

## DEPARTMENT OF COMPUTER SCIENCE & ENGINEERING



## **Data Parallel Training**

**COMP4901Y** 

Binhang Yuan



# Data Parallel Paradigm

## Recall the Mini-Batch SGD



- Basic ideas:
  - To reduce the variance of stochastic gradients;
  - Split the training data into smaller batches;
  - Sampling batch (usually without replacement)
- Suppose we have:
  - *B* is the batch size:
  - *K* is the number of GPUs.

$$w_{t+1} = w_t - \alpha_t \cdot \frac{1}{B} \sum_{i=1}^{B} \nabla f(w_t; \xi_i)$$

• Distribute the computation by the data dimension.

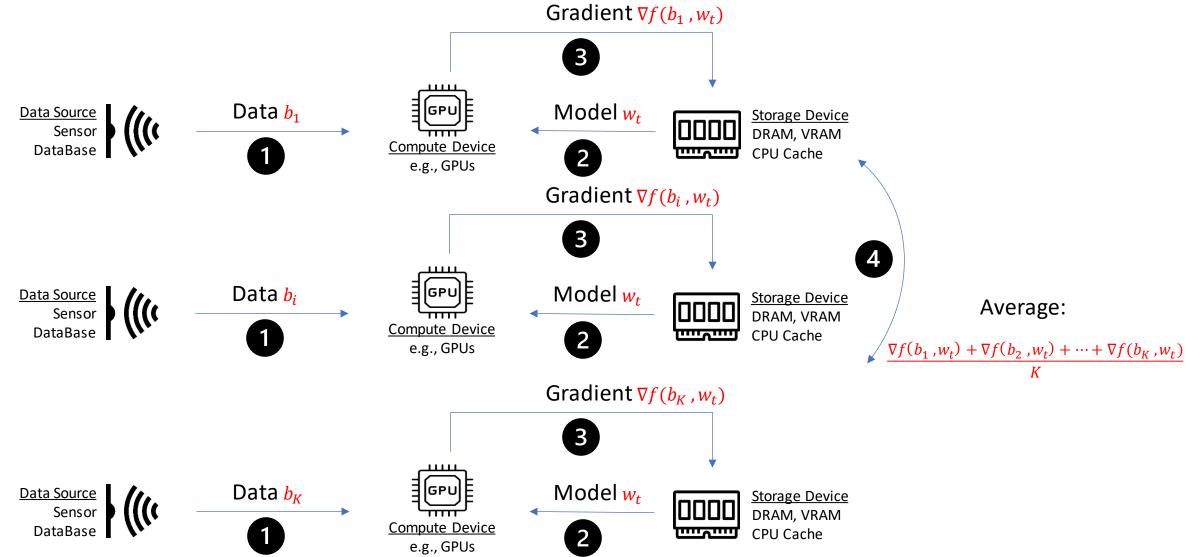
**Local Computation** 

$$w_{t+1} = w_t - \alpha_t \cdot \left(\frac{1}{B} \sum_{i_1=1}^{\frac{B}{K}} \nabla f(w_t; \xi_{i_1}) + \frac{1}{B} \sum_{i_2=1}^{\frac{B}{K}} \nabla f(w_t; \xi_{i_2}) + \dots + \frac{1}{B} \sum_{i_K=1}^{\frac{B}{K}} \nabla f(w_t; \xi_{i_K})\right)$$

$$\mathbf{GPU 1} \qquad \mathbf{GPU 2} \qquad \mathbf{GPU K}$$

### Data Parallel SGD



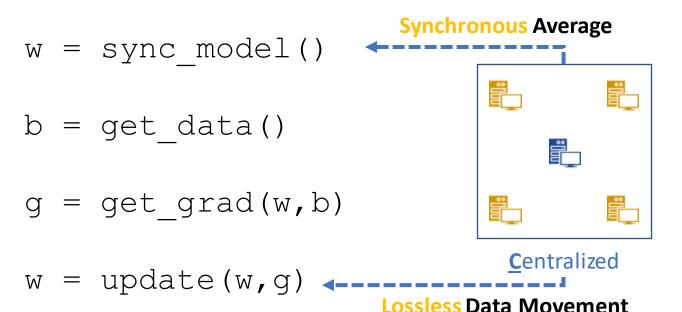


# **Basic Implementation**



### Core idea of data parallelism:

- Distribute batch gradient calculation to multiple workers;
- Synchronize gradients among GPU workers:
  - A central server (or AllReduce).
  - Synchronous average;
  - Lossless data movement.







- Suppose:
  - *C* is the total computation in the original forward and backward computation (counted by FLOPs);
  - D is the total number of parameters of the model;
  - *K* is the total number of GPUs in the cluster;
- Under Data parallel training:
  - What is the total computation on each device?  $\frac{C}{K}$
  - What is the communication volume? **ALLReduce** a *D* dimensional buffer.



# PyTorch-DDP Practice

## Overview



- PyTorch **DistributedDataParallel** (DDP) enables data parallel training in PyTorch.
- PyTorch **DistributedSampler** ensures each device gets a non-overlapping input batch.
- The model is replicated on all the devices.
- Each replica calculates gradients and synchronizes with the others using the **AllReduce** operation.

### **DDP API**

#### DISTRIBUTEDDATAPARALLEL

CLASS torch.nn.parallel.DistributedDataParallel(module, device\_ids=None, output\_device=None, dim=0, broadcast\_buffers=True, process\_group=None, bucket\_cap\_mb=25, find\_unused\_parameters=False, check\_reduction=False, gradient\_as\_bucket\_view=False, static\_graph=False, delay\_all\_reduce\_named\_params=None, param\_to\_hook\_all\_reduce=None, mixed\_precision=None, device\_mesh=None) [SOURCE]

Implement distributed data parallelism based on torch.distributed at module level.

This container provides data parallelism by synchronizing gradients across each model replica. The devices to synchronize across are specified by the input process\_group, which is the entire world by default. Note that DistributedDataParallel does not chunk or otherwise shard the input across participating GPUs; the user is responsible for defining how to do so, for example through the use of a DistributedSampler.

See also: Basics and Use nn.parallel.DistributedDataParallel instead of multiprocessing or nn.DataParallel. The same constraints on input as in torch.nn.DataParallel apply.

Creation of this class requires that torch.distributed to be already initialized, by calling torch.distributed.init\_process\_group().

DistributedDataParallel is proven to be significantly faster than torch.nn.DataParallel for single-node multi-GPU data parallel training.

To use DistributedDataParallel on a host with N GPUs, you should spawn up N processes, ensuring that each process exclusively works on a single GPU from 0 to N-1. This can be done by either setting CUDA\_VISIBLE\_DEVICES for every process or by calling:

```
>>> torch.cuda.set_device(i)
```

where i is from 0 to N-1. In each process, you should refer the following to construct this module:

```
>>> torch.distributed.init_process_group(
>>> backend='nccl', world_size=N, init_method='...'
>>> )
>>> model = DistributedDataParallel(model, device_ids=[i], output_device=i)
```







#### Initialize the Process Group

```
import os
import sys
import tempfile
import torch
import torch.distributed as dist
import torch.nn as nn
import torch.optim as optim
import torch.multiprocessing as mp
from torch.nn.parallel import DistributedDataParallel as DDP
def setup(rank, world size):
    os.environ['MASTER_ADDR'] = 'localhost'
    os.environ['MASTER_PORT'] = '12355'
    # initialize the process group
    dist.init process group("nccl", rank=rank, world size=world size)
def cleanup():
    dist.destroy process group()
```





### **Define the Toy Model**

```
class ToyModel(nn.Module):
    def __init__(self):
        super(ToyModel, self).__init__()
        self.net1 = nn.Linear(10, 10)
        self.relu = nn.ReLU()
        self.net2 = nn.Linear(10, 5)

def forward(self, x):
        return self.net2(self.relu(self.net1(x)))
```

### Use DDP API



#### Use DDP API

```
def demo_basic(rank, world_size):
    print(f"Running basic DDP example on rank {rank}.")
    setup(rank, world size)
    # create model and move it to GPU with id rank
    model = ToyModel().to(rank)
    ddp model = DDP(model, device ids=[rank])
    loss fn = nn.MSELoss()
    optimizer = optim.SGD(ddp model.parameters(), lr=0.001)
    optimizer.zero grad()
    outputs = ddp model(torch.randn(20, 10))
    labels = torch.randn(20, 5).to(rank)
    loss fn(outputs, labels).backward()
    optimizer.step()
    cleanup()
def run_demo(demo_fn, world_size):
    mp.spawn(demo_fn,
             args=(world_size,),
             nprocs=world size,
             join=True)
```





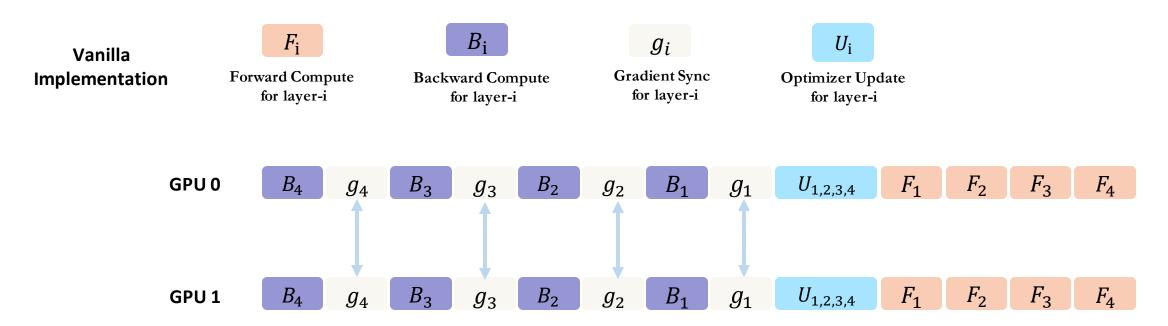
- Gradient Bucketing:
  - Collective communications are more efficient with large tensors.
  - Group gradient tensor to buckets on continuous memory (known as flatten.)
- Overlapping communication and computation during back-propagation:
  - The AllReduce operation on gradients can start before the local backward pass finishes.
  - With bucketing, DDP only needs to wait for all contents in the same bucket before launching communications.





### PyTorch System:

### Optimize the standard data-parallel computation:

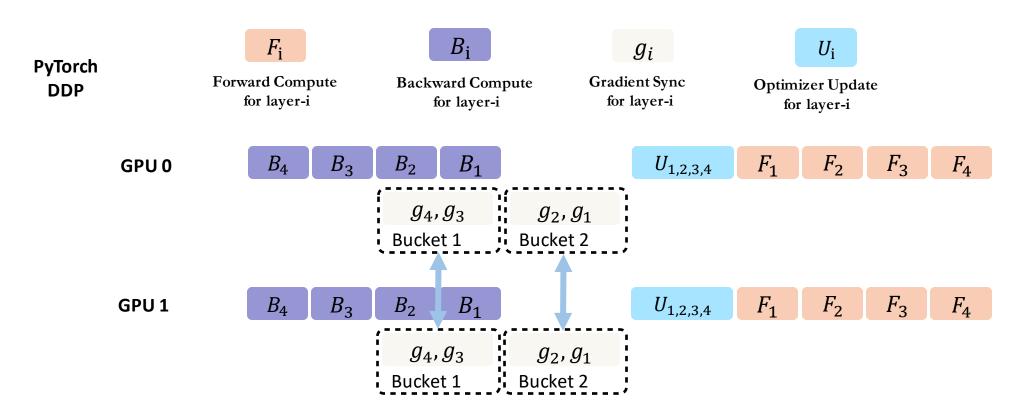






#### PyTorch System:

### Optimize the standard data-parallel computation:





# Algorithm Optimization

## Overview



- Three main categories of algorithm design:
  - Loss communication compression;
  - Asynchronous training:
  - Decentralized communication.

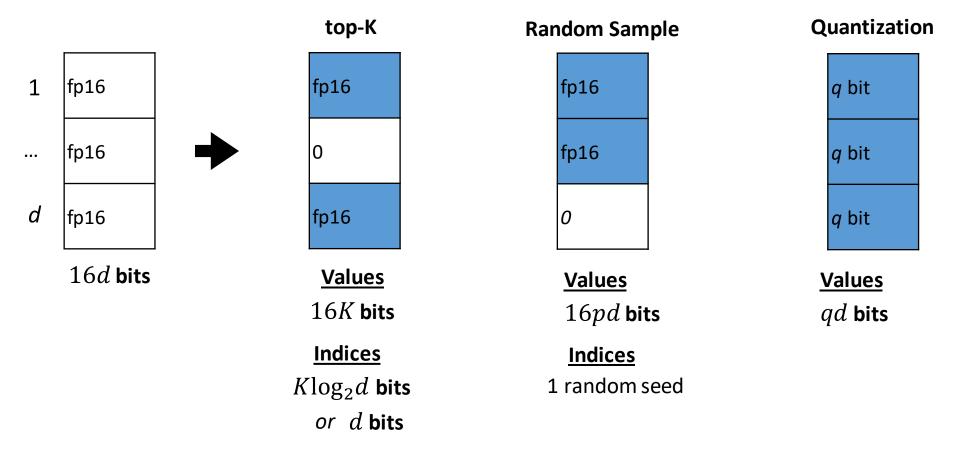




- Instead of exchanging the gradient as full precision floating point numbers;
- The system first compresses it into a lower precision representation before communication.
- Compression methods:
  - Top-K sparsification;
  - Random sparsification;
  - Quantization.

## **Compression Methods**





Expensive to compute and to encode Indices

Might not keep top values as in Top-K

Only provide up to 32x compression; hard to go aggressive





- Original paradigm: all workers conduct computation simultaneously and are all blocked until the communication among all machines is finished.
- If some machines are slower than other machines (i.e., there are stragglers), all machines need to wait for the slowest machine to finish the computation.
- Asynchronous training removes the synchronization barriers among all the machines with the consequence that each machine now has access to a staled model.

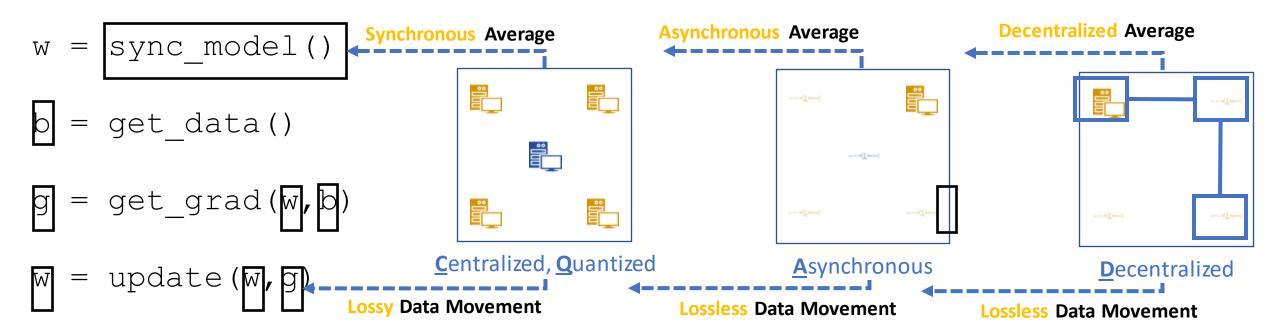
## **Decentralized Communication**



- Lossy communication compression is designed to alleviate system bottlenecks caused by network bandwidth.
- Another type of network bottleneck is caused by latency.
- In that case, when there are K workers, the ring -based AllReduce implementation has an  $\mathcal{O}(K)$  dependency on the network latency.
- Decentralized training:
  - Suppose a logical ring among *K* workers;
  - At every single iteration, a worker sends its model replica to the neighbour on its immediate left and neighbour on its immediate right;
  - Latency overhead becomes  $\mathcal{O}(1)$ .
  - On the downside, however, the information on a single worker only reaches its two adjacent neighbors in one round of communication.

# Algorithm Illustration.





Mathematical Formulation

$$w_{t+1} = w_t - \gamma \mathbf{C} \left( \sum_{i=1..n} \mathbf{C} (\nabla f_i(x_t, b_i)) \right)$$

$$w_{t+1} = w_t - \gamma \nabla f(x_{t-\tau_t}; b_i)$$
staleness caused by async

$$w_{t+1,i} = \frac{w_{t,i-1} + w_{t,i} + w_{t,i+1}}{3} - \gamma \nabla f(w_{t,i}; b_i)$$

### References



- https://dl.acm.org/doi/pdf/10.14778/3415478.3415530
- <a href="https://pytorch.org/tutorials/beginner/ddp\_series\_theory.html">https://pytorch.org/tutorials/beginner/ddp\_series\_theory.html</a>
- <a href="https://ethz.ch/content/dam/ethz/special-interest/infk/inst-cp/ds3lab-dam/documents/ZipML.pdf">https://ethz.ch/content/dam/ethz/special-interest/infk/inst-cp/ds3lab-dam/documents/ZipML.pdf</a>