EXPLORING THE HIDDEN DIMENSION IN ACCELERATING CONVOLUTIONAL NEURAL NETWORKS

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ABSTRACT

DeePa is a deep learning framework that explores parallelism in all parallelizable dimensions to accelerate the training process of convolution neural networks. DeePa optimizes parallelism at the granularity of each individual layer in the network. We propose a novel elimination-based algorithm that finds the optimal parallelism configuration for every layer in linear time. Our evaluation shows that DeePa achieves up to 6.5× speedup compared to state-of-the-art deep learning frameworks and reduces data transfers by up to 23×.

1 INTRODUCTION

Training convolutional neural networks (CNNs) is increasingly compute-intensive and time-consuming. It takes days or even weeks to train deep CNNs from scratch (Szegedy et al., 2014; Zeiler & Fergus, 2014; Simonyan & Zisserman, 2014; Szegedy et al., 2016). Existing deep learning frameworks such as TensorFlow, PyTorch, and Caffe2 parallelize the training process onto multiple processors (usually GPUs) using *data parallelism* dividing the entire image dataset into batches with the same number of images and assigning each batch to a dedicated processor.

The standard data parallel parallelization of CNN training only exploits parallelism across images. However, there are other dimensions that also parallelize the training process. For example, in CNNs for 2D images, data is commonly organized as 4-dimensional tensors (i.e., batch, height, width, channel). The *batch* dimension includes an index for each image in the input dataset. For each image, the *height* and *width* dimensions specify a position in that image. For a particular position, the *channel* dimension [1] indexes different neurons for that position. Exploring other parallelizable dimensions can potentially reduce the compute time and data transfer cost when training CNNs (see Section 2). Moreover, different layers in a CNN may prefer different parallelism configurations for achieving optimal performance.

We propose DeePa, a deep learning framework that explores parallelism in all parallelizable dimensions to accelerate the training of CNNs. DeePa can exploit different parallelism configurations for each individual layer. To the best of our knowledge, DeePa is the first system that models the parallelism of neural networks at the granularity of each individual layer. To generate a parallelism configuration for each layer, DeePa uses a novel elimination-based algorithm that automatically finds the configuration with the best estimated performance in linear time.

The main contributions of this paper are:

- We present DeePa, a deep learning framework that explores parallelism in all parallelizable dimensions to accelerate the training of CNNs.
- The parallelization strategy is selected at the granularity of each individual layer.
- We present a novel elimination-based algorithm for finding the parallelism configuration with optimal estimated performance for each layer.
- Our evaluation shows that, compared to state-of-the-art deep learning frameworks (e.g., TensorFlow and PyTorch), DeePa achieves 6.5×, 1.9×, and 1.5× speedup for AlexNet,

[1] Some papers use the term *depth* to refer to different neurons for a position. In this paper, depth refers to the number of layers for an entire neural network and we use *channel* for the neurons for a position.

Table 1: Detailed information for the example convolutional layers used in Figure 1 and 2.

Name	Input Channels	Output Channels	Height	Width	Kernel	Stride	Description
C1	128	128	112	112	3x3	1x1	Conv4 in VGG-16
C2	512	512	28	28	3x3	1x1	Conv8, Conv9, and Conv10 in VGG-16
C3	192	64	35	35	1x1	1x1	Conv1x1 in an Inception-v3 module
C4	48	64	35	35	5x5	1x1	Conv5x5 in an Inception-v3 module
C5	64	192	27	27	5x5	1x1	Conv2 in AlexNet
C6	256	256	13	13	3x3	1x1	Conv5 in AlexNet
C7	32	64	147	147	3x3	1x1	Conv3 in Inception-v3
C8	448	384	8	8	3x3	1x1	Conv3x3 in an Inception-v3 module

VGG-16, and Inception-v3, respectively. The performance improvement comes from reducing overall data transfers, automatically overlapping computation with data movement, and accelerating computation throughput.

2 MOTIVATION

This work is motivated by the following observations.

2.1 ACCELERATING COMPUTATION THROUGHPUT

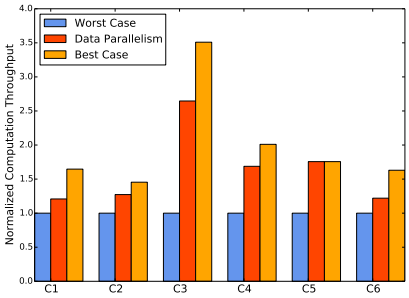


Figure 1: Relative performance for training different convolutional layers. Computation includes both forward processing and back propagation and is normalized by the worst case.

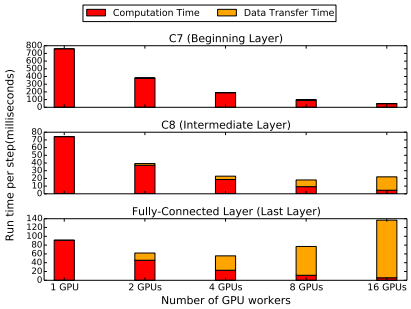
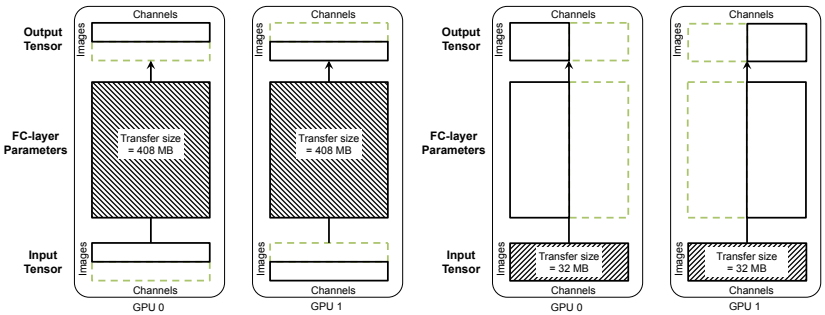


Figure 2: Computation and data transfer time to process a mini-batch of 512 images using data parallelism for the beginning layer, an intermediate layer, and the last layer of Inception-v3.

Convolutional layers generally consume the bulk of the training time in CNNs. Choosing parallelization strategies that exploit different data dimension can result in significantly different performance. Figure 1 shows the relative speed of training six different convolutional layers from AlexNet, VGG-16, and Inception-v3. The parameters of the convolutional layers are shown in Table 1. For each convolutional layer, we tried to parallelize the computation in each individual parallelizable dimension as well as combinations of different parallelizable dimensions, and we report the performance of the standard data parallel parallelization along with the worst and best parallelization strategies. Figure 1 shows that different parallelism configurations result in very different performance, and data parallelism generally achieves suboptimal performance. Therefore, exploring parallelism in other dimensions can potentially accelerate the training of convolutional layers.

2.2 REDUCING DATA TRANSFER COST

Different parallelism configurations can also result in significantly different amounts of data movement. Figure 3 shows an example of parallelizing the first fully-connected layer of VGG-16 on two GPUs through different dimensions. In data parallelism (Figure 3a), each GPU processes a batch of images and computes the gradient for the entire fully-connected layer. This requires each GPU synchronizing the gradients for the entire fully-connected layer (shown as the shadow rectangles) after each step. An alternative approach (Figure 3b) parallelizes in the channel dimension by assigning a subset of the output channels to each GPU. As a result, different GPUs compute the gradients for disjoint subsets of the fully-connected layer, which eliminates transferring the fully-connected layer



(a) Parallelism in the batch dimension. (b) Parallelism in the channel dimension.

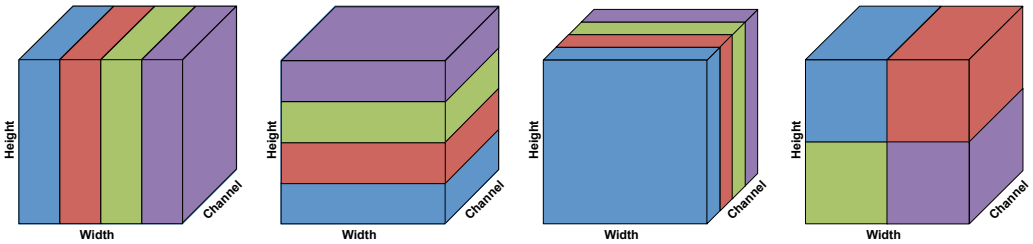
Figure 3: Different configurations for parallelizing the first fully-connected layer of VGG-16. Rectangles with solid lines indicate tensors managed by the local GPU, while rectangles with dot lines are tensors managed by a remote GPU. The shadow rectangles indicate data transfers for each step.

but introduces additional data transfers for input tensors (shown as the shadow rectangles). For this particular case, using parallelism in the channel dimension reduces data transfer costs by 12×.

2.3 Optimizing per-layer performance

To processing a certain mini-batch of images, increasing the number of workers does not always improve overall execution time, due to the data transfer overhead to synchronize gradients across different workers. Figure 2 shows the per-step training time for three different layers in Inception-v3 for a mini-batch size of 512 images on up to 16 GPUs. The training time includes forward processing, backward propagation, and gradient aggregation. The figure shows that different layers in a neural network may prefer different configurations for achieve optimal performance, and there is no single configuration that is optimal for all layers. For example, the beginning layer performs best on 16 GPUs while the last layer performs best on 4 GPUs.

3 DeePa



(a) (n=1, c=1, h=1, w=4) (b) (n=1, c=1, h=4, w=1) (c) (n=1, c=4, h=1, w=1) (d) (n=1, c=1, h=2, w=2)

Figure 4: Example configurations that parallelize an operation in a single dimension or combinations of multiple dimensions. The figure shows how each image is partitioned in different configurations.

In this section we describe DeePa, a framework that explores parallelism in all parallelizable dimensions to accelerate the training of CNNs. DeePa optimizes parallelism for each individual layer. Similar to TensorFlow and PyTorch, DeePa uses *computation graphs* to describe dependencies between operations. In a computation graph $G = (V, E)$, each node $n \in V$ is an operation (e.g., convolution or matrix-multiply), and each directed edge $(u, v) \in E$ represents a tensor that is an output of u and an input of v.

One key difference between DeePa and TensorFlow (or PyTorch) is that each node in the DeePa computation graph also includes a *configuration* that describes how the corresponding operation is parallelized across different workers. For each parallelizable dimension (i.e., batch, channel, height, and width), the configuration includes an integer that describes parallelism in that dimension. Figure 4 demonstrates some example configurations that explore parallelism in a single dimension as

well as combinations of different dimensions. DeePa assumes equal partitioning in each dimension. As a result, each worker receives the same size workload, which provides well-balanced workload distribution in our experiments.

For each node in the computation graph, its configuration describes how the *output* tensor is divided onto multiple workers. Each worker computes a *disjoint* subset of the output tensor, and thus each worker can process the operation in parallel without data dependencies. Given a node's configuration, DeePa calculates the input sets for each worker and automatically schedules proper data transfers between operations.

DeePa also provides three additional functions:

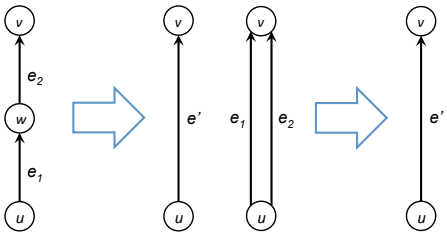
- For each node v and configuration c, $v.\mathit{compute}(c)$ estimates the time to process the corresponding operation under the parallelism configuration c. This includes both the forward processing and back propagation time and is estimated by running the operation multiple times on the device and measuring the average execution time.
- For each edge $e = (u, v)$, $e.\mathit{xfer}(c_u, c_v)$ estimates the time to transfer the input tensor e to each worker, using the size of the data to be moved and the known communication bandwidth. Note that $e.\mathit{xfer}(c_u, c_v)$ is zero if u and v have the same configuration (i.e., $c_u = c_v$), in which case no data is transferred.
- For each node v and configuration c, $v.\mathit{update}(c)$ estimates the time to update parameters for the corresponding operation. We use the data transfer time to approximate the update time, since the data transfer time is much longer than the compute time for updating parameters. Note that different configurations can have significantly different update time, as described in Section 2.2.

A *global configuration* g includes a parallelism configuration for each node in a computation graph: $g(v)$ describes the parallelism configuration for node v. With the above definitions in hand, we can model the per-step execution time for a computation graph:

$$\mathit{Cost}(g, (V, E)) = \sum_{v \in V} \{v.\mathit{compute}(g(v)) + v.\mathit{update}(g(v))\} + \sum_{e=(u,v) \in E} e.\mathit{xfer}(g(u), g(v)) \quad (1)$$

$\mathit{Cost}(g, (V, E))$ estimates the per-step execution time if the computation graph (V, E) is parallelized by using global configuration g. This execution time includes forwarding processing, backward propagation, and gradient aggregation. Equation 1 transforms the problem of finding the configuration for each individual node into an optimization problem.

4 Finding Optimal Global Configurations



(a) Node elimination. (b) Edge elimination.

Figure 5: Performing a node/edge elimination on a computation graph.

We now describe our algorithm for finding the global configuration that minimizes Equation 1. In DeePa, each node can select any parallelism configuration, therefore the number of potential global configurations is exponential in the number of nodes in a computation graph, which makes it impractical to enumerate all global configurations for deep CNNs such as VGG-16 and Inception-v3.

We observe *strong locality* in most CNNs: each node is only connected to a few nodes with similar depth in a computation graph. Based on this observation, we use the following two elimination strategies to adaptively simplify the computation graph.

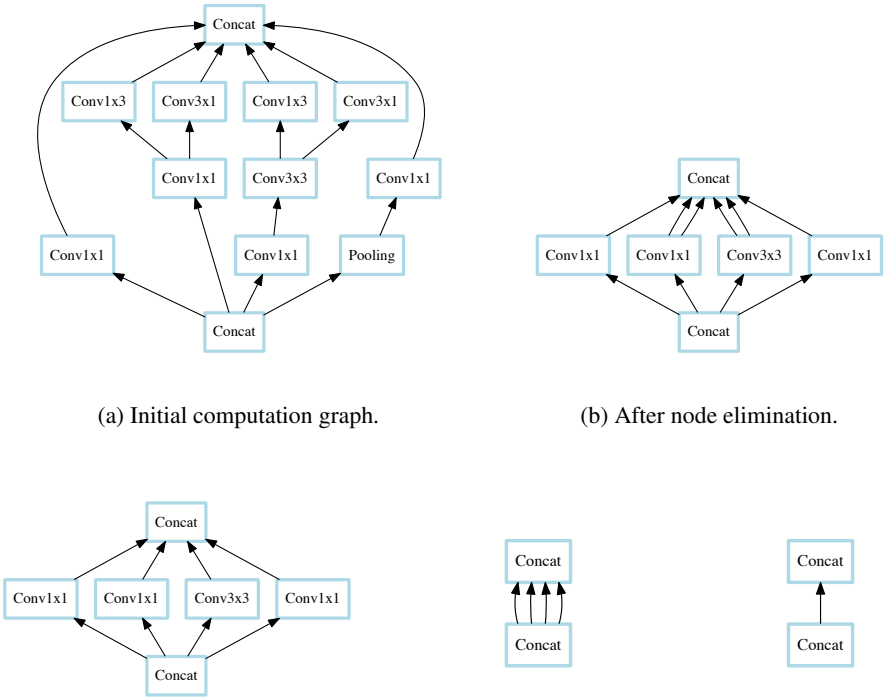
Node elimination. For each node w with a single in-edge $e_1 = (u, w)$ and a single out-edge $e_2 = (w, v)$, we can remove node w and the two edges e_1 and e_2 from the graph and insert a new edge $e' = (u, v)$ (shown in Figure 5a). The *xfer* function for node e' is

$$e'.\mathit{xfer}(c_u, c_v) = \min_{c_w}\{e_1.\mathit{xfer}(c_u, c_w) + w.\mathit{compute}(c_w) + w.\mathit{update}(c_w) + e_2.\mathit{xfer}(c_w, c_v)\} \quad (2)$$

For CNNs with a linear computation graph (e.g., AlexNet and VGG-16), node elimination is sufficient to reduce the original graph to a graph with only 2 nodes.

Edge elimination. For two edges with the same source and destination node (i.e., $e_1 = (u, v)$ and $e_2 = (u, v)$), we can remove e_1 and e_2 from the graph and insert a new edge $e' = (u, v)$ (shown in Figure 5b). The *xfer* function for node e' is

$$e'.\mathit{xfer}(c_u, c_v) = e_1.\mathit{xfer}(c_u, c_v) + e_2.\mathit{xfer}(c_u, c_v) \quad (3)$$



(a) Initial computation graph.

(b) After node elimination.

(c) After edge elimination.

(d) After node elimination.

(e) After edge elimination.

Figure 6: Iteratively performing node/edge eliminations on an Inception module.

Figure 6 shows how DeePa iteratively eliminates nodes and edges for an Inception-v3 module. For the whole Inception-v3 computation graph with 120 nodes, DeePa reduces it to a 2-node graph.

DeePa iteratively uses node and edge eliminations to simplify a computation graph until neither elimination can be applied. DeePa then enumerates all global configurations for the final graph and chooses the one that minimizes the *Cost* function in Equation 1.

After deciding the configuration for each node in the final graph, DeePa then decide the configuration for the eliminated nodes by undoing the node and edge eliminations in reverse order. When undoing a node elimination for node w, DeePa selects the configuration that minimizes Equation 2 for node w. Our algorithm is able to find a global configuration that minimizes the *Cost* function in linear time (proof in Appendix A.2). In our experiments, DeePa finds the global configuration for parallelizing Inception-v3 on 16 GPUs in approximately 100ms.

5 IMPLEMENTATION

We found that it is non-trivial to parallelize the training of CNNs in the height, width, and channel dimensions in existing frameworks (e.g., TensorFlow, PyTorch, and Caffe2), and none provides an interface for controlling per-operation parallelism. We implemented DeePa in Legion (Bauer et al., 2012), a high-performance parallel runtime for distributed heterogeneous architectures, and use cuDNN (Chetlur et al., 2014) and cuBLAS (cub, 2016) as the underlying libraries for processing neural network operations. The following Legion features significantly simplify our implementation for DeePa. First, Legion supports high-dimensional partitioning that allows us to parallelize any operation in any combination of the dimensions. Second, Legion allows DeePa to control parallelism at the granularity of each operation. Third, Legion allows fine-grain control over the placement of data in memory. Fourth, Legion's asynchronous tasking model makes it easy to exploit task as well as data parallelism. We also include two critical optimizations that help achieve good performance.

Overlapping computation with data transfers. Data parallelism in existing deep learning frameworks replicates the entire model on each worker and synchronizes *all* gradients on each worker after back propagation. This design disables opportunities for overlapping data transfer in gradient synchronization with computation in back propagation. DeePa manages the gradients of each operation separately and transfers an operation's gradients as long as its back propagation is completed. We have found that this can effectively hide the data transfer overhead for gradient synchronization. As a result, the synchronous training performance matches asynchronous training in DeePa, which allows users to aggressively use synchronous training with its better algorithmic efficiency.

Distributing parameter servers. Existing frameworks use *parameter servers* to store and update variables for a CNN model. Parameter servers are usually located in CPU memory due to the limited capacity of GPU memory. Since DeePa manages the parameters for each operation separately, this allows DeePa to opportunistically distribute the parameter server onto the GPU memories whenever possible. This eliminates data transfers for operations whose gradients and parameter server are located on the same GPU and transforms all GPU to CPU copies into faster GPU to GPU copies.

6 RELATED WORK

To the best of our knowledge, DeePa is the first deep learning framework that controls and optimizes the parallelism of neural networks in all dimensions at the granularity of each operation.

Existing frameworks such as TensorFlow (Abadi et al., 2016), Caffe2 (Caf, 2016), and PyTorch (Pyt, 2017) use data parallelism to distribute the training of CNNs and only explore parallelism in the batch dimension. The standard data parallel configuration maintains a replica of the entire network on each worker, which results in large data transfers for synchronizing the gradients in each step.

Mirhoseini et al. (2017) uses model parallelism for training Inception-v3 and uses a reinforcement learning algorithm to optimize the placement of each operation on a GPU device. The learned device placement on 4 GPUs achieves 19% speedup compared to the single-GPU performance. However, parallelism in each operation is not explored.

Krizhevsky (2014) introduces "one weird trick" (OWT) that combines data parallelism with model parallelism to accelerate the distributed training of AlexNet, which efficiently reduces the data transfer cost compared to the standard data parallel configuration. However, we have found that OWT still results in suboptimal performance for AlexNet. In Section 7.2, we show that DeePa further reduces the overall data transfers by $3\times$ and the per-step training time by $2.3\times$, compared to OWT.

Goyal et al. (2017) empirically shows no loss of accuracy for training ResNet-50 on the ImageNet dataset with a large minibatch size of 8192 images. It uses the standard data parallel configuration to distribute the training onto 256 GPUs and includes a number of optimizations for reducing communication overhead. As communication is a bottleneck in distributed deep learning, we believe our techniques for reducing data transfers can substantially benefit training on large numbers of GPUs.

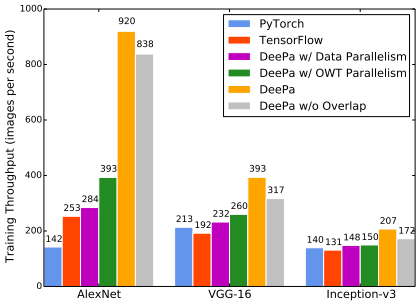


Figure 7: Training throughput (images/second) for AlexNet, VGG-16, and Inception-v3 on 16 GPUs. The purple and green bar shows the DeePa performance by restricting DeePa to use data and OWT parallelism, respectively.

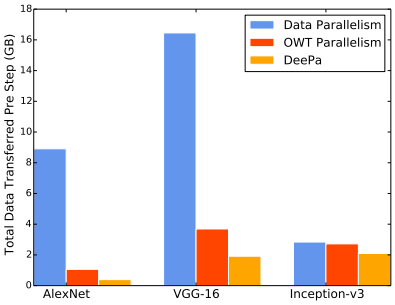


Figure 8: Total amount of data transferred in each step for training AlexNet, VGG-16, and Inception-v3 on 16 GPUs with a minibatch size of 512.

7 EVALUATION

We use AlexNet (Krizhevsky, 2014), VGG-16 (Simonyan & Zisserman, 2014), and Inception-v3 (Szegedy et al., 2016) as the benchmark CNNs and use the ImageNet dataset (Russakovsky et al., 2015) as the input. For each CNN, we compare the performance of DeePa against TensorFlow, PyTorch, and OWT. We implement OWT in DeePa by restricting all convolutional and pooling layers to use data parallelism and all fully-connected layers to use model parallelism. All experiments are performed on an XStream GPU machine (XSt, 2016), with two Intel 10-core E5-2680 Xeon processors, 256 GB main memory, and 16 NVIDIA Tesla K80 GPUs [2]. We use all 16 GPUs for training each CNN model with a minibatch size of 512 images. As a result, each GPU processes a batch of 32 images in data parallel configuration.

Figure 7 shows the training throughput for different frameworks. When enforcing DeePa to use data parallelism for all operations, DeePa achieves competitive performance compared to the best of TensorFlow and PyTorch. The OWT approach that uses model parallelism for fully-connected layers speeds up the training throughput by 1.4×, 1.2×, and 1.07×, compared to DeePa's data parallel approach. In addition, we measure the DeePa performance by using the best configuration found by the search algorithm in Section 4. DeePa achieves 6.5×, 1.9×, and 1.5× speedup, compared to TensorFlow and PyTorch.

Three main optimizations in DeePa achieve most of the performance benefit over other frameworks. First, DeePa significantly reduces data transfers in each step, as shown in Figure 8. Compared to data parallelism, the OWT approach reduces data transfers by 1.05-8.4×. However, the best configuration used by DeePa further reduces data transfers by 1.2-2.7× compared to OWT. Second, the optimization for overlapping computation with data transfers (described in Section 5) effectively hides data transfer latency and achieves better GPU utilization. The grey bars in Figure 7 illustrates DeePa's performance when the overlap optimization is disabled, which shows that overlapping computation with data transfers can improve the training throughput by 10%-30%. Third, DeePa also improves the performance by exploring parallelism in the height and width dimensions (see Section 7.3).

7.1 THE BEST CONFIGURATIONS

We describe the best configurations discovered for AlexNet, VGG-16, and Inception-v3 in Sections 7.2 to 7.4. The best configurations have several similarities.

First, for the beginning layers with large height/width dimensions and a small channel dimension, DeePa uses data parallelism on all available GPUs, since the data transfers for synchronizing gradients are much smaller than the data transfers for moving tensors between operations.

Second, deeper layers in CNNs tend to have smaller height/width dimensions and a larger channel dimension. As a result, the cost for moving tensors between different operations decreases, while the cost for synchronizing gradients increases. DeePa adaptively reduces the number of GPU workers

[2] The machine is equipped with 8 GPU cards, each of which has 2 Tesla K80 GPUs.

Table 2: The cost for different configurations for the first fully-connected of AlexNet.

Configuration	# GPU Workers	Transfer Cost	Compute Cost	Update Cost	Total Cost	Parallelism Configurations
{n=16, c=1}	16	0	+1.28	+1075	= 1076	Data Parallelism
{n=1, c=16}	16	134.4	+1.28	+0	= 135.7	OWT
{n=1, c=4}	4	33.6	+5.1	+0	= 38.7	
{n=1, c=2}	2	16.8	+10.2	+0	= **27.0**	DeePa
{n=1, c=1}	1	8.4	+20.4	+0	= 28.8	

for these layers to reduce the expensive data transfers for synchronizing gradients at the cost of introducing cheaper data transfers for moving tensors.

Third, DeePa uses model parallelism on a small number of GPU workers for fully-connected layers, because synchronizing gradients and moving tensors are both much more expensive than the compute time for fully-connected layers. DeePa reduces the data transfers for synchronizing gradients and moving tensors at the cost of using fewer GPUs.

7.2 AlexNet

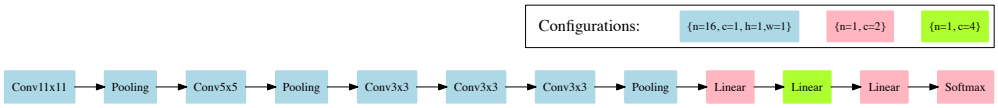


Figure 9: The global configuration for parallelizing AlexNet on 16 GPU workers.

Figure 9 shows the global configuration for AlexNet on 16 GPU workers. Note that DeePa selects the parallelism configuration that optimizes the performance for each layer. Table 2 lists the cost for different configurations of the first fully-connected layer. The standard data parallel configuration eliminates the cost for replicating the input tensors but introduces additional data transfers for synchronizing gradients. The OWT approach completely eliminates gradient synchronization at the cost of replicating the input tensors on every GPU worker. The configuration chosen by DeePa only uses 2 GPU workers for training the first fully-connected layer, which prolongs the compute time but significantly reduces the cost for both replicating input tensors and synchronizing gradients. As a result, DeePa reduces the total cost by 5× compared to other approaches.

DeePa uses data parallelism for all convolutional and pooling layers, because the additional data transfer cost introduced by transforming configurations outweighs potential performance benefits.

7.3 VGG-16

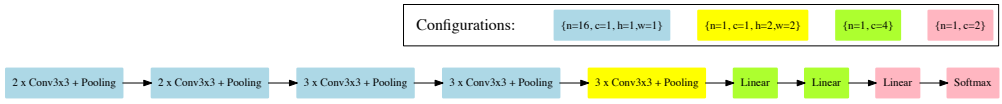


Figure 10: The global configuration for parallelizing VGG-16 on 16 GPU workers.

DeePa uses similar configurations for parallelizing the fully-connected layers in VGG-16 (Figure 10). In addition, DeePa also uses a different configuration to cooperatively accelerate the last three convolutional layers (the yellow node in Figure 10). Table 3 lists the cost for different parallelism configurations for the last three convolutional layers. The configuration with optimal total cost uses only four GPU workers for the last three convolutional layers to reduce data transfers for synchronizing gradients. DeePa also exploits parallelism in the height and width dimensions to further reduce the compute time.

7.4 Inception-v3

The Inception-v3 model has multiple Inception *modules* (Szegedy et al., 2016). Each module has several branches of convolutional and pooling layers, which are then concatenated as the output tensor of the module. Figure 11 shows the global configuration for Inception-v3. DeePa uses different configurations to parallelize different branches for the InceptionE1 module, as shown in Figure 12.

Table 3: The cost for different configurations for the last three convolutional layers of VGG-16.

Configuration	# GPU Workers	Transfer Cost	Compute Cost	Update Cost	Total Cost	Parallelism Configurations
{n=16, c=1, h=1, w=1}	16	0	+15.9	+134.4	= 150.3	Data Parallelism & OWT
{n=8, c=1, h=1, w=1}	8	39.2	+31.8	+67.2	= 138.2	
{n=4, c=1, h=1, w=1}	4	39.2	+63.7	+33.6	= 136.5	
{ n=1, c=1, h=2, w=2}	4	39.2	+54.7	+33.6	= **127.5**	DeePa
{n=2, c=1, h=1, w=1}	2	39.2	+127.4	+16.8	= 183.4	

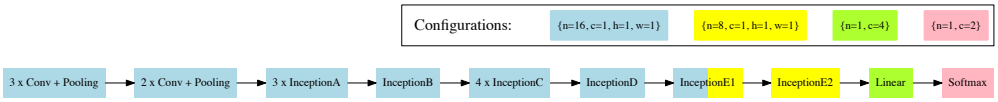


Figure 11: The global configuration for parallelizing Inception-v3 on 16 GPU workers. Each module is shown as a single node for simplicity.

We found that this configuration reduces data transfers by 30% in InceptionE1 and InceptionE2 and reduces overall data transfers by 20%.

8 Conclusion

We have presented DeePa, a deep learning framework that explores parallelism in all parallelizable dimensions to accelerate the training of CNNs. DeePa optimizes the parallelism configuration chosen at the granularity of individual layers. DeePa achieves up to 6.5× for training CNNs and reduces overall data transfers by up to 23× compared to state-of-the-art deep learning frameworks.

References

A New Lightweight, Modular, and Scalable Deep Learning Framework. https://caffe2.ai, 2016.

XStream Cray CS-Storm compute cluster. http://xstream.stanford.edu/, 2016.

Dense Linear Algebra on GPUs. https://developer.nvidia.com/cublas, 2016.

Tensors and Dynamic neural networks in Python with strong GPU acceleration. https://pytorch.org, 2017.

Martín Abadi, Paul Barham, Jianmin Chen, Zhifeng Chen, Andy Davis, Jeffrey Dean, Matthieu Devin, Sanjay Ghemawat, Geoffrey Irving, Michael Isard, et al. Tensorflow: A system for large-scale machine learning. In *OSDI*, volume 16, pp. 265–283, 2016.

Michael Bauer, Sean Treichler, Elliott Slaughter, and Alex Aiken. Legion: Expressing locality and independence with logical regions. In *Proceedings of the International Conference on High Performance Computing, Networking, Storage and Analysis*, 2012.

Sharan Chetlur, Cliff Woolley, Philippe Vandermersch, Jonathan Cohen, John Tran, Bryan Catanzaro, and Evan Shelhamer. cudnn: Efficient primitives for deep learning. *CoRR*, abs/1410.0759, 2014. URL http://arxiv.org/abs/1410.0759.

Priya Goyal, Piotr Dollár, Ross B. Girshick, Pieter Noordhuis, Lukasz Wesolowski, Aapo Kyrola, Andrew Tulloch, Yangqing Jia, and Kaiming He. Accurate, large minibatch SGD: training imagenet in 1 hour. *CoRR*, abs/1706.02677, 2017. URL http://arxiv.org/abs/1706.02677.

Alex Krizhevsky. One weird trick for parallelizing convolutional neural networks. *CoRR*, abs/1404.5997, 2014. URL http://arxiv.org/abs/1404.5997.

Azalia Mirhoseini, Hieu Pham, Quoc V Le, Benoit Steiner, Rasmus Larsen, Yuefeng Zhou, Naveen Kumar, Mohammad Norouzi, Samy Bengio, and Jeff Dean. Device placement optimization with reinforcement learning. *CoRR*, abs/1706.04972, 2017. URL http://arxiv.org/abs/1706.04972.

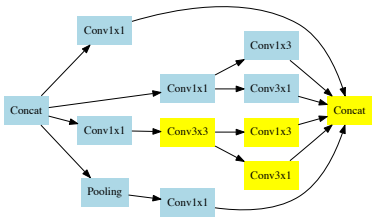


Figure 12: The configurations for parallelizing the InceptionE1 module.

Olga Russakovsky, Jia Deng, Hao Su, Jonathan Krause, Sanjeev Satheesh, Sean Ma, Zhiheng Huang, Andrej Karpathy, Aditya Khosla, Michael Bernstein, et al. Imagenet large scale visual recognition challenge. *International Journal of Computer Vision*, 2015.

Karen Simonyan and Andrew Zisserman. Very deep convolutional networks for large-scale image recognition. *CoRR*, abs/1409.1556, 2014. URL `http://arxiv.org/abs/1409.1556`.

Christian Szegedy, Wei Liu, Yangqing Jia, Pierre Sermanet, Scott E. Reed, Dragomir Anguelov, Dumitru Erhan, Vincent Vanhoucke, and Andrew Rabinovich. Going deeper with convolutions. *CoRR*, abs/1409.4842, 2014. URL `http://arxiv.org/abs/1409.4842`.

Christian Szegedy, Vincent Vanhoucke, Sergey Ioffe, Jon Shlens, and Zbigniew Wojna. Rethinking the inception architecture for computer vision. In *Proceedings of the IEEE Conference on Computer Vision and Pattern Recognition*, 2016.

Matthew D Zeiler and Rob Fergus. Visualizing and understanding convolutional networks. In *European conference on computer vision*, 2014.

A Appendix

A.1 Node and Edge elimination

We prove the correctness of the node and edge eliminations in Section 4. In particular, we prove that after applying node and edge eliminations, the modified graph has the same optimal configuration as the original graph.

A.1.1 Node elimination

For a given computation graph $G = (V, E)$, applying a node elimination on w requires w having a single in-edge $e_1 = (u, w)$ and a single out-edge $e_2 = (w, v)$. The node elimination results in a modified graph $G' = (V', E')$, where $V' = V - \{w\}$, $E' = E - e_1 - e_2 + e'$, and $e' = (u, v)$. We prove that G and G' has the same *Cost* function for the optimal global configuration g.

Theorem 1. *For the optimal global configuration g of graph $G = (V, E)$, $\mathit{Cost}(g, (V, E)) = \mathit{Cost}(g, (V', E'))$, where (V', E') is the modified graph of (V, E) after a node elimination.*

Proof. The *Cost* function is defined in Equation 1. We first compute the difference between $\mathit{Cost}(g, (V, E))$ and $\mathit{Cost}(g, (V', E'))$.

$$
\begin{aligned}
&\mathit{Cost}(g, (V, E)) - \mathit{Cost}(g, (V', E')) \\
=&\sum_{v \in V} \{v.\mathit{compute}(g(v)) + v.\mathit{update}(g(v))\} + \sum_{e=(u,v) \in E} e.\mathit{xfer}(g(u), g(v)) \\
&- \sum_{v \in V'} \{v.\mathit{compute}(g(v)) + v.\mathit{update}(g(v))\} + \sum_{e=(u,v) \in E'} e.\mathit{xfer}(g(u), g(v)) \\
=&w.\mathit{compute}(g(w)) + w.\mathit{update}(g(w)) \\
&+ e_1.\mathit{xfer}(g(u), g(w)) + e_2.\mathit{xfer}(g(w), g(v)) - e'.\mathit{xfer}(g(u), g(v))
\end{aligned} \tag{4}
$$

Since g is the optimal global configuration that minimize the *Cost* function, therefore

$$\begin{aligned} & w.\textit{compute}(g(w)) + w.\textit{update}(g(w)) + e_1.\textit{xfer}(g(u), g(w)) + e_2.\textit{xfer}(g(w), g(v)) \\ = & \min_{c_w}\{w.\textit{compute}(c_w) + w.\textit{update}(c_w) + e_1.\textit{xfer}(g(u), c_w) + e_2.\textit{xfer}(c_w, g(v))\} \end{aligned} \tag{5}$$

Combining Equation 2, 4,and 5, we get

$$\textit{Cost}(g, (V, E)) = \textit{Cost}(g, (V', E')) \tag{6}$$

□

A.1.2 EDGE ELIMINATION

For a given computation graph $G(V, E)$, applying an edge elimination on $e_1 = (u, v)$ and $e_2 = (u, v)$ results in a modified graph $G' = (V, E')$, where $E' = E - e_1 - e_2 + e'$ and $e' = (u, v)$. We prove that G and G' has the same *Cost* function for the optimal global configuration g.

Theorem 2. *For the optimal global configuration g of graph $G = (V, E)$, $\textit{Cost}(g, (V, E)) = \textit{Cost}(g, (V, E'))$, where (V, E') is the modified graph of (V, E) after an edge elimination.*

Proof. We compute the difference between $\textit{Cost}(g, (V, E))$ and $\textit{Cost}(g, (V, E'))$.

$$\begin{aligned} & \textit{Cost}(g, (V, E)) - \textit{Cost}(g, (V, E')) \\ = & e_1.\textit{xfer}(g(u), g(v)) - e_2.\textit{xfer}(g(u), g(v)) + e'.\textit{xfer}(g(u), g(v)) \\ = & 0 \end{aligned} \tag{7}$$

The last equation uses Equation 3. □

A.2 COMPLEXITY OF OUR ELIMINATION-BASED ALGORITHM

For any given computation graph $G = (V, E)$, we prove that our elimination-based algorithm can generate the final graph within $O(|V| + |E|)$ operations, with each operation either a node elimination or an edge elimination.

For any given graph $G = (V, E)$ we define a new function $\textit{Size}(G) = |V| + |E|$ that counts the number of nodes and edges in the graph. We first prove the following theorem.

Theorem 3. *For any given graph $G = (V, E)$, $\textit{Size}(G') \leq \textit{size}(G) - 1$, where $G' = (V', E')$ is the modified graph after a node or an edge elimination.*

Proof. A node elimination deletes two edges and a node from the original graph and inserts a new edge. Therefore, $\textit{Size}(G') = |V'| + |E'| = |V| - 1 + |E| - 1 < \textit{Size}(G) - 1$. A edge eliminate deletes two edges and inserts an edge. Therefore, $\textit{Size}(G') = |V'| + |E'| = |V| + |E| - 1 = \textit{Size}(G) - 1$. □

The above theorem shows that a node/edge elimination always reduces the value of the *Size* function by at least one. As a result, the total number of edge/node eliminations is smaller than $|V| + |E|$. Moreover, each elimination requires iterating over all possible parallelism configurations, which is a constant for a fixed number of workers. Therefore, we prove that our elimination-based algorithm generates the final graph within linear time complexity.

A.3 EXPERIMENT DETAILS

A.4 RELATED WORK ON OVERLAPPING COMMUNICATION WITH DATA TRANSFER

The overlap optimization in Section 5 is motivated by Goyal et al. (2017), which performs gradient aggregation in parallel with back propagation to scale *synchronous* training to large number of GPUs. We extend their design and implementation by also enabling the optimization for *asynchronous* training in DeePa.