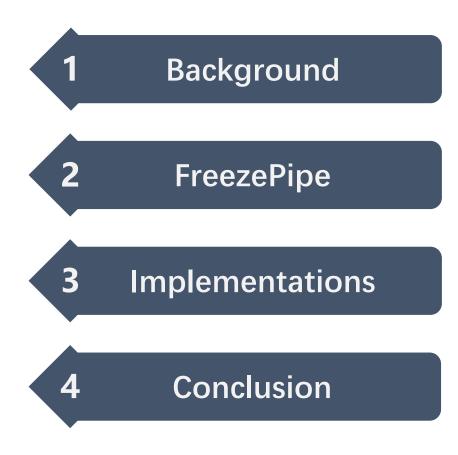


FreezePipe: An Efficient Dynamic Pipeline Parallel Approach Based on Freezing Mechanism for Distributed DNN Training

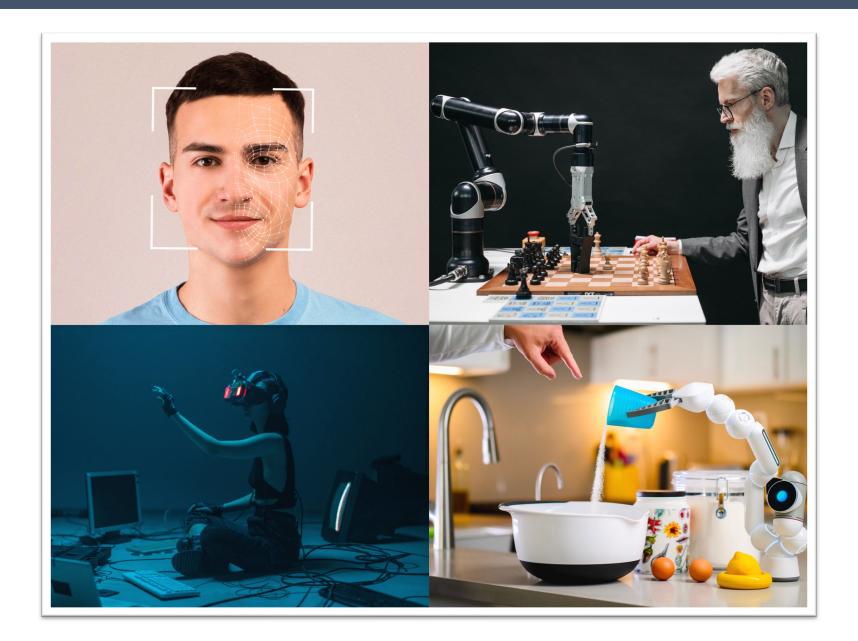
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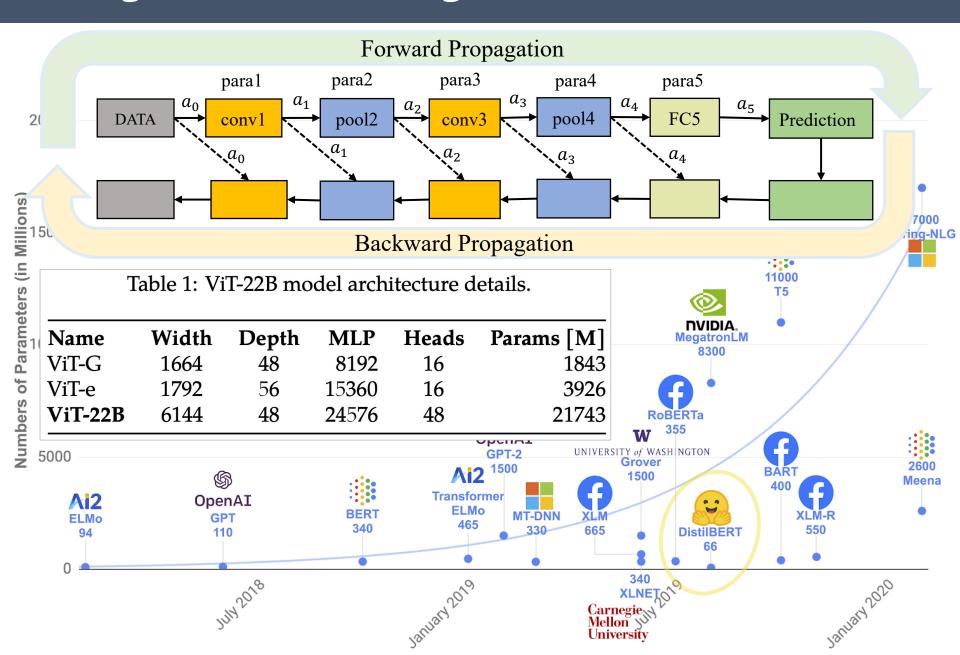
Outline



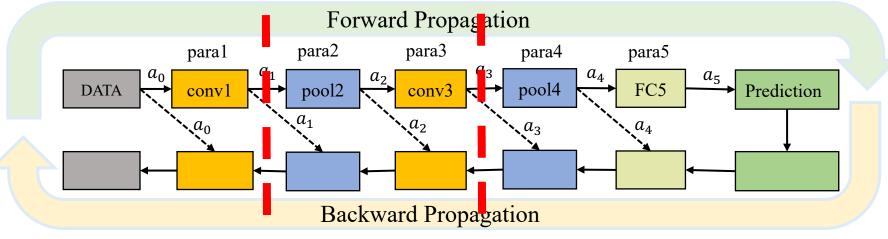
Background: Al impacts our life



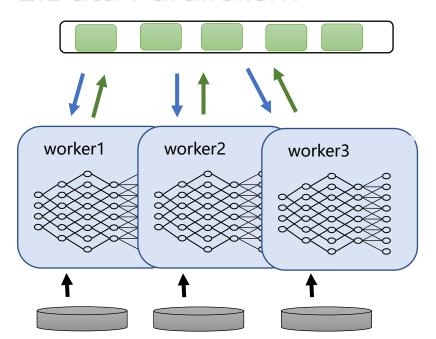
Background: Learning in distributed manner



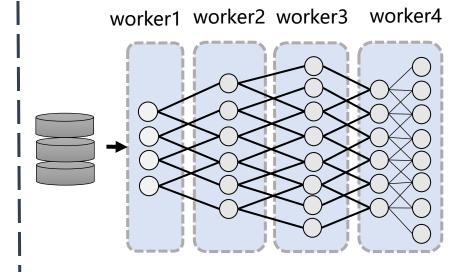
Background: Learning in distributed manner



1.Data Parallelism

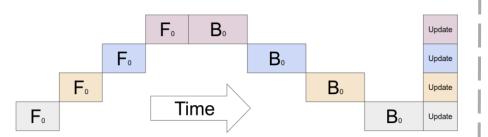


2.Model Parallelism



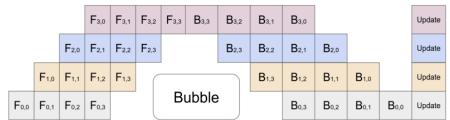
Background: Bubble time can be reduced

Model Parallelism



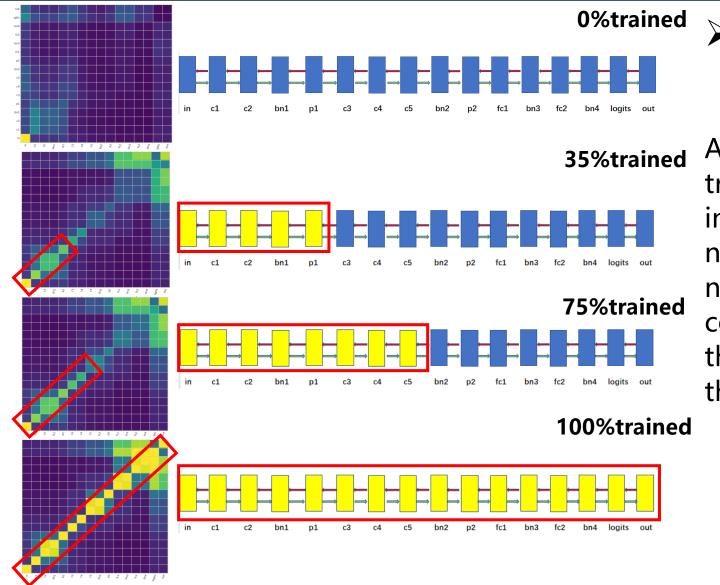
A low GPU utilization due to the neural network's computational dependency, which adversely impacts the efficiency of distributed training.

Pipeline Parallelism



➤ Using pipeline parallelism, each batch of data is further split into several micro-batches, and each micro-batch is processed by a pipeline composed of different parts of the model placed among the GPU cluster, so-called a stage, reducing GPU idle time.

Background: How the models converge?

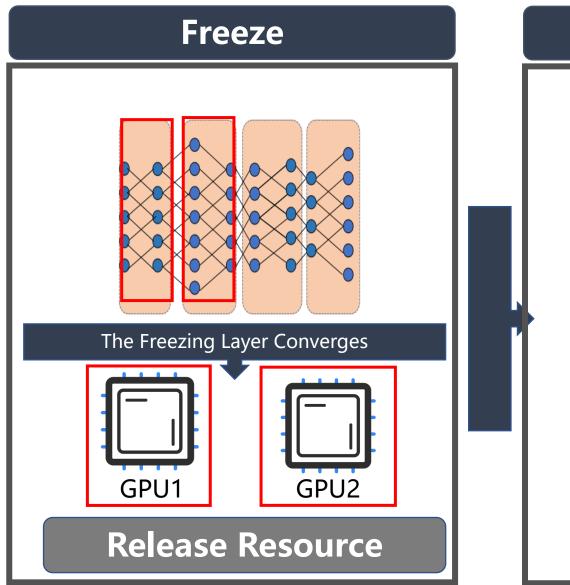


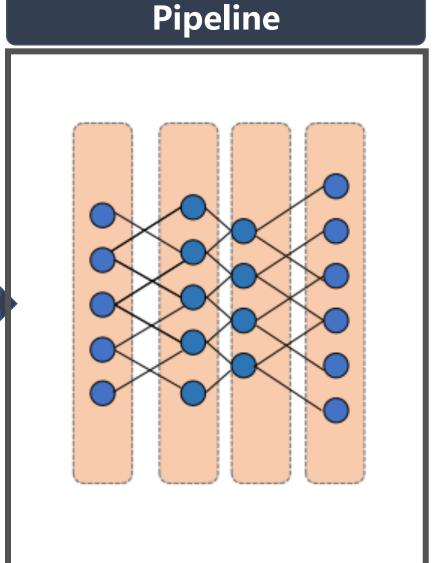
➤Why can be frozen?

As the number of training iterations increases, the neural network ,gradually converges from the input layer to the output layer.

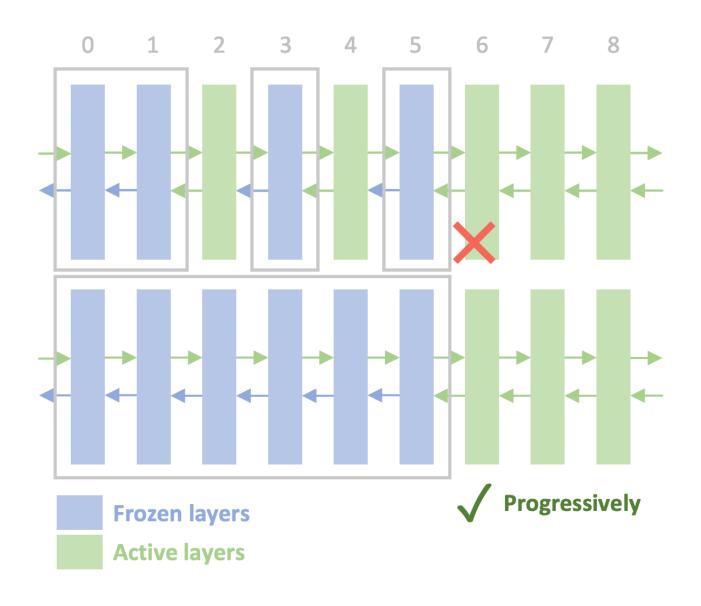
[1] Raghu M, Gilmer J, Yosinski J, et al. SVCCA: singular vector canonical correlation analysis for deep learning dynamics and interpretability[C]//Proceedings of the 31st International Conference on Neural Information Processing Systems. 2017: 6078-6087.

FreezePipe: Overview





FreezePipe: Freeze Decision



FreezePipe: Freeze Decision

Freezing rate:

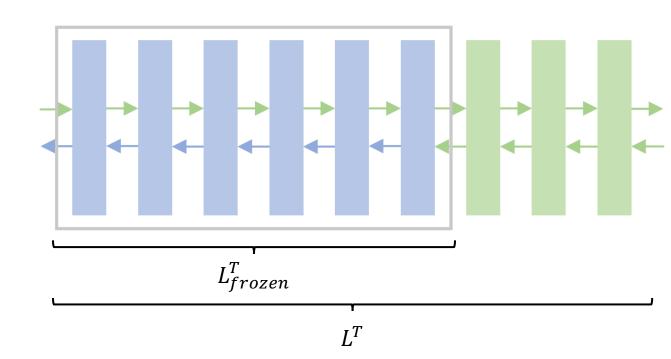
$$\beta^T = \frac{L_{frozen}^T}{L^T}$$

Cumulative gradient difference rate:

$$\eta_{l} = \frac{\left\|\sum_{i=1}^{m} G_{l_{i}}\right\|}{\left\|\sum_{i=1}^{m} |G_{l_{i}}|\right\|}$$

Freezing decision:

$$\eta_i \leq \eta_{\beta^T L^T}$$



FreezePipe: Freeze Decision

Freezing rate:

$$\beta^T = \frac{L_{frozen}^T}{L^T}$$

Cumulative gradient difference rate:

$$\eta_l = \frac{\left\|\sum_{i=1}^m G_{l_i}\right\|}{\left\|\sum_{i=1}^m \left|G_{l_i}\right|\right\|}$$

Freezing decision:

$$\eta_i \leq \eta_{\beta^T L^T}$$

Algorithm 1 Online Freezing

```
Require: Gradient difference list \{\eta_1, \eta_2, \dots, \eta_L\} ,percentile \beta^T
                             Ensure: Number of frozen layers l
                                 for i=1 to L do
                                     if \eta_i <= \eta_{\beta^T} then
                                         freeze layer_i
\eta_l = \frac{\left\|\sum_{i=1}^m G_{l_i}\right\|}{\left\|\sum_{i=1}^m |G_{l_i}|\right\|} \qquad \text{else} \\ l = i - 1, \text{break}
                                 end for
```

Lightweight Freeze Determination System:

We employ distributed gradient accumulators within each GPU to aggregate interlayer gradients, mitigating the impact of extensive gradient computations on GPU utilization. Subsequently, by calculating the gradient norms on the CPU, we determine the frozen layers

Dynamic Training Module: Model Partition

The computation time for forward propagation and backward propagation of the *i*-th layer partition:

$$comp_i = \sum (compF_i + compB_i)$$

Assuming N GPUs, M micro-batches, the total training time A for GPipe is:

$$A \approx (M + N - 1) * \max\{comp_i\} + T_{comm} + T_{update}$$

The GPipe schedule is dependent upon the maximum value of computing time at all stages:

$$\max\{comp_i\} \to A$$

Thus, the problem is transformed into a MiniMax problem:

$$\min \max\{comp_i\} \rightarrow \min A$$

Dynamic Training Module: Model Partition

Algorithm 2 Integer Binary Search

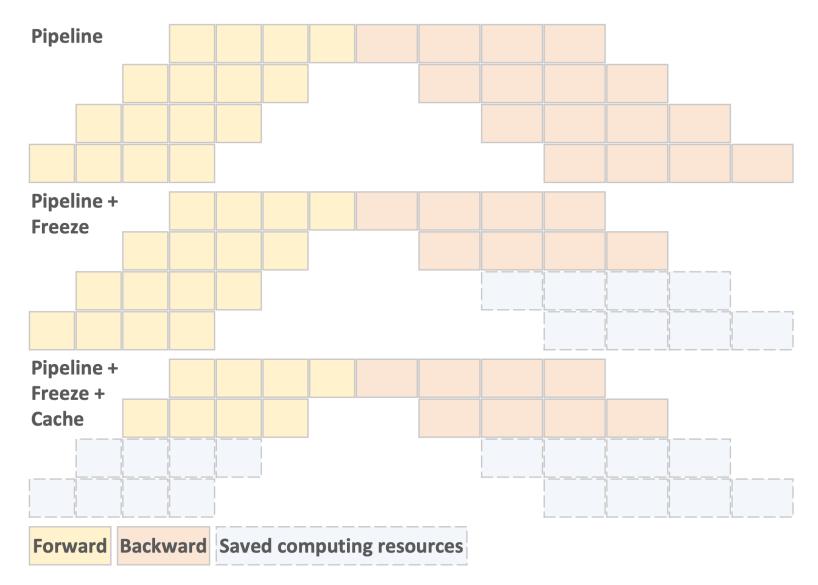
end for

```
Require: time costs list sequence = \{\mu_1, \mu_2, \dots, \mu_L\} ,the number
  of stages partitions
Ensure: partitionslist
  for i=1 to L do
    \mu_i = [\mu_i * 1000]
  end for
  low = max(sequence)
  high = sum(sequence)
  while low < high do
     mid = low + (high - low)//2
     nowpartitions = getpartition(sequence, mid)
    if nowpartitions <= partitions then
       high = mid
     else
       low = mid + 1
     end if
  end while
  maxvalue = low
  for i = 1 to Len(sequence) do
    update\ partitions list
```

2) Optimal Solution: An integer binary search method is designed to solve the Min-Max problem. Input to the algorithm is a list of time costs for each layer of the neural network and the number of stages in the pipeline, as follows: This algorithm has a time complexity of $O(N \log(\sum_{i=1}^{n} \mu_i))$, which is less than the conventional dynamic programming algorithm $O(kn^2)$ and the Torchgpipe [12] algorithm "Block Partitions" of Sequences" $O(kn^3)$.

Dynamic Training Module: Forward Cache

Automatic Caching:



Repartition after caching

After caching the intermediate activations of the frozen layers, as a result of the freed GPU computational resources, the pipeline needs to be repartitioned:

$$A = \begin{cases} A^{AfterCaching}, max\{comp_i^G\} > max\{comp_i^{AfterCaching}\} \\ A^G, otherwise \end{cases}$$

Implementation

Benchmarks:

Our primary performance metrics focus on the training time and accuracy of a deep neural network (DNN) model. In this section, we compare FreezePipe to TorchGPipe^[1] through experiments. We utilize the VGG-16 architecture with the CIFAR-10 dataset. A batch size of 64 is employed for each DNN model.

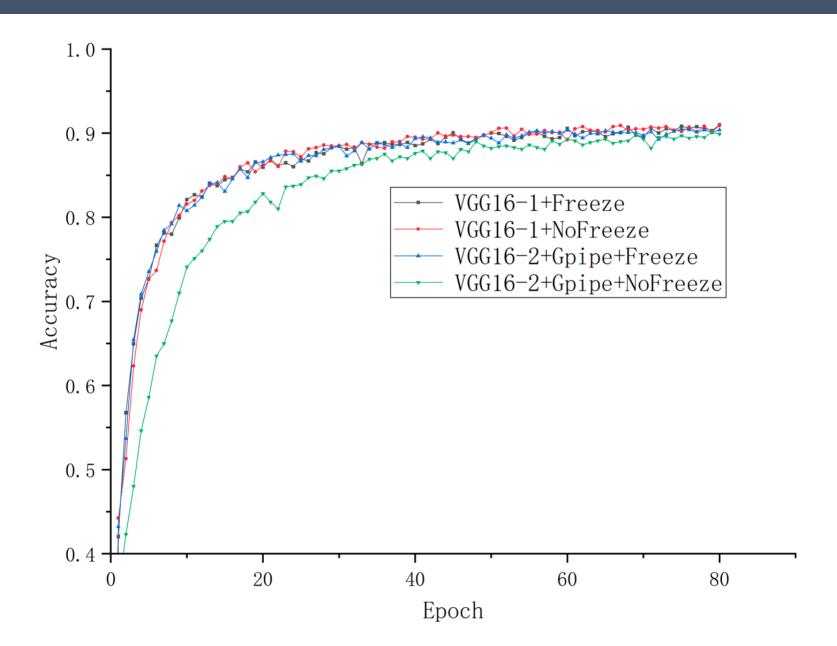
Hardware configuration:

Our primary experiments utilize a GPU cluster comprising individual machines, each outfitted with an NVIDIA GeForce RTX 3090, boasting 24GB of GPU memory. The inter-GPU bandwidth within a machine, utilizing the PCIE protocol, is 10GB/s. All machines operate on the 64-bit Ubuntu 20.04 operating system, equipped with NVIDIA-SMI 470.141.03, Driver Version 470.141.03, and CUDA Version 11.4.

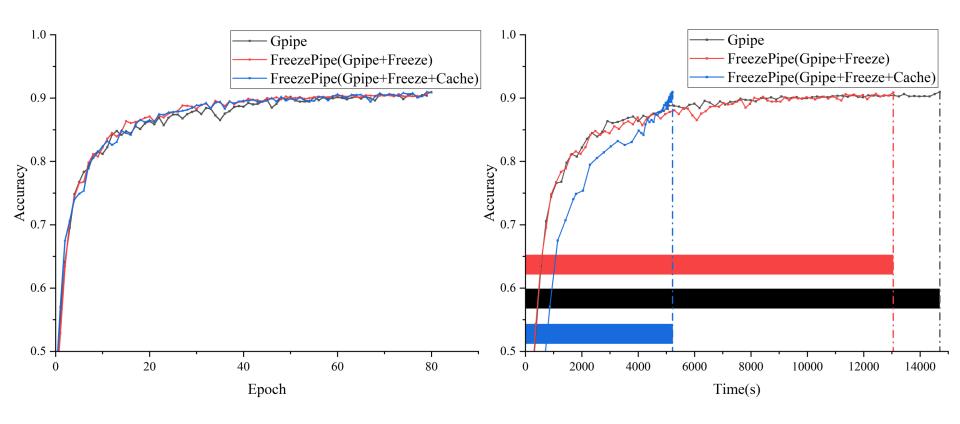
Hyperparameters: In the experiment, we fixed the freezing rate to 0.5 and performed online freezing every ten epochs, ideal results yielded.

[1] Kim C, Lee H, Jeong M, et al. torchgpipe: On-the-fly Pipeline Parallelism for Training Giant Models[J]. arXiv e-prints, 2020: arXiv: 2004.09910.

Conclusion



Conclusion



Thank You For Your Listening

