ZeRO: Memory Optimizations Toward

Training Trillion Parameter Models

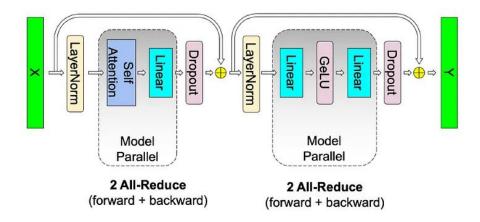
Related Works

Model Parallelism

Model Parallel: memory usage and computation of a model is distributed across multiple workers

Megatron-LM:

- + Split the matrix into multiple parts and do matmul separately.
- + No sync point within Linear and Self-attn.



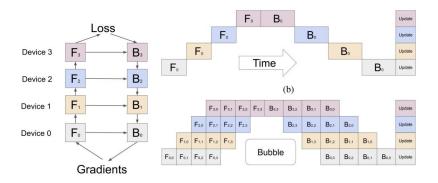
Pipeline Parallelism

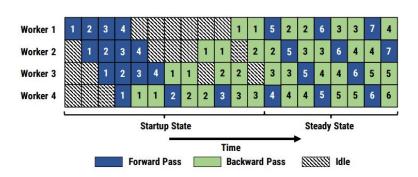
The model is distributed across multiple GPUs over layers.

Devices can be idle while waiting for others

Pipeline Parallelism is a specific form of model parallelism

- + **GPipe:** Divides input data mini-batches into smaller micro-batches.
- + **PipeDream:** Start backward as soon as possible. Do async gradient update.
- + Note that PipeDream is not equivalent to traditional DL.

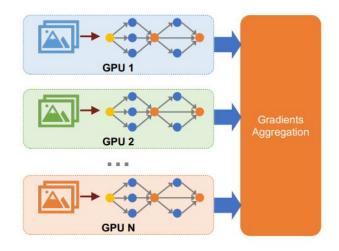




Data Parallelism

Each device has the same model and do forward and backward on a mini-batch separately. Quite easy and intuitive, but ...

- Cannot train LLM that cannot fit into one device.
- 2. Each device has the whole replica of the model.



Reduce

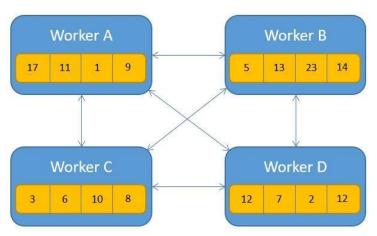
Goal: In data parallelism, it is essential to ensure that each device is updated coherently, therefore we need to aggregate(reduce) gradients across different devices.

 Centralized Reduce: all workers communicate with parameter servers for weights update; cannot scale to large numbers of workers

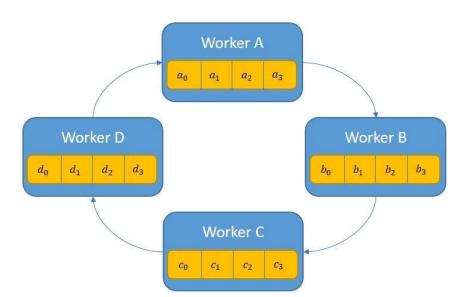
- All Reduce
- Naïve AllReduce
- Ring AllReduce

Naïve AllReduce

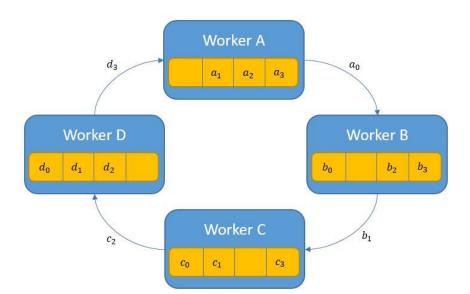
- Each worker can send its local gradients to all other workers
- If we have N workers and each worker contains M parameters
- Overall communication: N * (N-1) * M parameters
- Issue: each worker communicates with all other workers; same scalability issue as parameter server



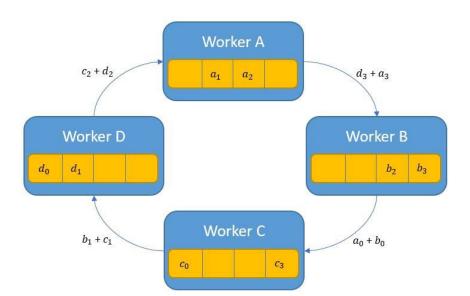
- Construct a ring of N workers, divide M parameters into N slices
- Step 1 (Aggregation): each worker send one slice (M/N parameters) to the next worker on the ring; repeat N times



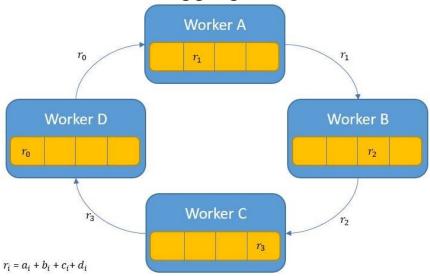
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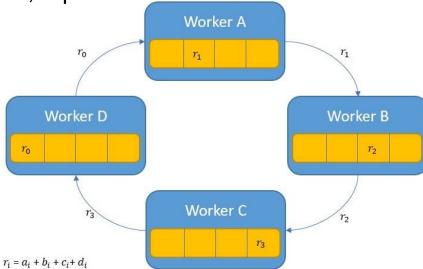


- Construct a ring of N workers, divide M parameters into N slices
- Step 1 (Aggregation): each worker send one slice (M/N parameters) to the next worker on the ring; repeat N times
- After step 1, each worker has the aggregated version of M/N parameters



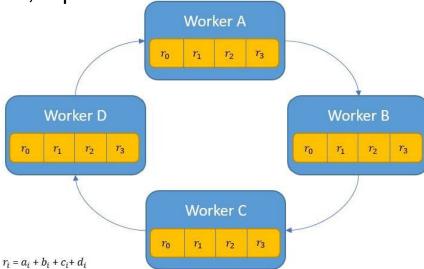
- Construct a ring of N workers, divide M parameters into N slices
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 Step 2 (Broadcast): each worker send one slice of aggregated parameters to the next worker; repeat N times



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- Step 1 (Aggregation): each worker send one slice (M/N parameters) to the next worker on the ring; repeat N times
- Step 2 (Broadcast): each worker send one slice of aggregated parameters to the next worker; repeat N times
- Overall communication: 2 * M * N parameters
 - Aggregation: M * N parameters
 - Broadcast: M * N parameters

Summary

| | Pros | Cons |
|-------------------|---------------------------------------|--|
| Model Parallelism | Good memory efficiency | Poor compute /communication efficiency (5% of peak perf in training 40B model with Megatron) |
| Data parallelism | Good compute/communication efficiency | Poor memory efficiency (Every device has one copy of model) |

Question: How can we reduce memory footprint of DP?

Understanding Memory Consumption

The GPUs need to store

- Model weights
- Forward activation
- Backward gradient
- Optimizer state

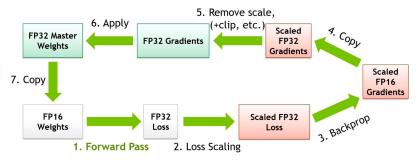
Memory Usage

Common methods in optimization: Adam + Mixed-precision Training

- + Optimizer States: Momentum + Variance
- + Model: Parameters and Gradients

MIXED PRECISION TRAINING

while θ_t not converged do $t\leftarrow t+1$ $g_t\leftarrow \nabla_{\theta}f_t(\theta_{t-1})$ (Get gradients w.r.t. stochastic objective at timestep t) $m_t\leftarrow \beta_1\cdot m_{t-1}+(1-\beta_1)\cdot g_t$ (Update biased first moment estimate) $v_t\leftarrow \beta_2\cdot v_{t-1}+(1-\beta_2)\cdot g_t^2$ (Update biased second raw moment estimate) $\widehat{m}_t\leftarrow m_t/(1-\beta_1^t)$ (Compute bias-corrected first moment estimate) $\widehat{v}_t\leftarrow v_t/(1-\beta_2^t)$ (Compute bias-corrected second raw moment estimate) $\theta_t\leftarrow \theta_{t-1}-\alpha\cdot \widehat{m}_t/(\sqrt{\widehat{v}_t}+\epsilon)$ (Update parameters) end while



Adam Optimizer

Mixed-precision Training

Memory Usage

Example: Adam as optim, and Mixed-precision Training. N parameters

- + FP32 master parameters: 4N Bytes
- + FP32 optimizer states: 4N * 2 Bytes (Momentum and Variance)
- + FP16 model parameters: 2N Bytes
- + FP16 optimizer states: 2N Bytes (Momentum only)

16N Bytes in total!

For 1.5B GPT-2, 24GB vMem

For 175B GPT-3, 2800GB vMem

Other Memory Usages

+ Activations:

+ As a concrete example, the 1.5B parameter GPT-2 model trained with sequence length of 1K and batch size of 32 requires about 60 GB of memory

+ Temporary Buffers:

+ Storing intermediate results. Operations such as gradient all-reduce, or gradient norm computation tend to fuse all the gradients into a single flattened buffer before applying the operation in an effort to improve throughput.

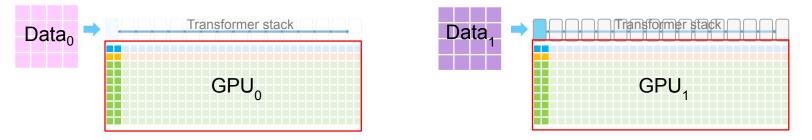
+ Memory Fragmentation:

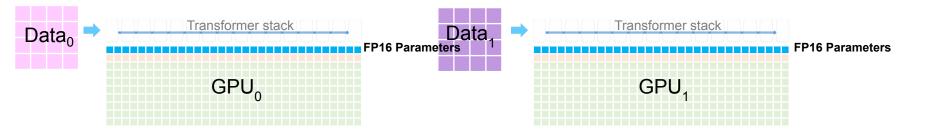
+ In extreme cases can be 30%.



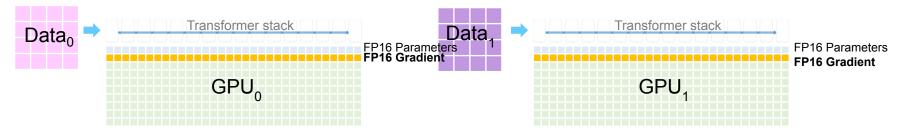
Suppose there are

- Two data splits: Data₀ and Data₁
- Two GPUs: GPU₀ and GPU₁
- 16 layer Transformer Model

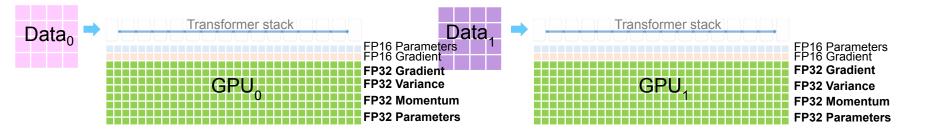




- FP16 parameters
- FP16 Gradients
- FP32 Optimizer States (Gradients, Variance, Momentum, Parameters)



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Common Approaches to Reduce Memory

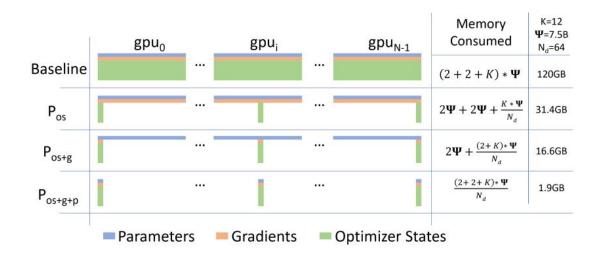
- + Reducing Activation Memory
 - + Activation Checkpoint, Compression
 - + All Work in parallel with ZeRO
- CPU Offload
 - + Requires CPU-GPU-CPU transfer, which can take 50% time
- + Memory Efficient Optimizer
 - + Maintaining coarser-grained statistics of model parameters and gradients.
 - + Works in parallel with ZeRO

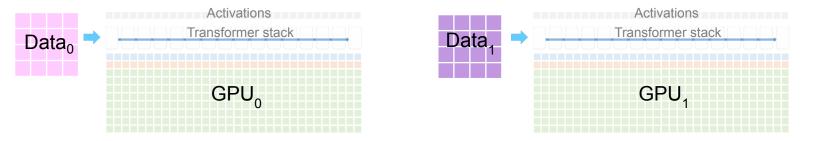
ZeRO - Zero Redundancy Optimizer

Work done by Microsoft, implemented in Deepspeed.

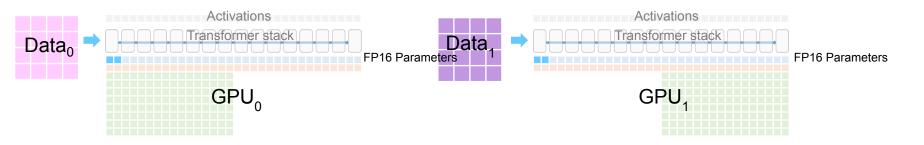
Features:

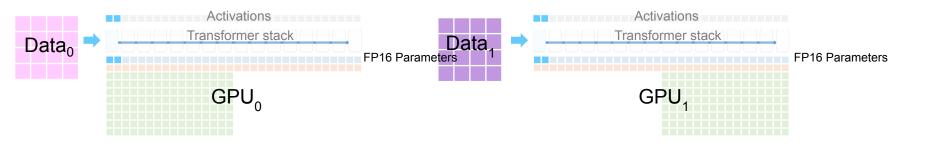
- + Eliminating data redundancy in data parallel training
- + Can be widely used in large language model training.

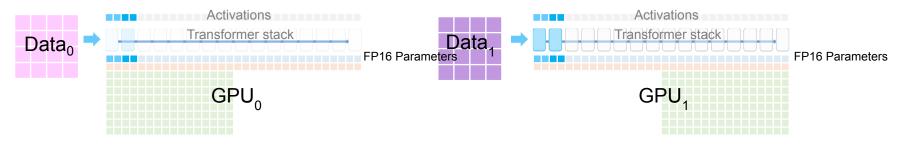


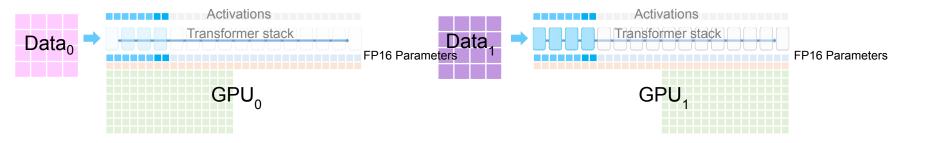


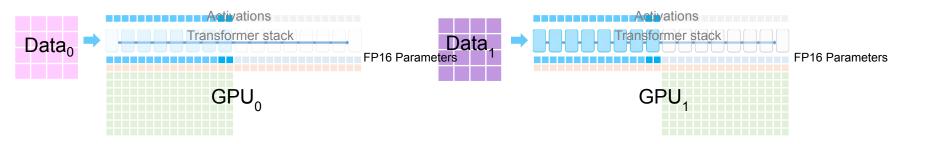
Question: How can we partition optimizer states?

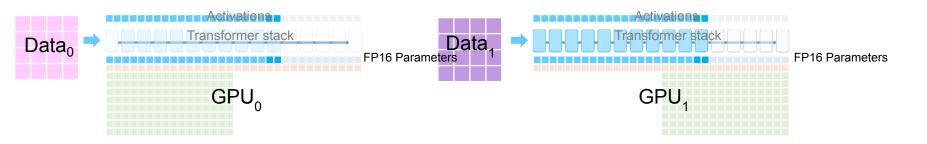


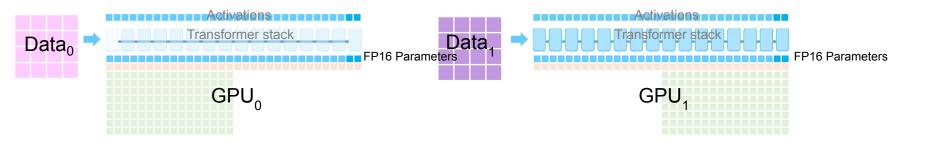


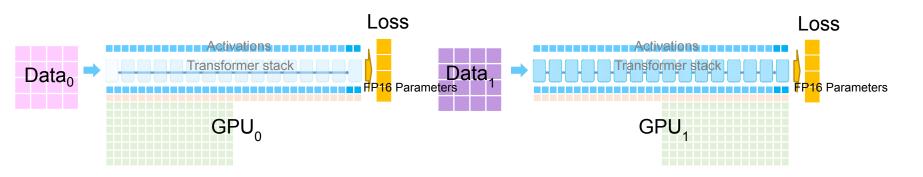


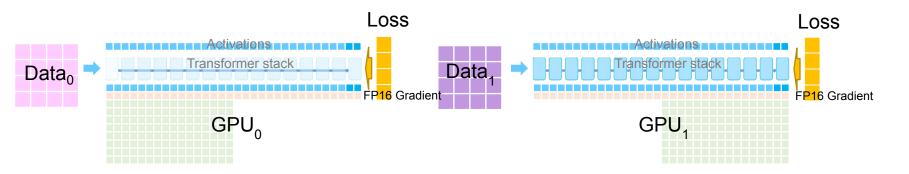




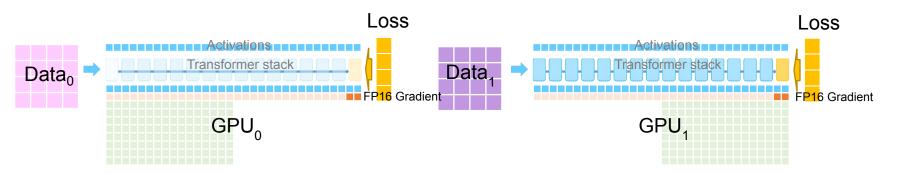


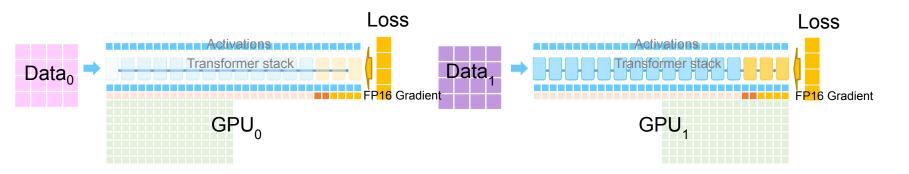


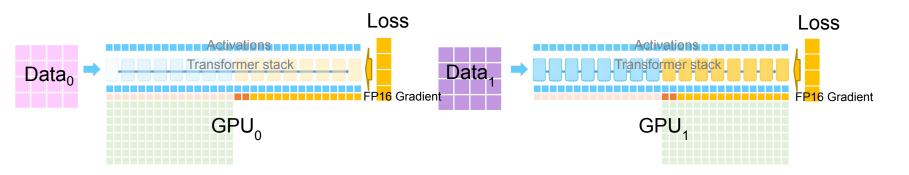


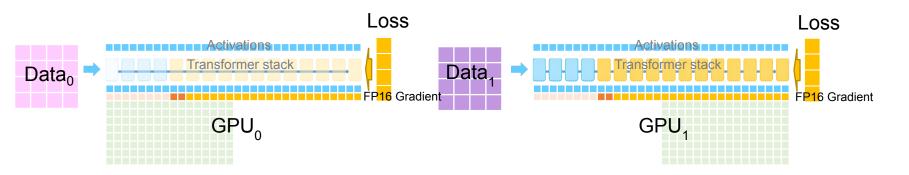


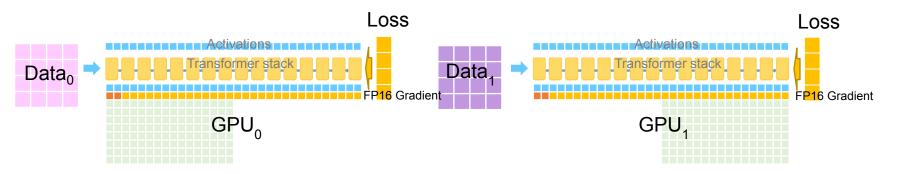
loss backward to calculate fp16 gradients

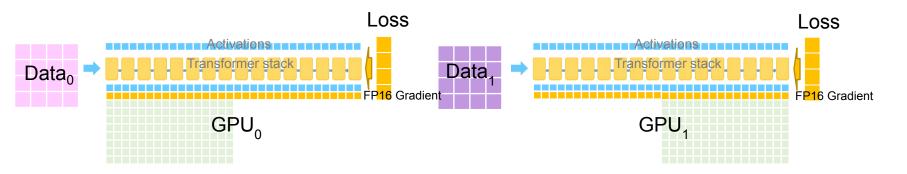


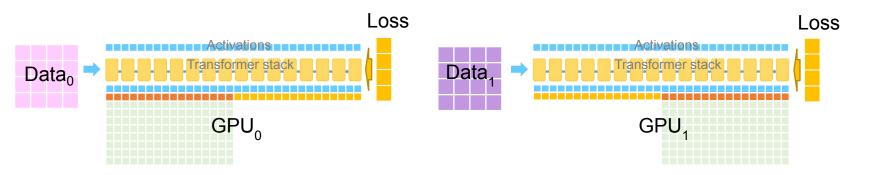




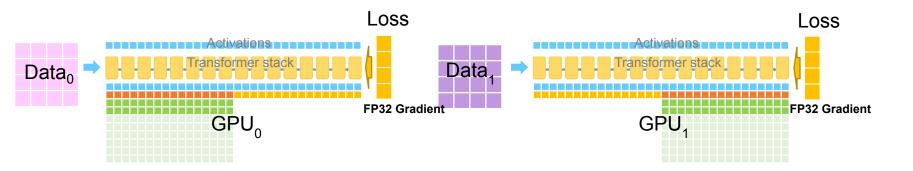




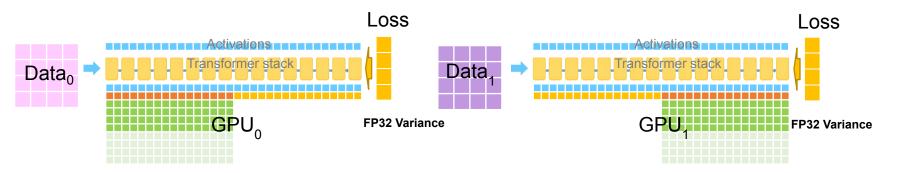




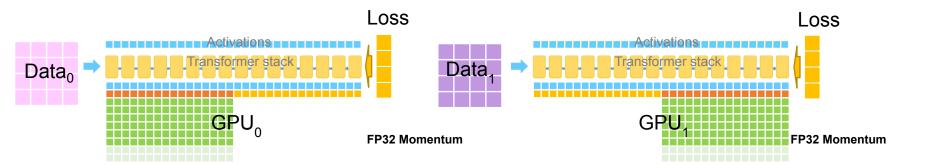
- gradient gathering from another GPU and average gradient calculation



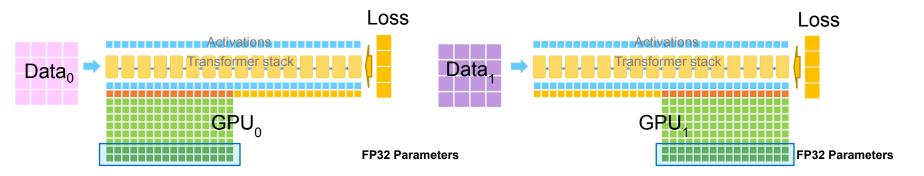
- fp32 gradient update



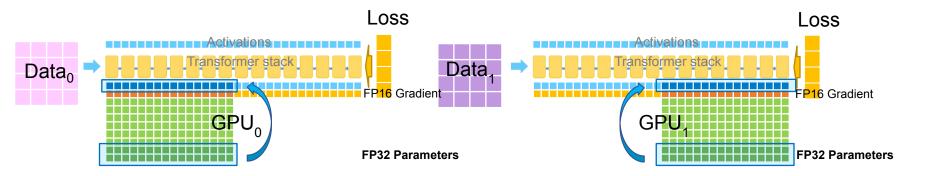
- fp32 variance update



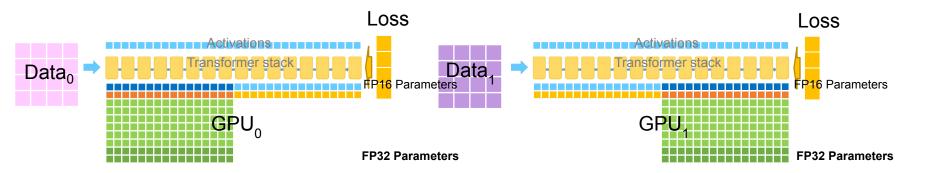
- fp32 momentum update



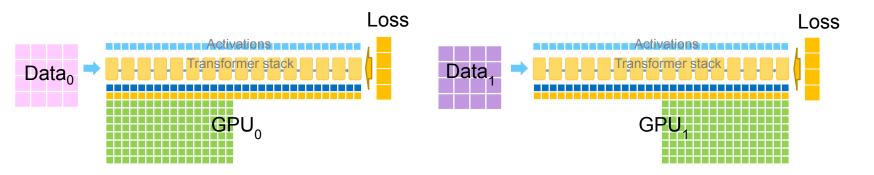
- fp32 parameters update



- fp32 parameters update using fp16 gradient



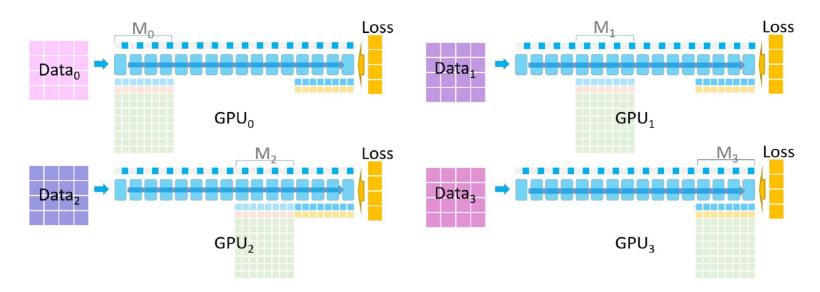
- fp16 parameters update using fp32 parameters



all gather the gp16 weights to complete the iteration

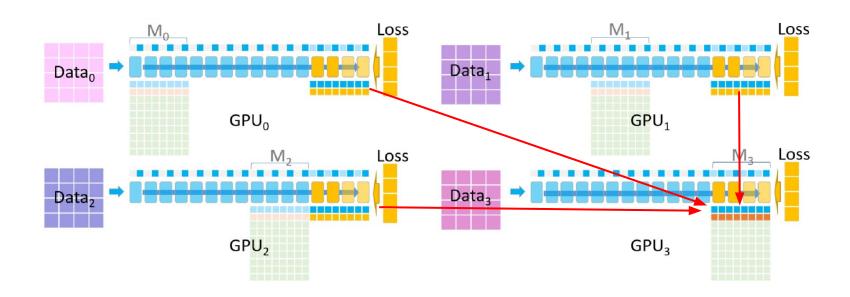
Key idea:

- Each GPU is only responsible for one partition of the parameters, so it should also only be responsible for one partition of gradients that are corresponding to their designated parameters.
- But different GPUs are responsible for different data, meaning we still need to run all GPUs for all gradients.
- Therefore, during backward pass, a GPU can immediately delete the gradients that it's not responsible for, after it passes those gradients(computed with its data) to the GPU responsible for those gradients.
- The result is each GPU can hold less memory for gradients(linear to # of GPUs)

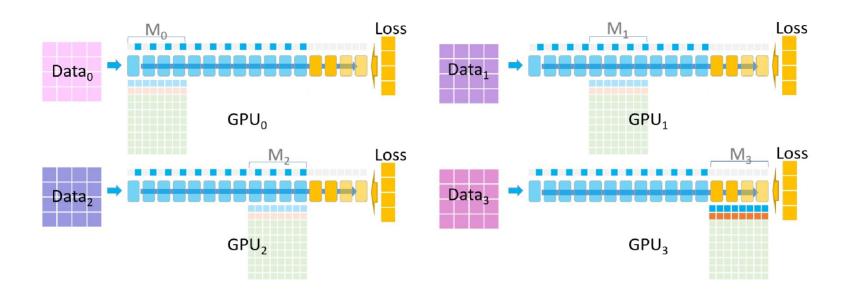


The backward pass starts

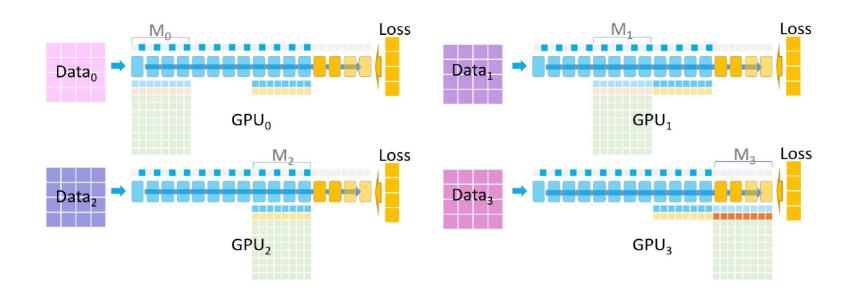
GPU 0,1,2 hold temporary buffers for the gradients that GPU 3 is responsible for (M3)



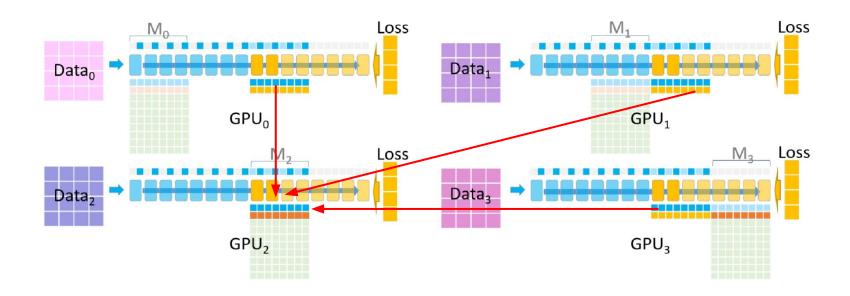
GPU 0,1,2 pass the M3 gradients to GPU 3



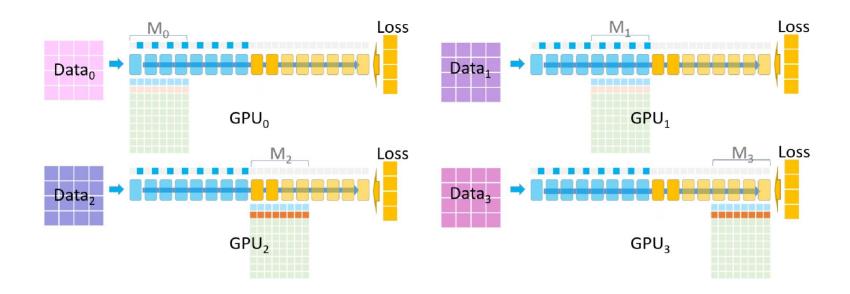
Then they delete M3 gradients, GPU 3 will keep M3 gradients.



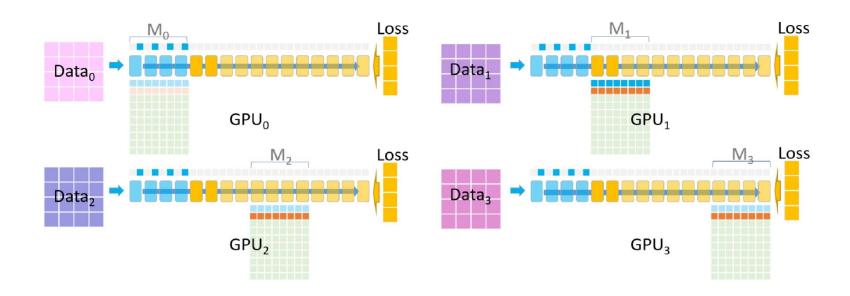
GPU 0,2,3 hold temporary buffers for the gradients that GPU 2 is responsible for (M2)



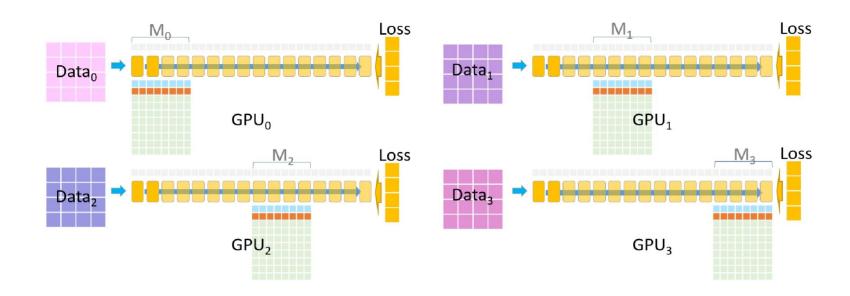
GPU 0,2,3 pass the M2 gradients to GPU 2



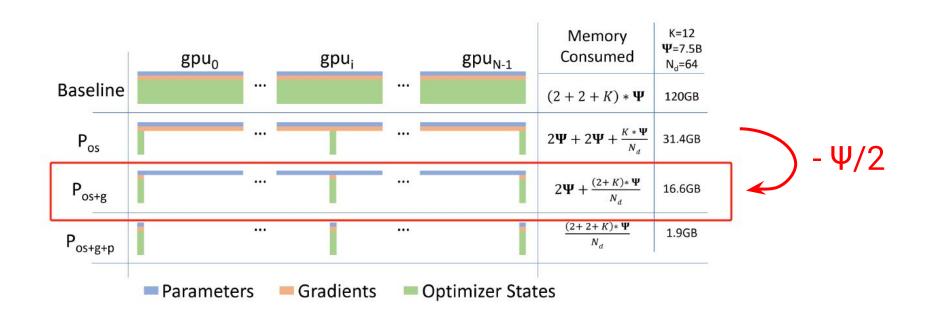
Then they delete M2 gradients, GPU 2 will keep M2 gradients.



Same thing for GPU1/M1



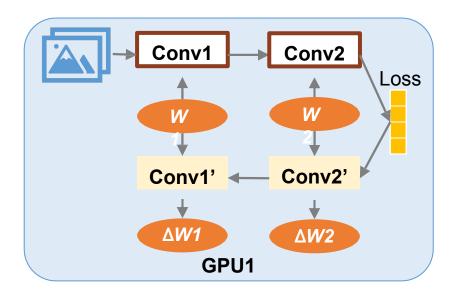
Same thing for GPU0/M0

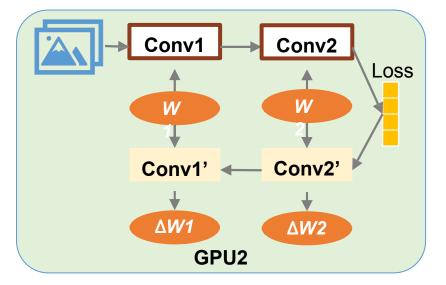


| DP | 7.5 B Model (GB) | | | 128 B Model (GB) | | | 1T Model (GB) | | |
|------|-------------------------|------------|--------------|-------------------------|------------|--------------|---------------|------------|--------------|
| | P_{os} | P_{os+g} | P_{os+g+p} | P_{os} | P_{os+g} | P_{os+g+p} | P_{os} | P_{os+g} | P_{os+g+p} |
| 1 | 120 | 120 | 120 | 2048 | 2048 | 2048 | 16000 | 16000 | 16000 |
| 4 | 52.5 | 41.3 | 30 | 896 | 704 | 512 | 7000 | 5500 | 4000 |
| 16 | 35.6 | 21.6 | 7.5 | 608 | 368 | 128 | 4750 | 2875 | 1000 |
| 64 | 31.4 | 16.6 | 1.88 | 536 | 284 | 32 | 4187 | 2218 | 250 |
| 256 | 30.4 | 15.4 | 0.47 | 518 | 263 | 8 | 4046 | 2054 | 62.5 |
| 1024 | 30.1 | 15.1 | 0.12 | 513 | 257 | 2 | 4011 | 2013 | 15.6 |
| | | | 200 | | | | | | |

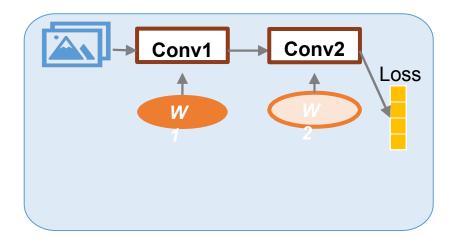
Looks pretty good!

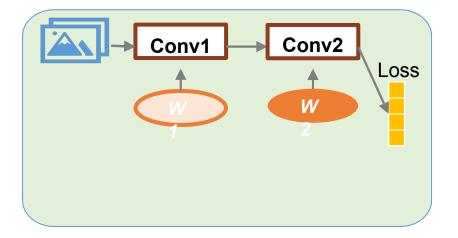
• In data parallel training, all GPUs keep all parameters during training





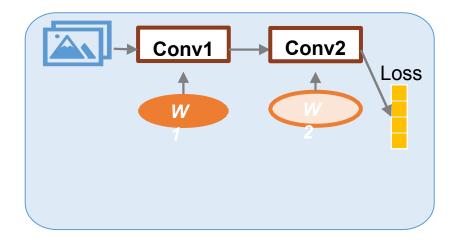
• In ZeRO, model parameters are partitioned across GPUs

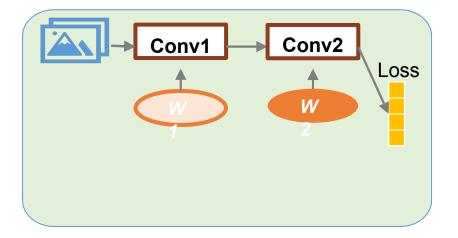




GPU1 GPU2

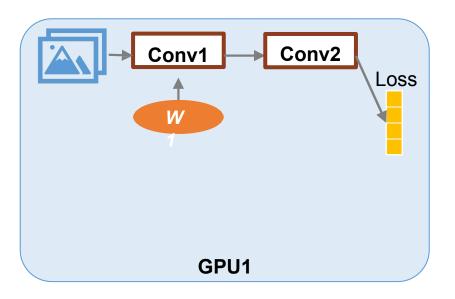
- In ZeRO, model parameters are partitioned across GPUs
- GPUs broadcast their parameters during forward

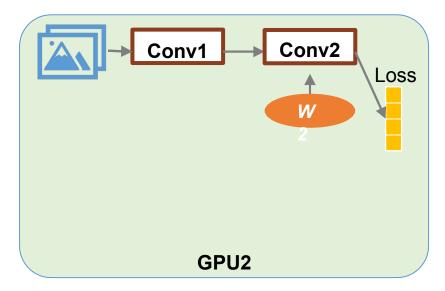




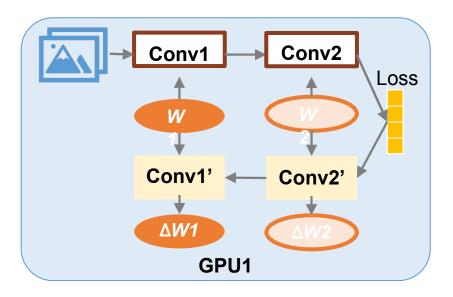
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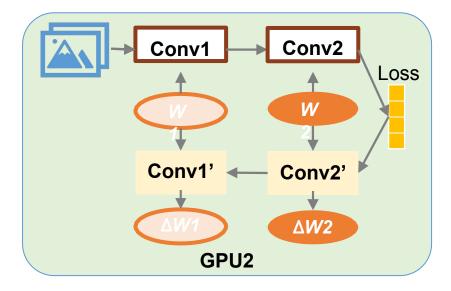
- In ZeRO, model parameters are partitioned across GPUs
- Parameters are discarded right after use





- In ZeRO, model parameters are partitioned across GPUs
- GPUs broadcast their parameters again during backward





Zero-DP Summary

- Zero-DP stage 1 and 2 (optimizer state and gradient) doesn't additional communication, while enabling up to 8x memory reduction
- Zero-DP stage 3 (parameter) incurs a maximum of 1.5x communication

ZeRO - R

- Partitioned Activation Checkpointing
 - Split every activation to different devices. Gather them when needed.
- Constant Size Buffers
 - Buffer is used in doing all-reduce to improve bandwidth.
 - Modern implementations fuses all the parameters into a single buffer.
 - ZeRO uses constant size buffers to be more efficient for a large model.
- 3. Memory Defragmentation
 - Long-lived memory (Model parameters, Optimizer state): Store together
 - Short-lived memory (Discarded activations)

Results

Theoretical: On a 32GB V100 clusters (Up to 1024 V100),

- 1. Enable the training of a model with 1 Trillion (1000B) parameters using 1024 V100.
- 2. There is no limit to the number of GPUs. (So probably more)

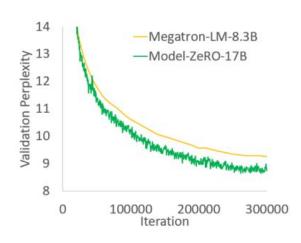
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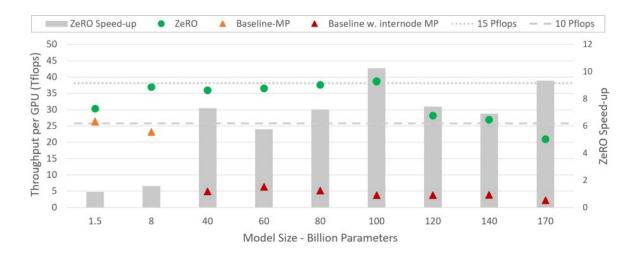
Per-device memory consumption of different optimizations

Results

Practical:

- 1. Train a 17B model (Turing-NLG. The largest as of 2020.1) and has SOTA perplexity in Webtext-103.
- 2. Train a 100B model on 400 GPUs, achieving high throughput over baseline (~10x, 30% of the theoretical peak).





Summarization

- 1. ZeRO is a distributed learning framework with data parallelization.
- 2. ZeRO partitions model states across devices.
- ZeRO trains a new SOTA model with 17B models in 2019.