## **Large Scale Training**

argonne-lcf / ai-science-training-series

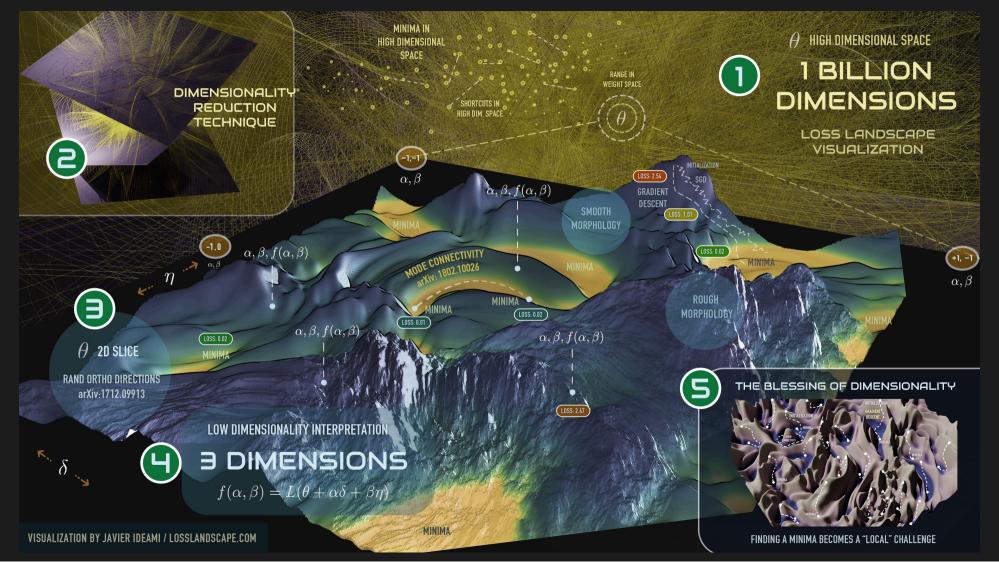
**Sam Foreman** 2022-11-01



## Why Distributed Training?

- Large batches may not fit in GPU memory
- Splitting data across workers → larger batch size
- Smooth loss landscape
- Improved gradient estimators
- Less iterations needed for same number of epochs
  - May need to train for more epochs if another change is not made
  - e.g. scaling learning rate
- See Large Batch Training of Convolutional Networks





## **Recent Progress**

Year	Author	Batch Size	Processor	DL Library	Time	Accuracy
2016	He et al. [1]	256	Tesla P100 x8	Caffe	29 Hrs	75.3%
	Goyal et al. [2]	8192	Tesla P100	Caffe 2	1 hour	76.3%
	Smith et al. [3]	8192 → 16,384	full TPU pod	TensorFlow	30 mins	76.1%
	Akiba et al. [4]	32,768	Tesla P100 x1024	Chainer	15 mins	74.9%
	Jia et al. [5]	65,536	Tesla P40 x2048	TensorFLow	6.6 mins	75.8%
	Ying et al. [6]	65,536	TPU v3 x1024	TensorFlow	1.8 mins	75.2%
	Mikami et al. [7]	55,296	Tesla V100 x3456	NNL	2.0 mins	75.29%
2019	Yamazaki et al	81,920	Tesla V100 x 2048	MXNet	1.2 mins	75.08%

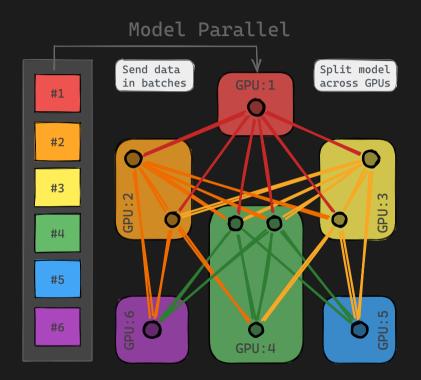


# **Model Parallel Training**



## **Model Parallel Training**

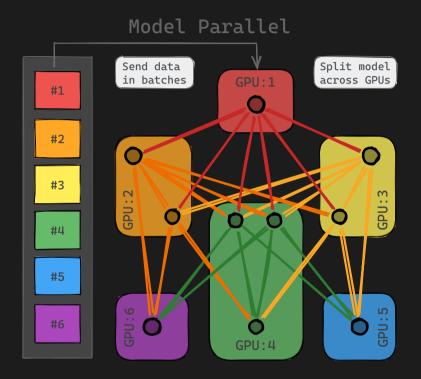
- Split up network over multiple workers
  - Each receives disjoint subset
  - All communication associated with subsets are distributed
- Communication whenever dataflow between two subsets
- Typically more complicated to implement than data parallel training
- Suitable when the model is too large to fit onto a single device (CPU / GPU)





## **Model Parallel Training**

- Suitable when the model is too large to fit onto a single device
  - Partitioning the model into different subsets is not an easy task
  - Might introduce load imbalancing issues limiting scale efficiency
- huggingface/transformers useful reference
  - Excellent series of posts in their documentation on Model Parallelism

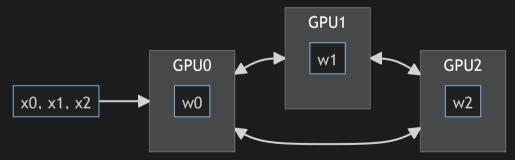




## **Model Parallel Training: Example**

$$y = w_0 * x_0 + w_1 * x_1 + w_2 * x_2$$

- 1. Compute  $y_0 = w_0 * x_0$  and send to o GPU1
- 2. Compute  $y_1 = y_0 + w_1 * x_1$  and send to ightarrow GPU2
- 3. Compute  $y=y_1*w_2*x_2$   ${f extstyle 
  abla}$

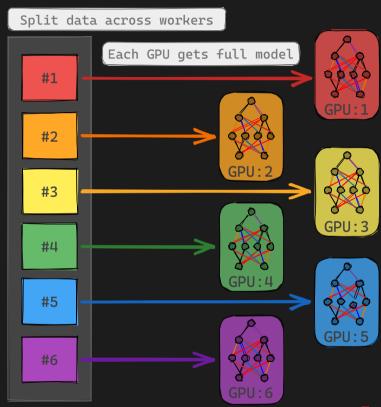






- Typically easier to implement
- Existing frameworks (Horovod, DeepSpeed, DDP, etc)
- Relatively simple to get up and running (minor modifications to code)<sup>1</sup>
- Recent presentation on data-parallel training available on YouTube

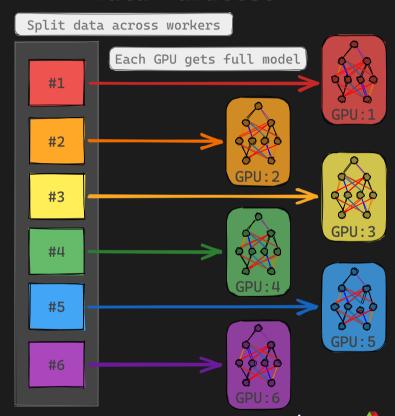
#### Data Parallel

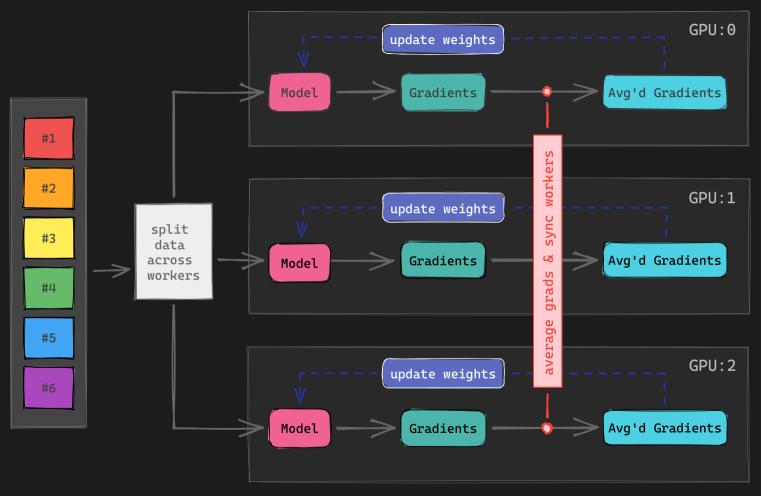




- Each worker has copy of complete model
- Global batch of data split into multiple minibatches
  - Each worker computes the corresponding loss and gradients from local data
- Before updating parameters, loss and gradients averaged across workers

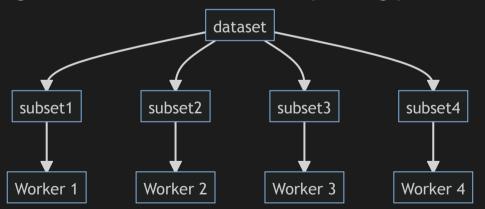
#### Data Parallel







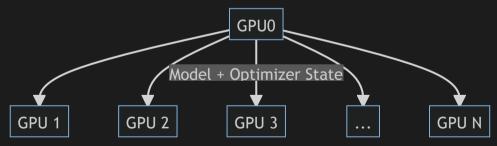
- Each worker has identical copy of model
- Global batch of data split across workers
- Loss + Grads averaged across workers before updating parameters





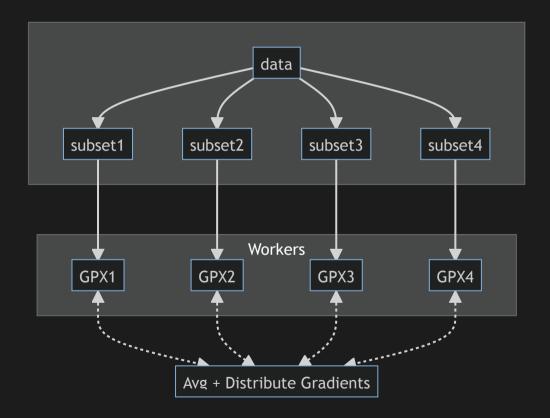
#### **Broadcast Initial State**

- At the start of training (or when loading from a checkpoint), we want all of our workers to be initialized consistently
  - Broadcast the model and optimizer states from hvd.rank() == 0 worker

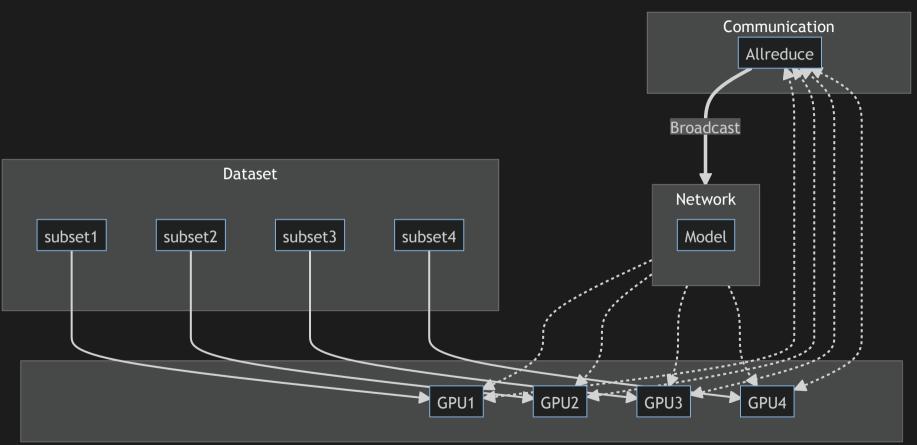




- Disjoint subsets of a neural network are assigned to different devices
- Each worker receives:
  - identical copy of model
  - unique subset of data









#### **Best Practices**

- Use parallel IO whenever possible
  - Feed each rank from different files
  - Use MPI IO to have each rank read its own batch from a file
  - Use several ranks to read data, MPI to scatter to remaining ranks
    - Most practical in big at-scale training

### **▲** Computation stalls during communication!

Keeping the communication to computation ratio small is important for effective scaling

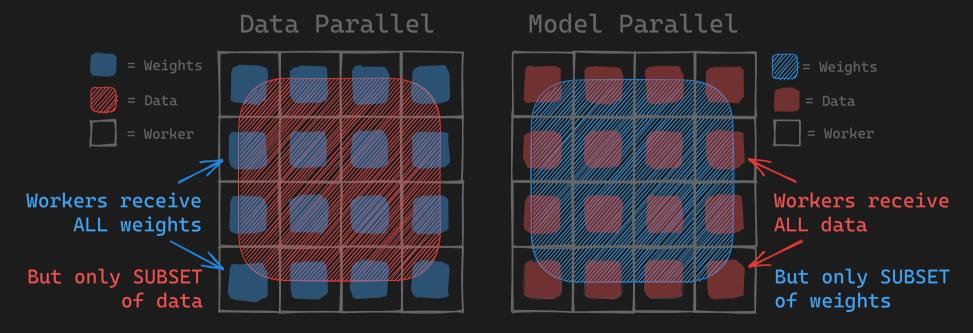


#### **Best Practices**

- Take advantage of data storage
  - Use striping on lustre
  - Use the right optimizations for Aurora, Polaris, etc.
- Preload data when possible
  - Offloading to a GPU frees CPU cycles for loading the next batch of data
    - minimize IO latency this way



## Comparison





### **Horovod: Overview**

- 1. Initialize Horovod<sup>1</sup>
- 2. Assign GPUs to each rank
- 3. Scale the initial learning rate by num. workers
- 4. Distribute gradients + broadcast state
  - Distribute gradients by wrapping tf.GradientTape with hvd.DistributedGradientTape
  - Ensure consistent initialization by broadcasting model weights and optimizer state from rank == 0 to other workers
- 5. Ensure workers are always receiving unique data
- 6. Take global averages when calculating loss, acc, etc. using hvd.allreduce  $(\ldots)$
- 7. Save checkpoints only from rank == 0 to prevent race conditions



#### TensorFlow + Horovod

Initialize Horovod:

```
import horovod.tensorflow as hvd
hvd.init()
```

Set one GPU per process ID (hvd.local\_rank())

```
gpus = tf.config.experimental.list_physical_devices('GPU')
for gpu in gpus:
    tf.config.experimental.set_memory_growth(gpu, True)
if gpus:
    local_rank = hvd.local_rank()
    tf.config.experimental.set_visible_devices(gpus[local_rank], 'GPU')
```



## **Scale the Learning Rate**

1. Scale by the number of workers to account for increased batch size

```
import horovod.tensorflow as hvd
optimizer = tf.optimizers.Adam(lr_init * hvd.size())
```



#### TensorFlow + Horovod

Training step then looks like:

```
@tf.function
def train step(data, model, loss fn, optimizer, first batch):
    batch, target = data
   with tf.GradientTape() as tape:
        output = model(batch, training=True)
        loss = loss fn(target, output)
tape = hvd.DistributedGradientTape(tape)
grads = tape.gradient(loss, model.trainable variables)
optimizer.apply gradients(zip(grads, model.trainable variables))
if first batch:
    hvd.broadcast variables(model.variables, root rank=0)
    hvd.broadcast variables(optimizer.variables, root rank=0)
return loss, output
```



#### **Deal with Data**

- At each training step, we want to ensure that each worker receives unique data
- This can be done in one of two ways:
  - 1. Manually partition data (ahead of time) and assign different sections to different workers
    - 1. Each worker can only see their local portion of the data
  - 2. From each worker, randomly select a mini-batch
    - 1. Each worker can see the full dataset

#### **▲** Don't forget your seed!

When randomly selecting, it is important that each worker uses different seeds to ensure they receive unique data



#### **Deal with Data**

```
(images, labels), (xtest, ytest) = tf.keras.datasets.mnist.load data(path='mnist.npz')
dataset = tf.data.Dataset.from tensor slices(
    (tf.cast(images[..., None] / 255.0, tf.float32),
    tf.cast(labels, tf.int64))
test dataset = tf.data.Dataset.from tensor slices(
    (tf.cast(xtest[..., None] / 255.0, tf.float32),
    tf.cast(ytest, tf.int64)))
nsamples = len(list(dataset))
ntest = len(list(test dataset))
dataset = dataset.repeat().shuffle(1000).batch(args.batch size)
test dataset = test dataset.shard(num shards=hvd.size(), index=hvd.rank()).repeat().batch(args.batch
```



## **Average Across Workers**

 Typically, we will want to take the global average of the loss across all our workers, for example:

```
global_loss = hvd.allreduce(loss, average=True)
global_acc = hvd.allreduce(acc, average=True)
```



## Thank you!

- Organizers
- ALCF Data Science & Operations
- Feel free to reach out!



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