

# Distributed Deep Learning

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# **Introduction**

# Distributed computing is scalability



In previous lectures, we mainly looked at single-node computation.

# Distributed computing is scalability

## User API

High level Packages

Programming API

Gradient Calculation (Differentiation API)

## System Components

Computational Graph Optimization and Execution

Runtime Parallel Scheduling

## Architecture

GPU Kernels, Optimizing Device Code

Accelerators and Hardwares

# Distributed Computing is Scalability

- Servers are stored in racks (42U rack, normally).
- Rack servers (typically 3U) are installed in these racks, each server normally has 4-8 GPU cards, interconnected through PCIe.
- Servers in the same rack normally has a faster network (ToR).
- Racks are also connected, but a lot slower in terms of point-to-point latency.

How do we map AI workloads into such systems?

# **AI Compute Parallelism**

# Canonical view of ML parallelism

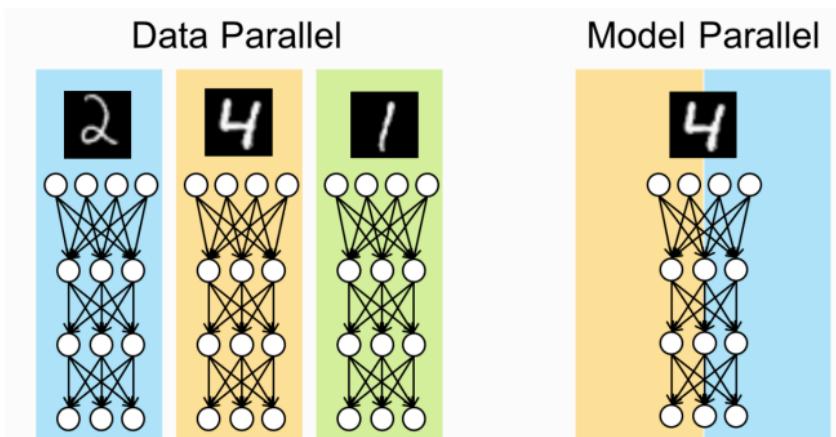
- **Data Parallelism:** data is partitioned across distributed workers, but the model is replicated.
- **Operator parallelism:** partition the computation of a specific operator, such as matmul, along non-batch axes, and compute each part of the operator in parallel across multiple devices.
- **Pipeline parallelism:** places different groups of ops from the model graph, referred as stages, on different workers

These parallelism can also be mixed.

We replicate the same model to various workers ( $N$  GPUs). The data is then split into  $N$  parts to be deployed on these  $N$  GPUs.

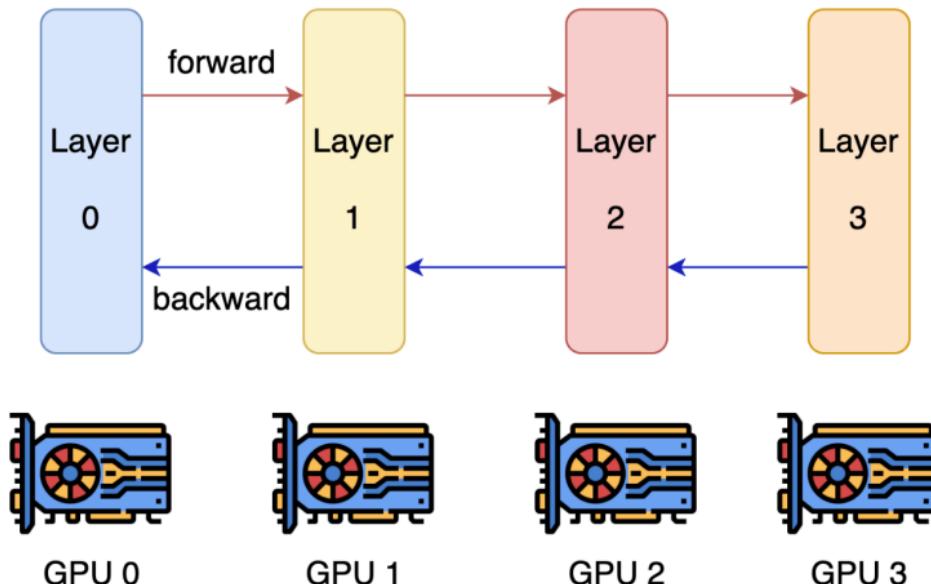
# Operator and model parallelism

- Different colors represent different GPU devices
- We can separate the model and put different portions of the model to different devices
- Some people say tensor parallelism – this is if you chop inside an operator, and operator/model parallelism normally refers to the fact that you are chopping at the operator granularity.



# Pipeline parallelism

- Very similar to CPU pipelining
- Works both in inference and training



# An alternative view of ML parallelism

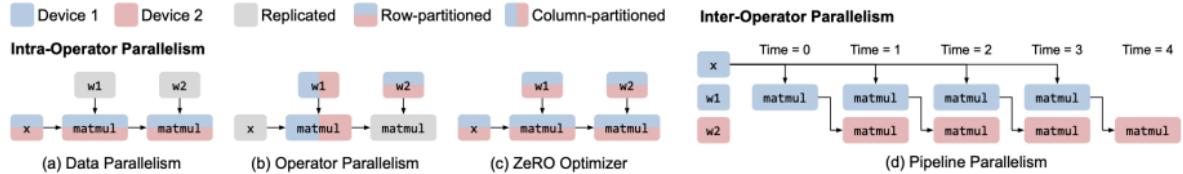
- **Intra-operator parallelism:**

An operator works on multi-dimensional tensors. We can partition the tensor along some dimensions, assign the resulting partitioned computations to multiple devices, and let them execute different portions of the operator at the same time.

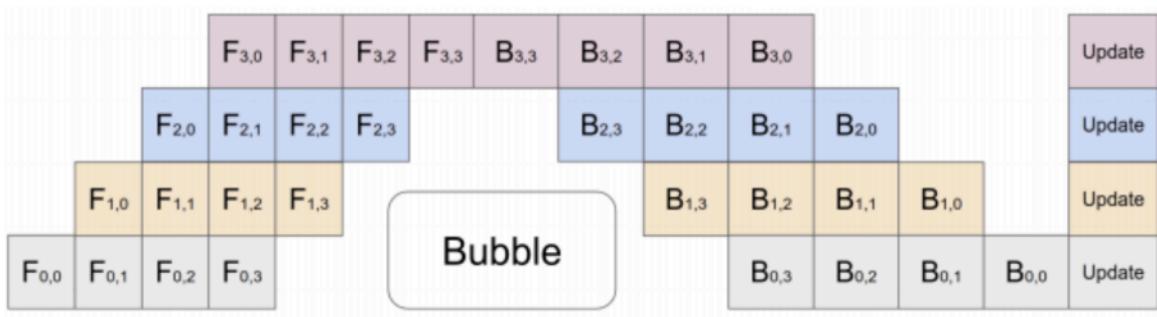
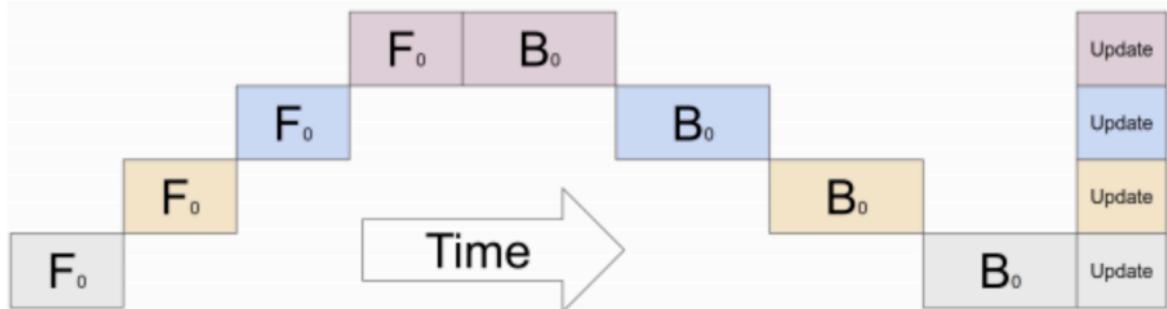
- **Inter-operator parallelism:**

We define inter-operator parallelism as the orthogonal class of approaches that do not perform operator partitioning, but instead, assign different operators of the graph to execute on distributed devices.

# ML parallelism



# ML Parallelism Model parallelism



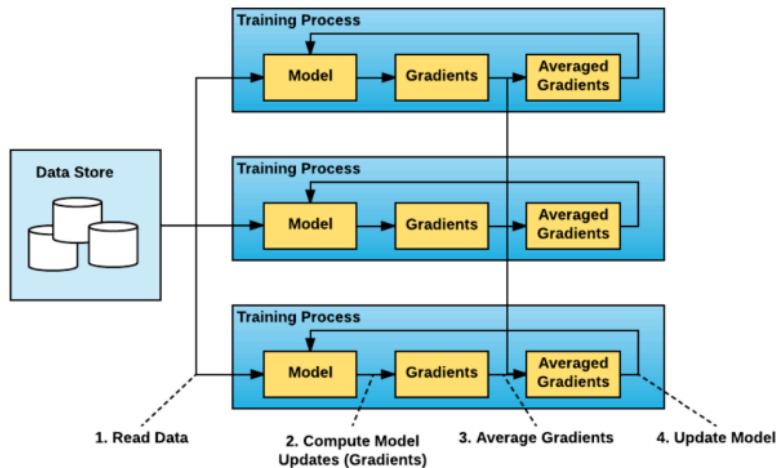
# **Distributed Training**

# Parameter server

It follows the data parallel approach.

In training, we split the data, and compute update at each local replica.

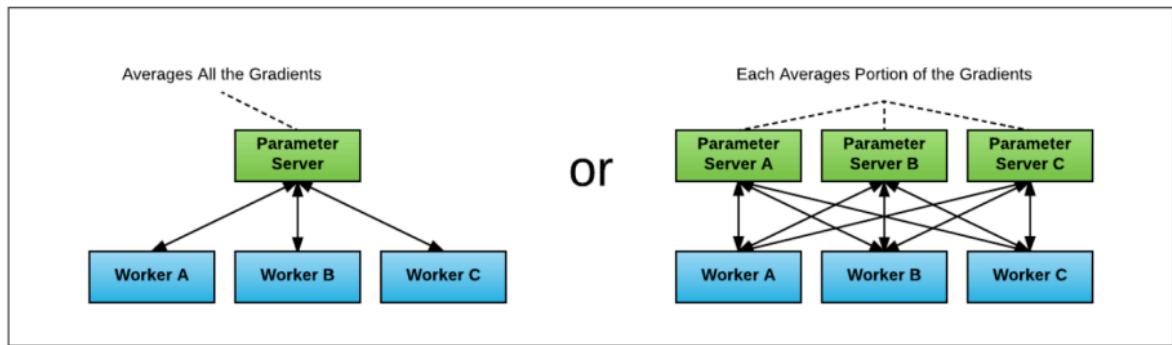
We then aggregate the gradients and propagate back the updated weights.



## Parameter server (ii)

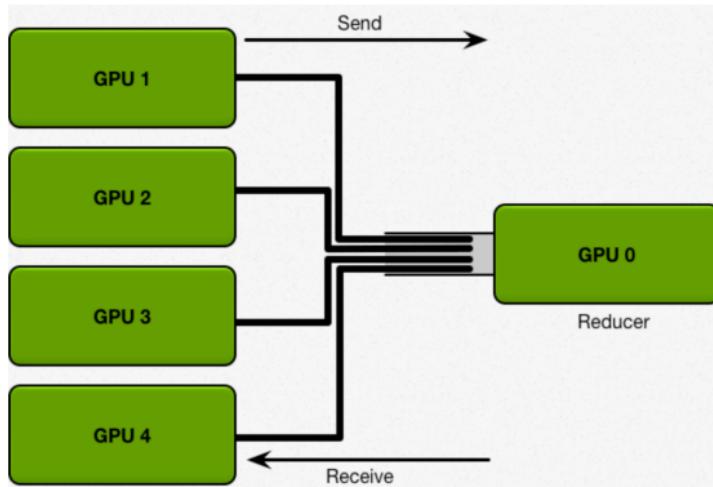
We need to sync the gradients and updated weights.

We gather the gradients in parameter server or parameter servers.



# The ring all reduce pattern

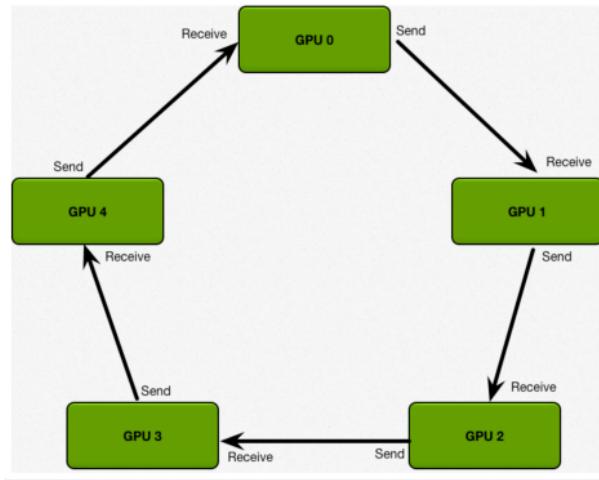
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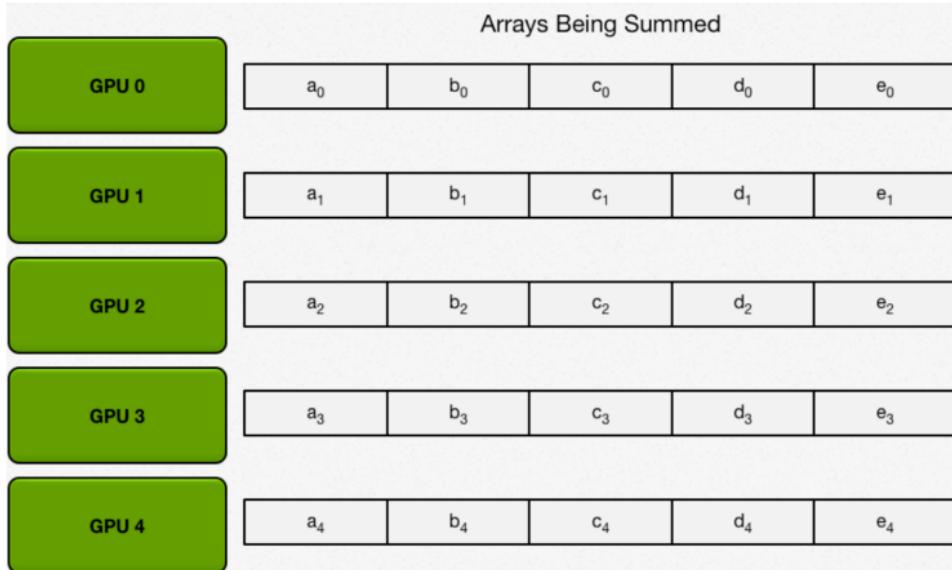
The most obvious communication pattern is to allow a crossbar connection, however, this is not very scalable.

# The ring all reduce pattern

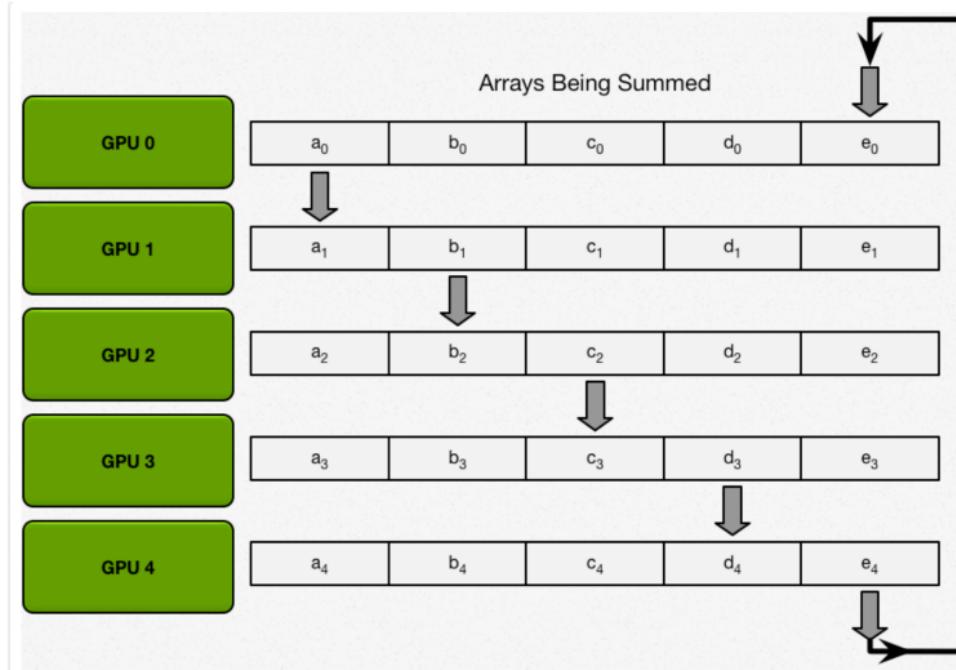
With the following physical connections



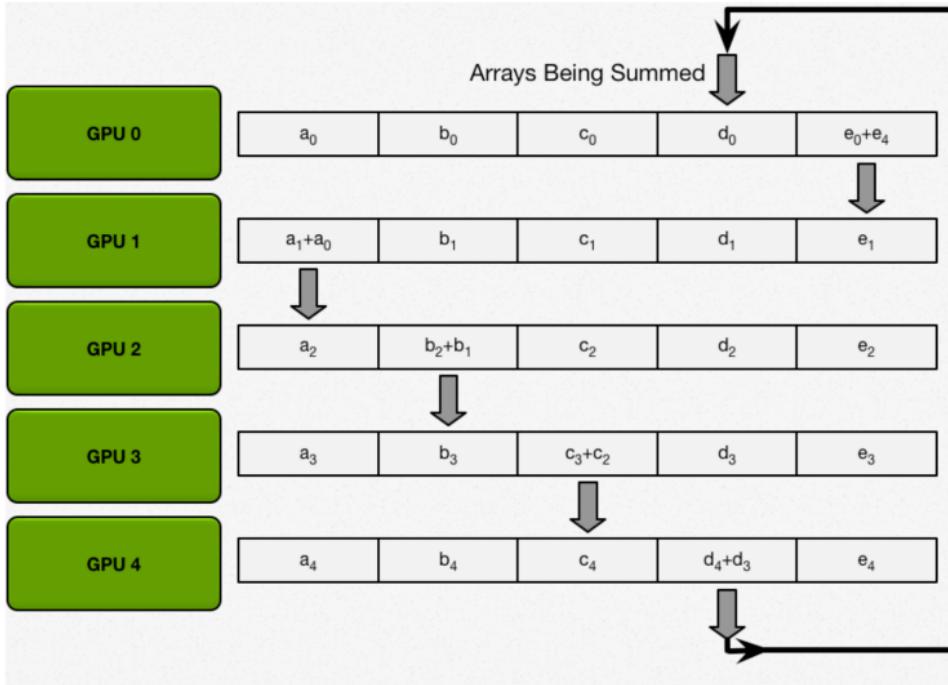
# The ring all reduce pattern



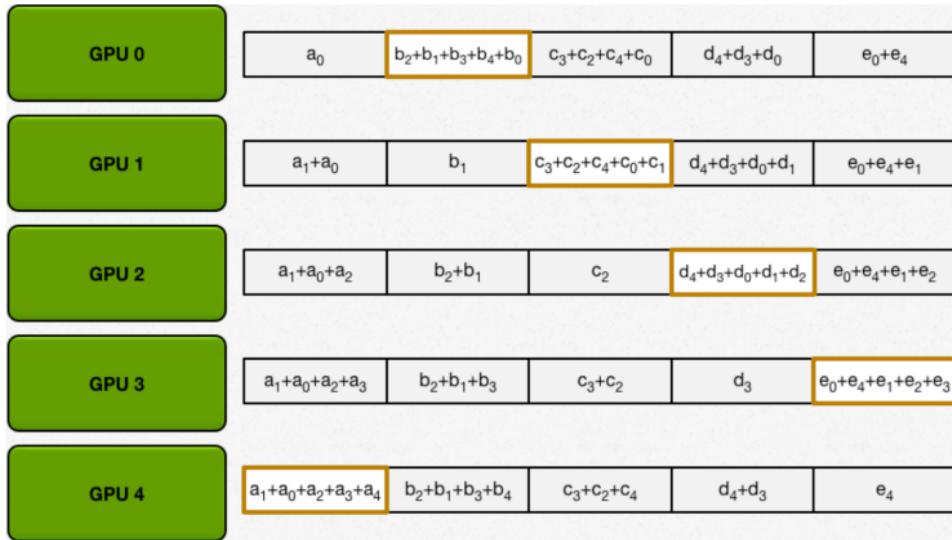
## The ring all reduce pattern (ii)



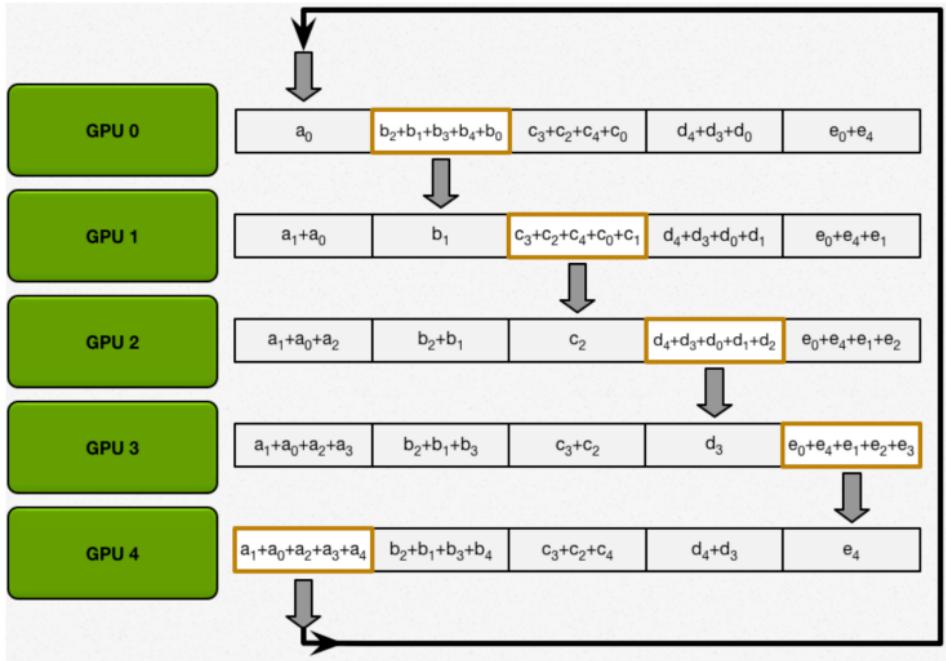
## The ring all reduce pattern (iii)



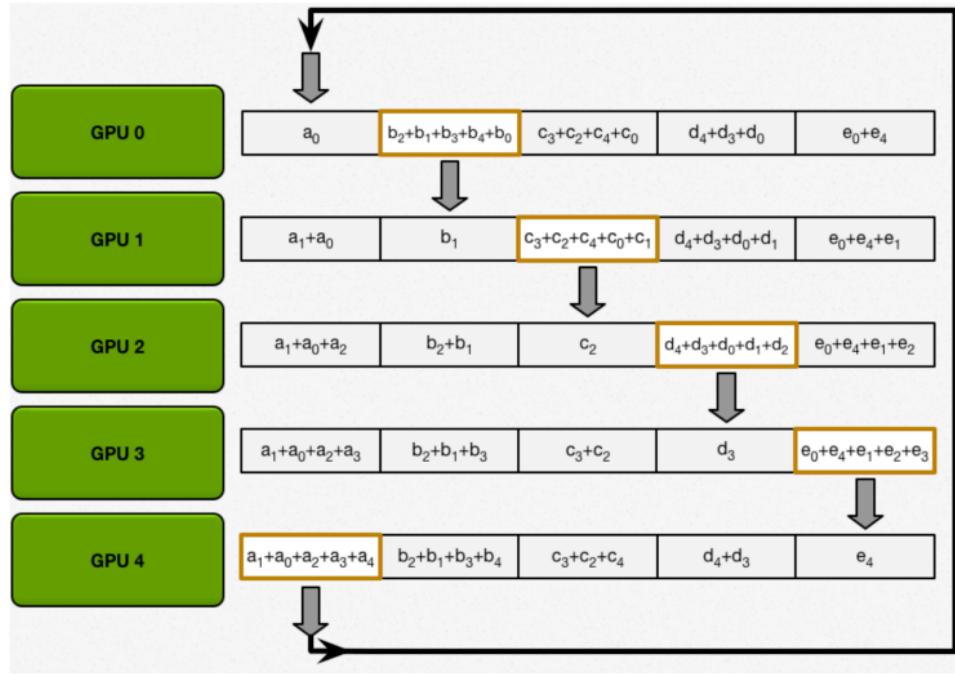
## The ring all reduce pattern (iv)



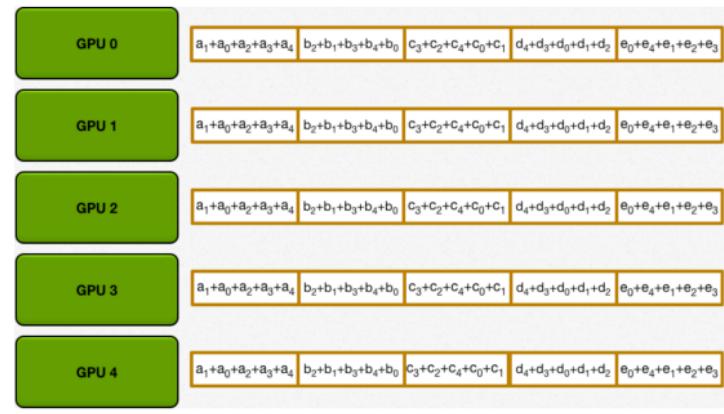
# The ring all reduce pattern (v)



# The ring all reduce pattern (vi)



## The ring all reduce pattern (vii)



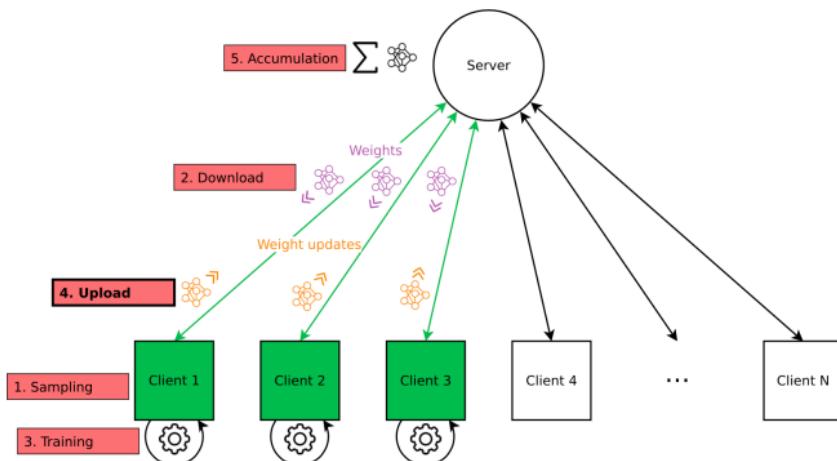
Normally you do not need to worry. Typically, you don't need to be concerned about the distributed training strategy, as PyTorch, NCCL, and MPI collectively manage most of it for you. However, you must consider the distributed training strategy if you are using more than 8 GPUs, which exceeds the capacity of a single node.

# **Federated Learning**

# The concept

Let's do not let the user data leave their phone (tricky!).

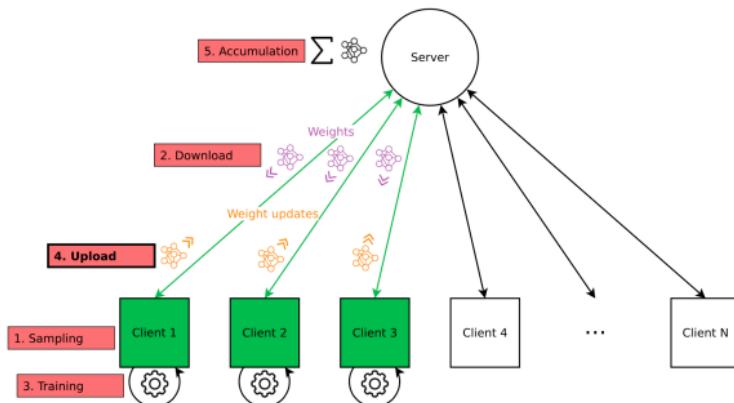
Phase 1 and 2: The server selects a subset of clients, with each algorithm employing a distinct sampling strategy. These clients update their local weights from the server.



## The concept (ii)

Let's do not let the user data leave their phone (tricky!).

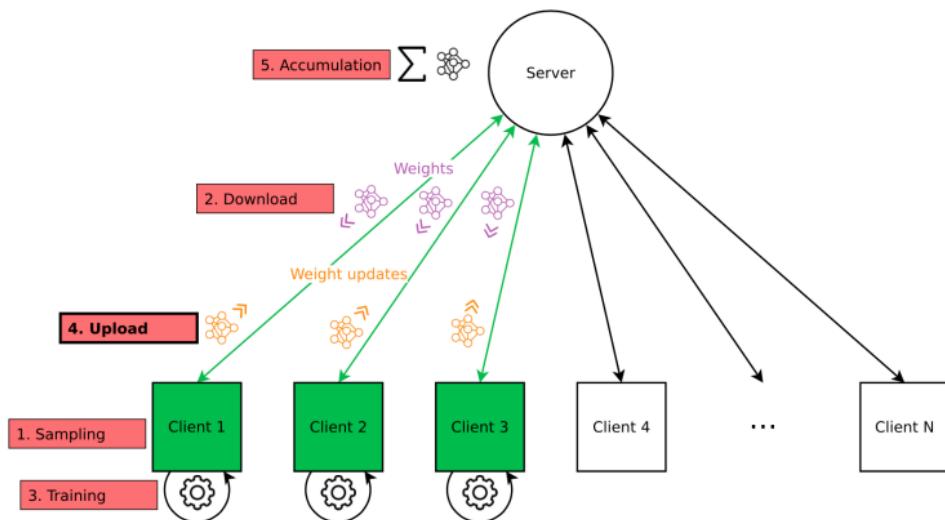
Phase 3: each client perform training on a model replica locally, using its own data. This training is always running concurrently, no matter a device is chosen or not. The server also normally gives a time constraint on this training, if a client did not finish its training in time, it is ignored (this is known as stragglers).



## The concept (iii)

Let's do not let the user data leave their phone (tricky!).

Phase 4 and 5: Clients now transfer back weight updates across  $E$  epochs back to the server. The server accumulate these updates.



# Pros and Cons

## **Biggest advantage**

No “data” leaves the user device!

## **Other advantages**

- Some level of privacy protections
- Give people an opportunity to do differential privacy

## **Disadvantages**

- Can take very long to train since effectively there is a subsampling.
- If server is not honest, bad things can happen.

# Summary

- Different parallelisms (eg. data, model, pipeline)
- Distributed Training, concepts on parameter servers and Ring All Reduce.
- Other learning paradigms: federated learning.