Multi-Node Training



Bottlenecks of Single Node Jobs

- limited number of GPUs per node
- limited PCI bus bandwidth
- limited CPU power
- limited local storage



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Different Kinds of "Multi-Node Jobs"

Model Size:

- model fits into single GPU
- model needs multiple GPUs but fits on single node
- model needs multiple nodes

Model Replication:

- model is replicated across multiple GPUs on a single node
- model is replicated across multiple nodes

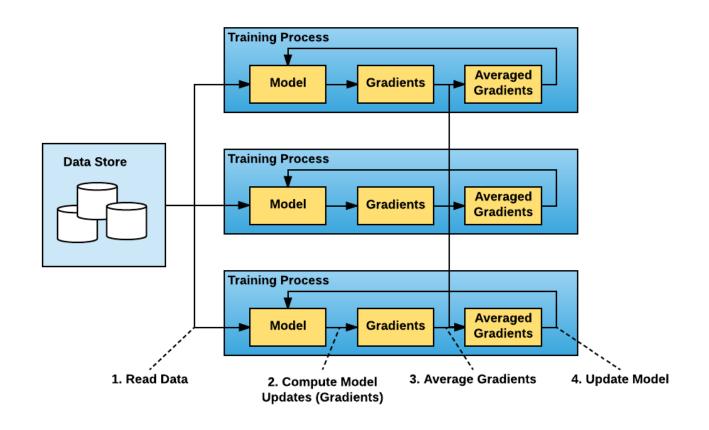


Multiple Communications Channels

- PCI bus for CPU, CPU
 ⇔GPU (computations involving single model)
- RDMA to CPU memory (data)
- GPUdirect for I/O to GPU, RDMA (data loading, inter-node parameters)
- NVLINK between GPUs (multi-GPU SGD within a node)
- standard Ethernet to CPU memory (control messages, maybe data)



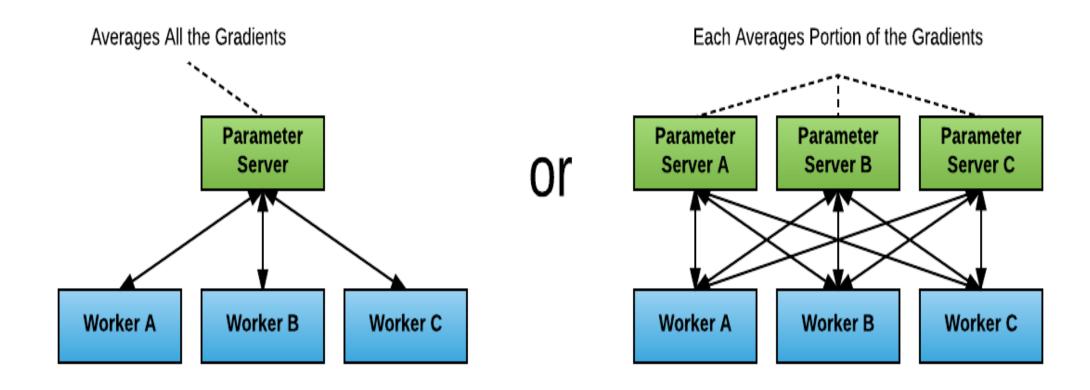
Common Multi-GPU / Multi-Node Training



(source: Horovod)



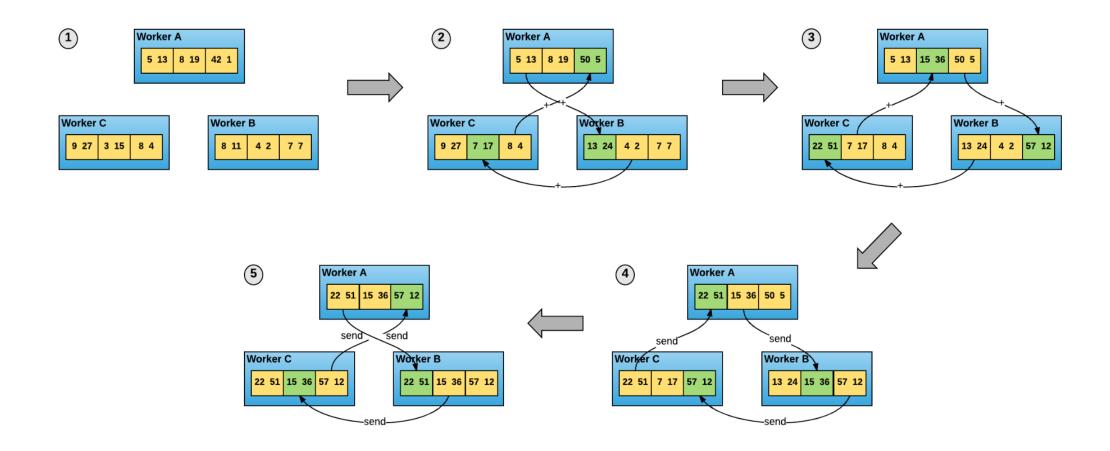
Parameter Servers



(source: Horovod)



Direct Parameter Exchanges



(source: Horovod)



PyTorch Support for Multi-GPU Training

- DataParallel, multiple I/O streams
- DistributedDataParallel, single node, multiple I/O streams
- DistributedDataParallel, multiple nodes, multiple I/O streams

These are wrappers around models and will take care of synchronizing gradients/weights across model instances.



How do you distribute?

- starting up jobs across clusters / machines
 - Ansible
 - Kubernetes
 - Slurm
- communications libraries
 - sockets, ZMQ
 - NCCL, Gloo
 - torch.distributed, Horovod
- all-in-one
 - o MPI



Kubernetes

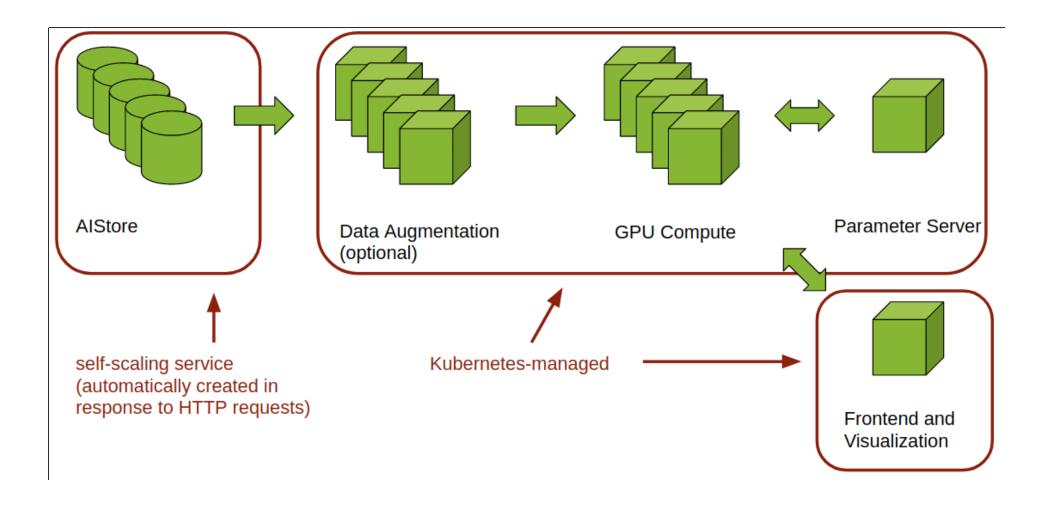
- start and stop processes on a cluster of machines
- encapsulate processes in containers
- provide a virtual network overlay
- provide access to hardware resources
- provide job and pod management



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Kubernetes for DL





Kubernetes Pod

- Kubernetes jobs/daemons/services are specified in YAML.
- They are applied with the kubectl apply command
- Basic unit is a Pod, a collection of Docker containers ("a pod of whales")

```
apiVersion: v1
  kind: Pod
  metadata:
  name: mypod
  spec:
  containers:
  - name: mypod
       image: gcr.io/research-191823/bigdata19
      command: ["nvidia-smi"]
       resources:
       limits:
           nvidia.com/gpu: "1"
  restartPolicy: Never
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```



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Kubernetes Setup with kubetpl

Job specs are complex and contain many repeats; templating helps:

GPU-based Job:

```
$ kubetpl job -G 1 -c nvidia-smi | kubectl apply -f -
```

CPU-based Pod:

\$ kubetpl pod -l tmbdev/redis-server -c redis-server | kubectl apply -f -



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Distributed Training

(notebook)

