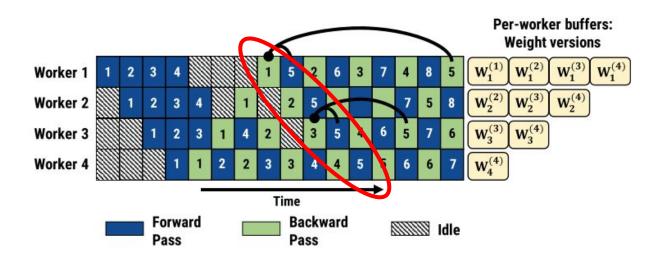
- Weight version mismatch challenge. The model parameters used for forward and backward should be the same for one minibatch
- Weight stashing: PipeDream will keep multiple weight versions one for each active minibatch.

• 
$$w^{(t+1)} = w^{(t)} - v \cdot \nabla f(w_1^{(t-n+1)}, w_2^{(t-n+2)}, \dots, w_n^{(t)})$$



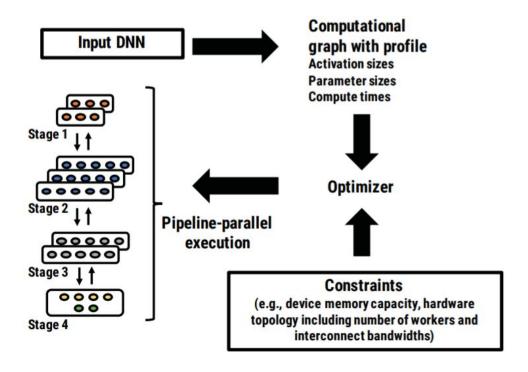
# PipeDream Challenge 3: Effective Learning

- Weight stashing does not guarantee the consistency of parameter versions for one minibatch across stages (Vertical Sync is an optional technique for this)
- Weight stashing does not significantly increase the per-worker memory

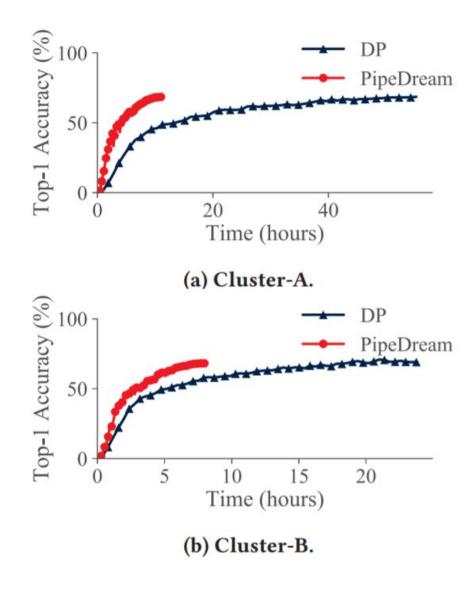


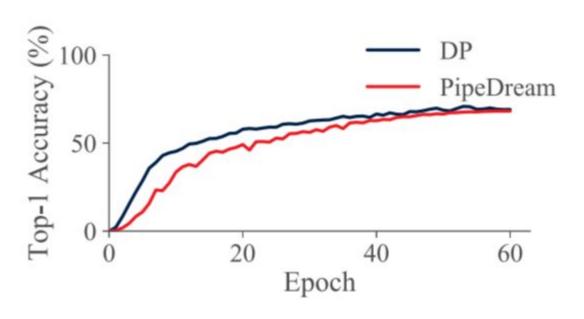
#### PipeDream Implementation

- A standalone Python library
- Uses PyTorch but can be integrated with other frameworks
- Some highlights:
  - Parameter State
  - Intermediate State
  - Stage Replication
  - Checkpointing



# PipeDream Evaluation: Image Classification





# PipeDream Evaluation: Other Tasks

Task	Model	Dataset	Accuracy Threshold	# Servers × # GPUs per server (Cluster)	PipeDream Config	Speedup over DP	
				, , ,		Epoch time	TTA
Translation	GNMT-16 [55]	WMT16 EN-De	21.8 BLEU	1x4 (A) 4x4 (A) 2x8 (B)	Straight Straight Straight	1.46× 2.34× 3.14×	2.2× 2.92× 3.14×
	GNMT-8 [55]	WMT16 EN-De	21.8 BLEU	1x4 (A) 3x4 (A) 2x8 (B)	Straight Straight 16		1.5× 2.95× 1×
Language Model	AWD LM [40]	Penn Treebank [41]	98 perplexity	1x4 (A)	Straight	4.25×	4.25×
Video Captioning	S2VT [54]	MSVD [11]	0.294 METEOR	4x1 (C)	2-1-1	3.01×	3.01×

#### PipeDream Evaluation: Comparison To GPipe

- Pipeline depth equal to PipeDream optimal:
  - 55% throughput slowdown on Cluster-A
  - 71% slowdown on Cluster-B
- Pipeline depth to max:
  - 35% slowdown on Cluster-A
  - 42% slowdown on Cluster-B
- Due to more frequent pipeline flushes

# Questions and Break

# A Unified Architecture for Accelerating Distributed DNN Training in Heterogeneous GPU/CPU Clusters

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OSDI'20 - the 14th USENIX Symposium on Operating Systems Design and Implementation

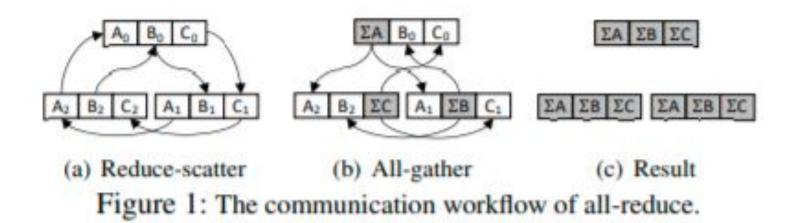
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#### Overview

- Recap
  - All-reduce
  - Parameter Server
- BytePS
  - Inter-machine communication
  - o Intra-machine communication
- Implementation and Evaluation

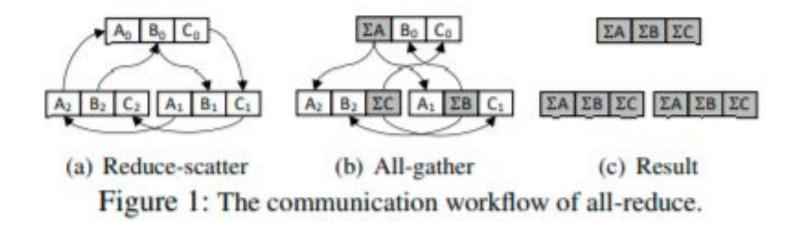
# Recap

#### All-reduce



- utilize a ring-based structure to enable parallel communication
- each node serves as a starter of a ring and an end of another ring

#### **Analysis**

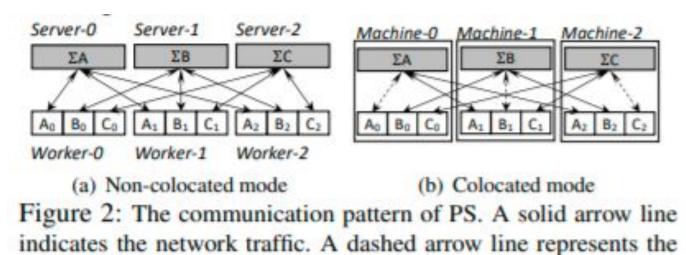


- Reduce-scatter: Each node sends (and receives)  $\frac{(n-1)M}{n}$  sytes to (and from) other nodes
- All-gather: Each node sends (and receives)  $\frac{(n-1)M}{n}$  sytes to (and from) other nodes
- Total traffic:  $\frac{2(n-1)M}{n}$  sytes. Total communication time:  $\frac{2(n-1)M}{nB}$

Notation: M: model size n: number of nodes B: network bandwidth

#### Continue recapping...

#### Parameter Server(PS)

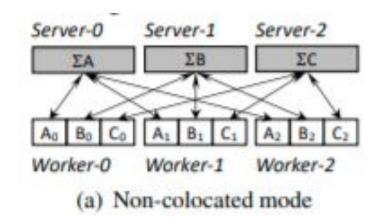


consists of two types of nodes: servers and workers

loop-back (local) traffic.

- workers running on GPU machines push gradients to servers
- servers aggregate gradients and update the parameters
- workers pull the latest parameters from servers for next iteration

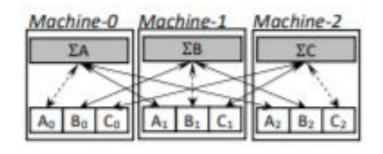
# **Analysis**



- (a) Non-colocated mode
  - server processes deployed on dedicated CPU machines
  - ullet model partitioned into k parts on k different CPU machines
- Each GPU worker sends and receives M bytes
- Each CPU server receives and sends  $\frac{nM}{k}$  bytes
- Communication time per iteration:  $\max(\frac{M}{B}, \frac{nM}{kB})$

Notation: M: model size n: number of nodes B: network bandwidth k: number of dedicated CPU machines

# **Analysis**



(b) Colocated mode

- (b) Colocated mode
- server processes deployed on GPU machines
- model partitioned into k parts on k different CPU machines
- Each node sends and receives M bytes
- Total traffic:  $\frac{2(n-1)M}{n}$  bytes. Total communication time:  $\frac{2(n-1)M}{nB}$

PS in colocated mode has the same communication time per iteration as all-reduce!

# However, both all-reduce and PS architecture have limitations...

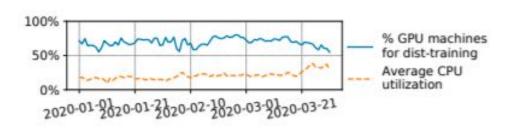


Figure 3: Daily statistics of our internal DNN training clusters from 2020-01-01 to 2020-03-31.

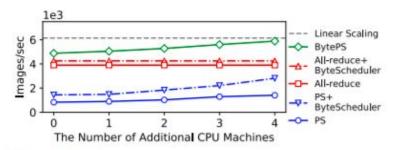
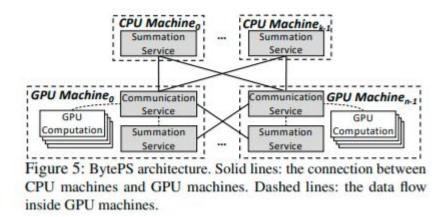


Figure 4: VGG-16 training performance of different architectures. We use 4 GPU machines with 32 GPUs in total. *Linear Scaling* represents the maximal performance (in theory) of using 32 GPUs.

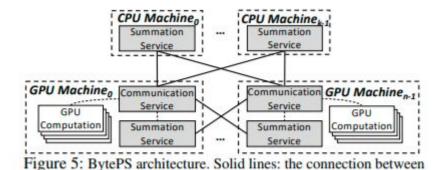
- Spare CPUs and bandwidth are not fully utilized on production GPU clusters
- Existing all-reduce and PS architectures not insufficient in performance
- This is why we need BytePS architecture, which mitigates the low utilization of these spare CPUs and bandwidth.

#### BytePS Architecture



- Two main modules: Communication Service(CS) and Summation Service(SS)
- SS runs on the CPUs of both CPU and GPU machines
- CS performs GPU computation on every GPU machine

# How to achieve optimal inter-machine communication for the BytePS architecture?



CPU machines and GPU machines. Dashed lines: the data flow

- An optimal communication requires wise SS workload assignment among CPU machines and GPU machines
- In the next slide, we use the following notation:

inside GPU machines.

Mss<sub>cpu</sub>: workload of SS on CPU machines

M<sub>SSGPU</sub>: workload of SS on GPU machines

#### **Analysis**

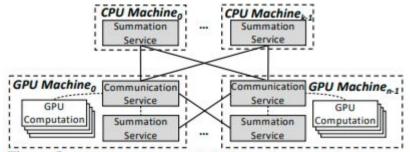


Figure 5: BytePS architecture. Solid lines: the connection between CPU machines and GPU machines. Dashed lines: the data flow inside GPU machines.

- a CS module sends and receives  $M M_{SS_{GPU}}$  bytes
- a SS module on a GPU machine receives and sends  $M_{SS_{GPU}}$  bytes from other n-1 GPU machines
- a GPU machine requires communication time  $t_g = \frac{M + (n-2)M_{SS_{GPU}}}{B}$
- a CPU machine requires communication time  $t_c = M_{SS_{CPU}}/B$
- the sum of all SS workloads should be M

#### **Analysis**

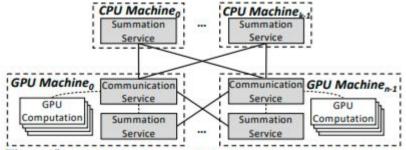


Figure 5: BytePS architecture. Solid lines: the connection between CPU machines and GPU machines. Dashed lines: the data flow inside GPU machines.

- GPU machine communication time  $t_0$  and CPU machine communication time  $t_0$  need to be equal to prevent one of them from being the bottleneck
- optimal workload assignment is given as

$$M_{SS_{CPU}} = \frac{2(n-1)}{n^2 + kn - 2k}M$$
  $M_{SS_{GPU}} = \frac{n-k}{n^2 + kn - 2k}M$ 

#### Intra-machine communication

#### PCIe-only topology

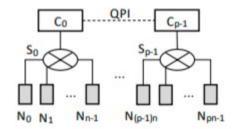


Figure 7: Notations of the PCIe-only topology.

- CPUs connected by QPI
- GPUs split into groups and connected to PCIe switches
- GPU-to-GPU memory copy across PCle switches is more of a bottleneck

#### CPU-assisted aggregation(PCle-only topology)

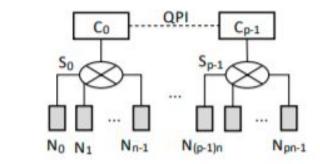
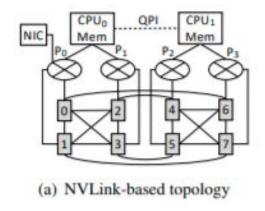


Figure 7: Notations of the PCIe-only topology.

Reduce-Scatter  $\implies$  GPU-CPU copy  $\implies$  CPU-reduce  $\implies$  Networking  $\implies$  CPU-GPU copy  $\implies$  All-Gather

 CPU-assisted aggregation is proved to be near-optimal in both theory and practice

#### **NVLink-based** topology



- GPUs split to different PCle switches
- All GPUs interconnected by NVLinks
- CPU-assisted aggregation no longer needed
- Asymmetric architecture causes competition of PCIe bandwidth

#### BytePS's solution of Summation Service(SS)

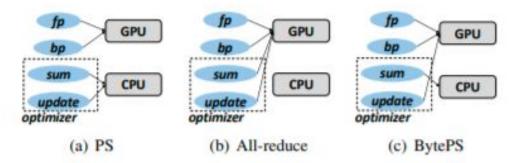


Figure 10: Component placement comparison between all-reduce, PS and BytePS.

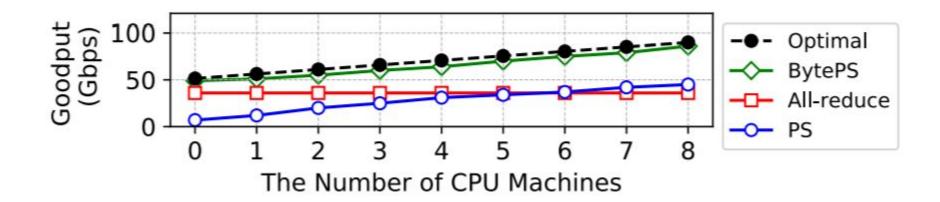
- CPU becomes a bottleneck as it cannot match increasing network bandwidth
- BytePS moves parameter update step to GPU machines

# BytePS: Implementation

- Multi-stage Pipeline
  - Tensor partition and pipelining
- RDMA Performance Optimization
  - Reduces RDMA WRITE operation to one round-trip by storing remote buffer address
  - Eliminates internal loopback on GPU machines
  - Uses page-aligned memory
- Provides native Python interfaces

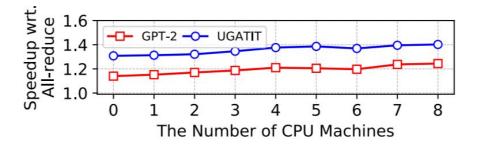
# BytePS: Evaluation

- Performance is close to optimal, better than all-reduce/PS

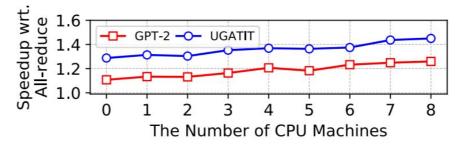


# BytePS: Evaluation

- Running more CPU machines achieves up to 20% speed increase with <<10% cost, since GPU is much more expensive



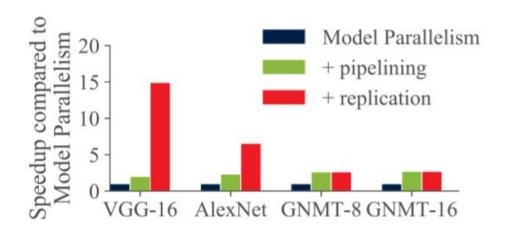
(a) PCIe-only GPU machines



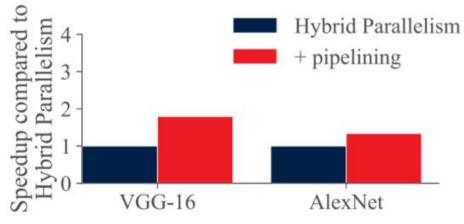
(b) NVLink-based GPU machines

# Thank you!

#### PipeDream Evaluation: Other Intra-batch Schemes







(b) Hybrid Parallelism.

(red is Pipedream)

# PipeDream Evaluation: Setup

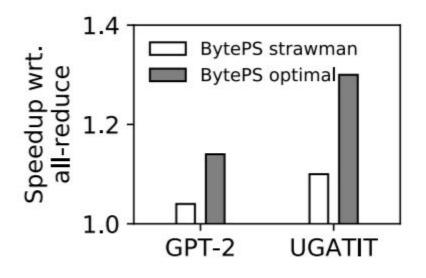
- Tasks:
  - Image Classification
  - Translation
  - Language Modeling
  - Video Captioning
- Clusters:

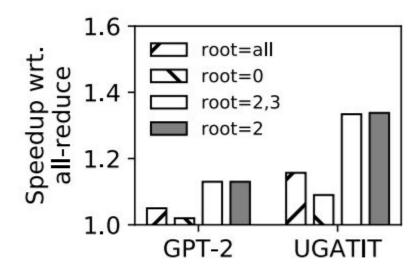
Cluster name	Server SKU	GPUs per server	Interconnects Intra-, Inter-server
Cluster-A	Azure NC24 v3	4x V100	PCIe, 10 Gbps
Cluster-B	AWS p3.16xlarge	8x V100	NVLink, 25 Gbps
Cluster-C	Private Cluster	1 Titan X	N/A, 40 Gbps

- Models:
  - VGG-16
  - ResNet-50
  - AlexNet
  - GNMT
  - AWD
  - S2VT
- Measure time trained to top-I accuracy

# BytePS: Evaluation

 Adaptation to intra-machine topology without optimization from CPU machines



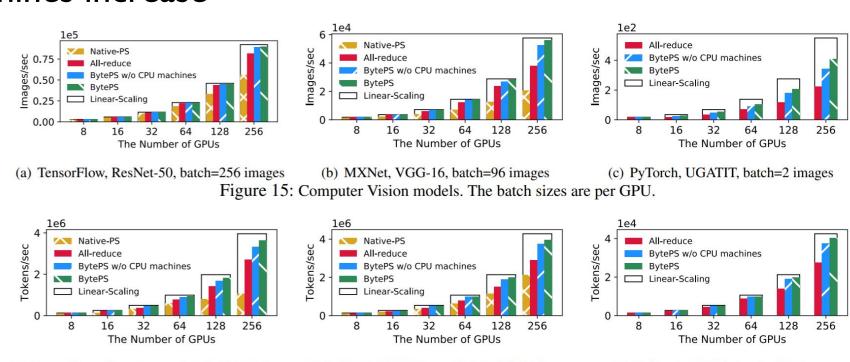


(a) PCIe-only GPU machines

(b) NVLink-based GPU machines

# BytePS: Evaluation

High scalability, speedup becomes larger as number of GPU machines increase



(a) TensorFlow, Transformer, batch=3072 tokens (b) MXNet, BERT-Large, batch=8192 tokens (c) PyTorch, GPT-2, batch=80 tokens Figure 16: NLP models. The batch sizes are per GPU.