

Parallel computing for Deep Learning

Goals: to understand...

- benefits of parallellization
- parallellization strategies
- parallel stochastic gradient descent (SGD)
- synchronous and asynchronous parallel SGD
- communication backends
- frameworks for distributed deep learning
- documentation of distributed DL frameworks



Parallel computing for Deep Learning

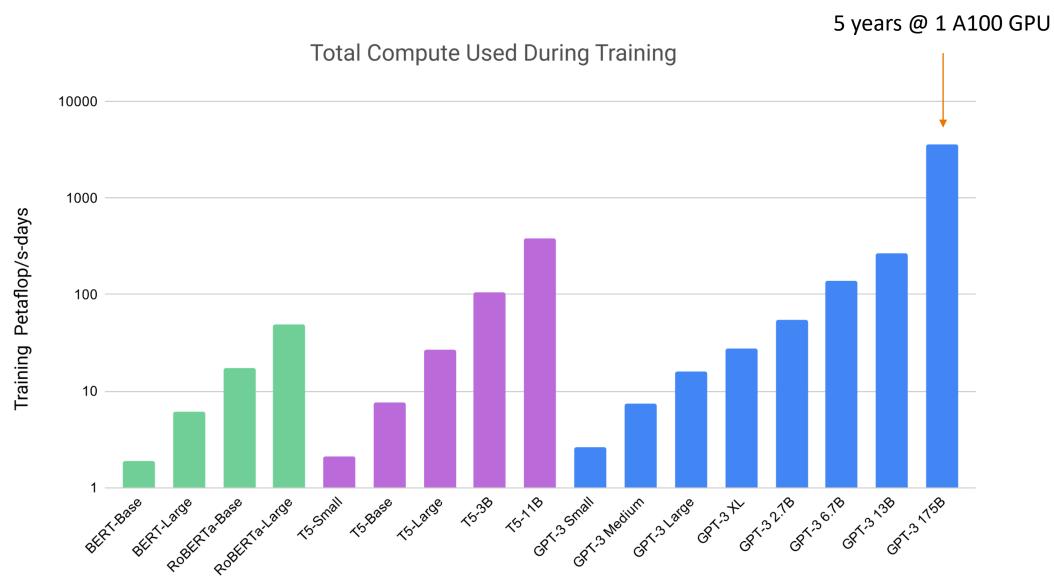
Goal: understand the documentation of distributed DL frameworks.

From TensorFlow docs on "distribution strategy":

- "tf.distribute.Strategy intends to cover a number of use cases along different axes... Synchronous vs asynchronous training: These are two common ways of distributing training with data parallelism. In sync training, all workers train over different slices of input data in sync, and aggregating gradients at each step. In async training, all workers are independently training over the input data and updating variables asynchronously. Typically sync training is supported via all-reduce and async through parameter server architecture."
- "MultiWorkerMirroredStrategy currently allows you to choose between two different implementations of <u>collective ops</u>. CollectiveCommunication.RING implements ring-based collectives using gRPC as the communication layer. <u>CollectiveCommunication.NCCL uses</u> <u>Nvidia's NCCL to implement collectives</u>."



Parallelization: why?





Parallelization: why?

Faster trainings ...

- Enables learning on larger datasets
- Enables improved accuracy through better hyperparemeter tuning
- Enables larger, more complex models

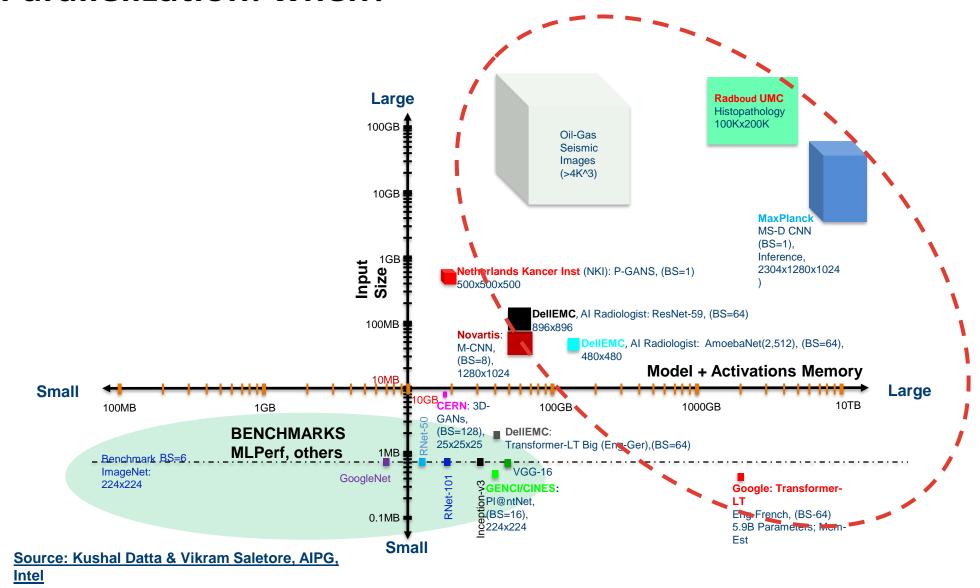
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Bigger models (high memory requirement) ...

Enables larger, more complex models



Parallelization: when?



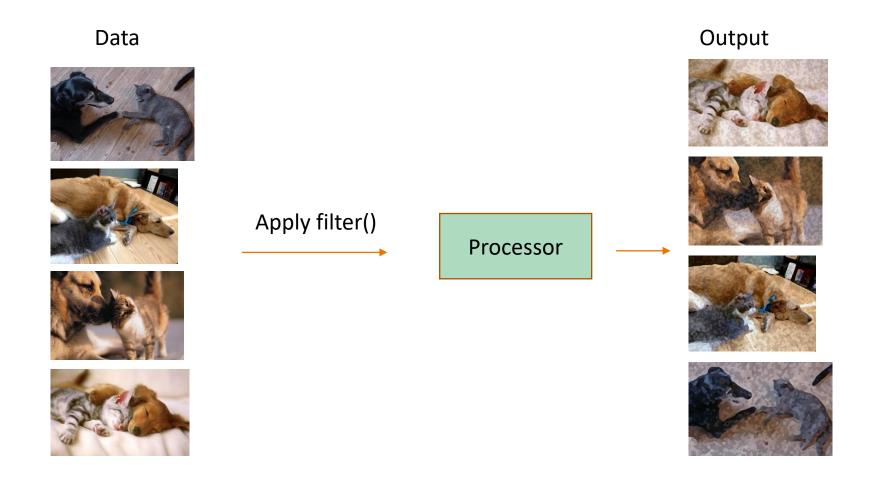
SURF

What is parallel computing?

Multiple processors or computers working on a single computational problem

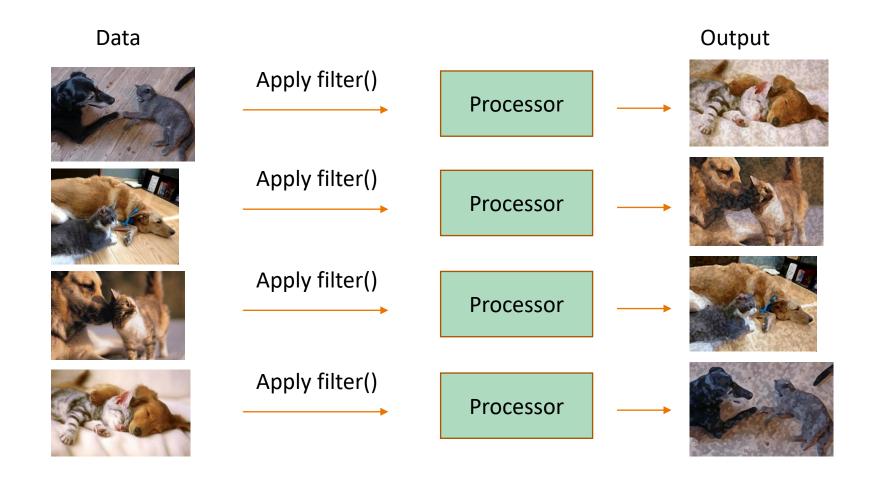


Serial computing





Parallel computing





Benefits:

- Solve computationally intensive problems (speedup)
- Solve problems that don't fit a single memory (multiple computers)

Requirements:

Problem should be divisible in smaller tasks



Parallelization for deep learning

How does one 'divide' DL tasks?

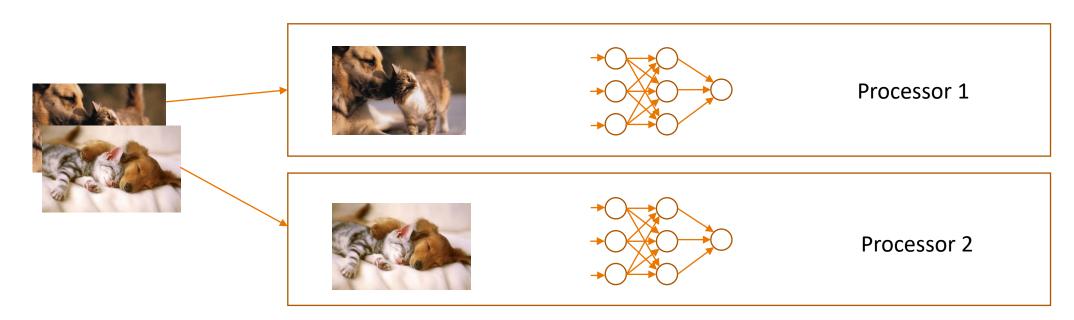
- Data parallelism
- Model parallelism
- Hybrid data/model parallelism
- Pipeline parallelism

Increasing complexity



Data Parallelism

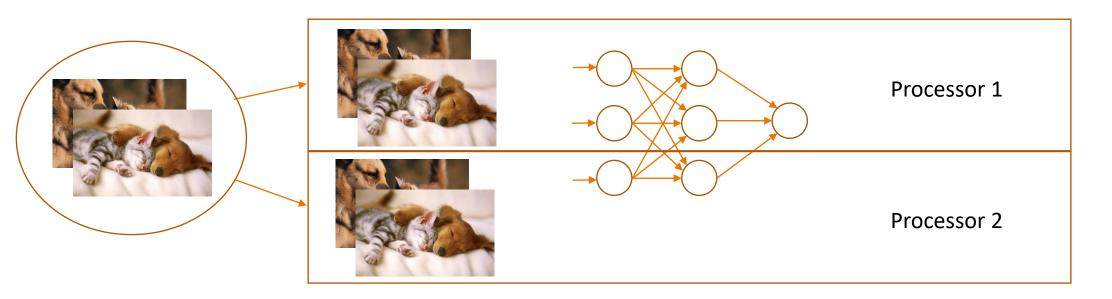
- Split the batch over multiple processors (CPUs/GPUs)
- Each processor holds a copy of the model
- Forward pass: calculated by each of the workers
- Backward pass: gradients computed (per worker), communicated and aggregated





Model parallelism

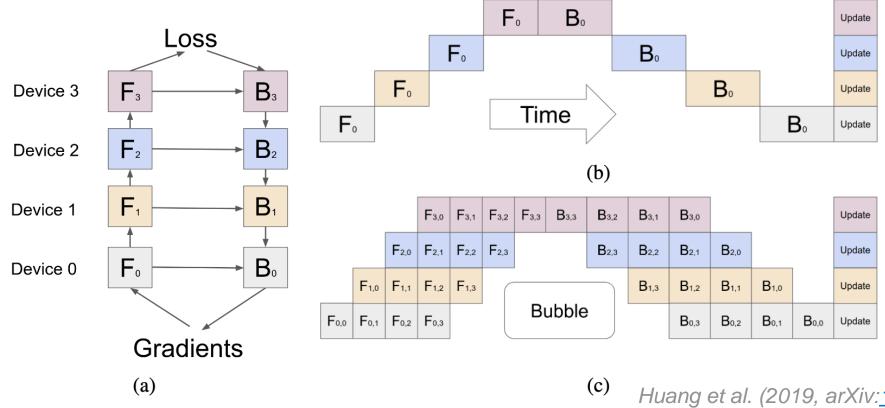
- Split the model over multiple processors (CPUs/GPUs)
- Each processor sees all the data
- Communication needed both during forward and backward pass!





Pipeline parallelism

- Model parallelism, executed in a pipelined fashion over multiple micro-batches.
- More efficient than model parallelism: hides communication with computation
- E.g. Gpipe (for PyTorch / TensorFlow), PipeDream (PyTorch)





Parallelization for deep learning

Very generic statements (there are exceptions) on model vs data parallelism

	Data parallel	Model parallel
Throughput	Increases	Decreases
Implementation	Relatively straigtforward	More difficult
Communication	Low to moderate	High
Typical use case	Speedup training	Models with large memory requirement

Use data parallel whenever you can. Use model parallel if you *really* need to. Even then, consider alternatives:

- Model pruning
- Use different hardware architecture (e.g. data parallel @ CPU)
- Use pipeline parallelism



Data parallel stochastic gradient descent

- Most networks are trained using stochastic gradient descent (SGD)
- Distributed stochastic gradient can be done in two ways
 - Synchronous SGD
 - Asynchronous SGD

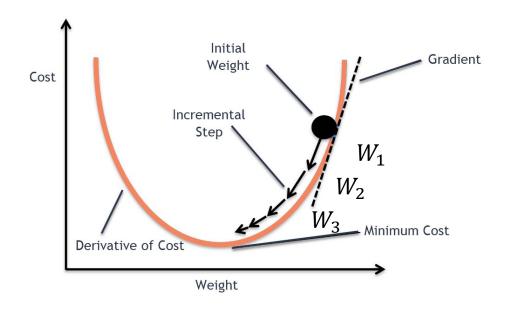


Stochastic gradient descent (SGD)

SGD: find optimum by following the slope

$$w = w - \eta \nabla Q(w)$$

w = weights, η = learning rate, $\nabla Q(w)$ = gradient for current batch.





Data parallel synchronous SGD

- Each device (j) computes the gradients ($\nabla Q_i(w)$) based on its own batch!
- Needs to be aggregated before updating weights

Device 1
$$g_1 = \nabla Q_1(w)$$

Device 2
$$g_2 = \nabla Q_2(w)$$

Device 3
$$g_1 = \nabla Q_3(w)$$

Device 4
$$g_1 = \nabla Q_4(w)$$

$$\nabla Q(\mathbf{w}) = \sum_{j} \nabla Q_{j}(\mathbf{w})$$



$$w = w - \eta \nabla Q(w)$$



Data parallel synchronous SGD

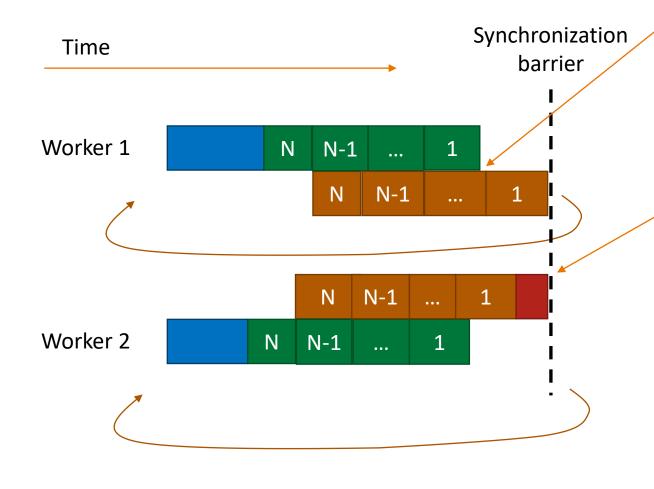
Effect on batch size:

- For N workers that each see n examples: batch size effectively $n \times N$.
- Larger batch => generally needs to be compensated by higher learning rate.
- No exact science!
 - Some use $\eta_{distributed} = \eta_{serial} \cdot N$
 - Some use $\eta_{distributed} = \eta_{serial} \cdot \sqrt{N}$
 - Experiment!



Data parallel synchronous SGD

A different view...



Pro tip:

- Overlap communication and computation (don't waste compute cycles waiting for communication!)
- *Most* (distributed) DL frameworks already take care of this for you ©

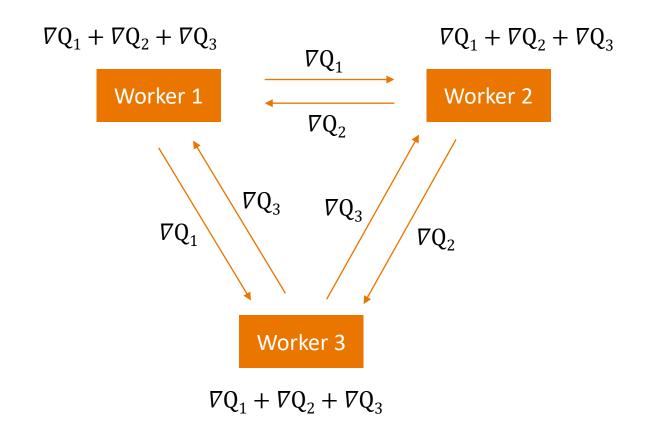
Pro tip 2:

- Penalty for *synchronous* SGD: you have to wait for the slowest worker, before next iteration.
- Make sure all workers are equally fast!
- Forward pass
- N Backward pass, compute gradient of layer N
- N Backward pass, communicate gradient N
- Wait



Decentralized data parallel synchronous SGD

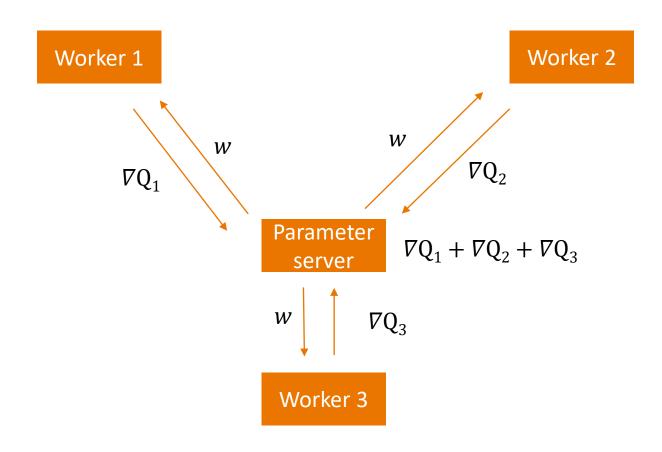
Gradients are communicated and aggregated by all workers





Centralized data parallel synchronous SGD

There is also an alternative, where a parameter server is used to aggregate the gradients, and distribute the updated model:





Centralized vs decentralized

- Centralized approach does not scale well: parameter servers create a communication bottleneck
- More info, see e.g. https://arxiv.org/pdf/1705.09056.pdf



Data parallel asynchronous SGD

We skip this, you can look at the slides later if interested...



Data parallel asynchronous SGD

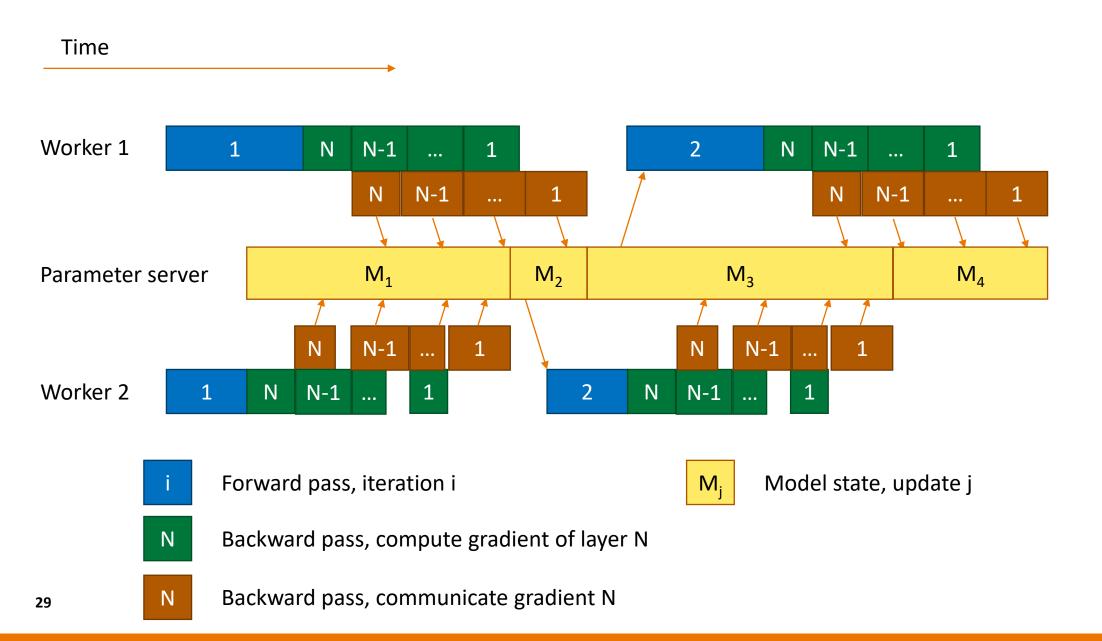
Gradients are not aggregated, weights are updated as soon as *one* worker (*j*) finishes, with the gradient of that worker:

$$w = w - \eta \nabla Q_{j}(w)$$

- If another worker finishes, it does an update on the *current* set of weights (even though the gradient may have be computed based on an earlier version of the weights)
- Asynchronous SGD does not have the same convergence guarantees as synchronous SGD
- Generally scales well, because there is no 'barrier' that induces wait time
- More info, see e.g. https://ai.google/research/pubs/pub45187



Data parallel asynchronous SGD



Data parallel SGD

Word of warning:

- Data parallel *asynchronous* SGD is an active area of research
- Data parallel synchronous SGD is well understood and is currently the accepted and advised approach



Communicating gradients...

Ok, so the most widely accepted approach is

distributed...

data parallel...

synchronous...

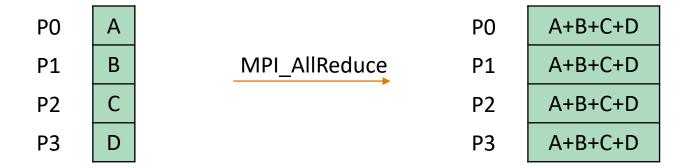
SGD...

... but how do distributed deep learning frameworks aggregate their gradients in such a setup?



MPI / NCCL

- MPI often used to communicate gradients
- MPI_AllReduce aggregates and sums gradients (remember: $\nabla Q(w) = \sum_j \nabla Q_j(w)$)
- NVIDIA's NCCL library contains an implementation of MPI routines optimized for GPU
 GPU communication





Frameworks for distributed learning

- TensorFlow's tf.distribute: quite tricky to program. Lot's of code changes needed from serial to distributed (https://www.tensorflow.org/guide/distributed training)
- TensorFlow + Horovod: serial => distributed with minimal code changes (https://horovod.readthedocs.io/en/stable/tensorflow.html)
- PyTorch's torch.distributed (https://pytorch.org/tutorials/intermediate/dist-tuto.html)
- PyTorch + Horovod: serial => distributed with minimal code changes (https://horovod.readthedocs.io/en/stable/pytorch.html)
- PyTorch Lightning: hides a lot of boiler plate code (also nice for serial training). Very little
 changes needed between serial & parallel execution, especially on a SLURM cluster
 (https://pytorch-lightning.readthedocs.io/en/latest/clouds/cluster.html#slurm-managed-cluster)



Performance considerations

Bottlenecks can be:

- I/O, especially with distributed training (many processes reading the same files)
 - Tip: don't read many small files, pack samples together into a few large files
- Communication
 - Don't use parameter servers. Use NCCL backend on NVIDIA GPUs. Don't try to parallelize very light workloads – communication overhead too large
- Memory bandwidth
 - High level frameworks (PyTorch, TensorFlow) use optimized low level libraries (cuDNN, MKL-DNN, etc) for optimized



Recap of goal: understand docs of DL frameworks

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Practical tips & take home messages

- If increased throughput is the goal, use data parallelism
- If a large model is the goal, use model (or hybrid or pipeline) parallelism, but consider the consequences (slower training) and alternatives (model pruning, CPUbased training, etc)
- Account for the difference in convergence behavior of data parallel SGD, e.g. by adjusting & experimenting with the learning rate.
- Synchronous parallel SGD is the most common approach for distributed learning, because it is well understood. Asynchronous parallel SGD can scale very well, but convergence behavior is less clear.
- Use an efficient backend for collective communications.



Further reading

- Distributed TensorFlow using Horovod:
 https://towardsdatascience.com/distributed-tensorflow-using-horovod-6d572f8790c4
- Demystifying Parallel and Distributed Deep Learning: An In-Depth Concurrency Analysis: https://arxiv.org/pdf/1802.09941.pdf
- Prace best practice guide for Deep Learning: http://www.prace-ri.eu/IMG/pdf/Best-Practice-Guide-Deep-Learning.pdf
- Technologies behind Distributed Deep Learning:
 https://preferredresearch.jp/2018/07/10/technologies-behind-distributed-deep-learning-allreduce/

