

High Performance Machine Learning and Deep Learning with MVAPICH

Tutorial at MUG '24

by

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Outline

• Introduction

- Machine Learning
	- Distributed K-Means
	- ML Solutions
- Deep Learning
	- Deep Neural Networks
	- Distributed Deep Learning
	- DL Solutions
- Conclusion

What is Machine Learning and Deep Learning?

- Machine Learning (ML)
	- "the study of computer algorithms to improve automatically through experience and use of data"
- Deep Learning (DL) a subset of ML
	- Uses Deep Neural Networks (DNNs)
	- **Perhaps, the most revolutionary subset!**
- Based on learning data representation
- DNN Examples: Convolutional Neural Networks, Recurrent Neural Networks, Hybrid Networks
- Data Scientist or Developer Perspective for using DNNs
	- 1. Identify DL as solution to a problem
	- 2. Determine Data Set
	- 3. Select Deep Learning Algorithm to Use
	- 4. Use a large data set to train an algorithm

Courtesy: [https://hackernoon.com/difference-between-artificial-intelligence-machine-learning-and](https://hackernoon.com/difference-between-artificial-intelligence-machine-learning-and-deep-learning-1pcv3zeg)[deep-learning-1pcv3zeg](https://hackernoon.com/difference-between-artificial-intelligence-machine-learning-and-deep-learning-1pcv3zeg), [https://blog.dataiku.com/ai-vs.-machine-learning-vs.-deep-learning,](https://blog.dataiku.com/ai-vs.-machine-learning-vs.-deep-learning) https://en.wikipedia.org/wiki/Machine_learning

History: Milestones in the Development of ML/DL

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Three Main Types of Machine Learning

Courtesy: <https://bigdata-madesimple.com/machine-learning-explained-understanding-supervised-unsupervised-and-reinforcement-learning/>

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Support for Parallel and Distributed Execution

- Scikit-learn:
	- Supports execution via Joblib (<https://joblib.readthedocs.io/en/latest/>)
	- Joblib supports multi-threaded and multi-process execution (on multiple nodes)
- XGBoost:
	- Multiple ways to run on cluster of nodes:
		- Dask [\(http://dask.org](http://dask.org/))
		- Ray ([https://ray.io/\)](https://ray.io/)
		- AWS YARN
		- Apache Spark [\(https://spark.apache.org/\)](https://spark.apache.org/) using XGBoost4J-Spark
- cuML:
	- Execution is supposed on multiple nodes using Dask ([http://dask.org\)](http://dask.org/) and NVIDIA's NCCL

Parallelizing the K-means Algorithm

Step 0: Initialize centroids

- Assign initial cluster means randomly
- Step 1: Data Division
	- Distribute elements among GPUs
- Step 2: Assign elements (color)
	- Assign each element to the cluster with the closest mean
- Step 3: Update local cluster mean
	- Calculate local cluster means for all local points
- Step 4: Update global cluster mean*
	- Calculate global cluster means by calling Allreduce()
- Step 5: Repeat steps 2-4 until convergence

Set of input elements

Step 5: Repeat 2-4 until convergence

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The cuML Library: Accelerating ML on GPUs

- The NVIDIA RAPIDS project aims to build end-to-end data science analytic pipelines on GPUs
- An important component is the cuML library:
	- GPU-accelerated ML library
	- GPU-counterpart of Scikit-learn
	- Supports the execution of ML workloads on Multi-Node Multi-GPUs (MNMG) systems
- Most existing ML libraries, including Scikit-learn and Apache Spark's MLlib, only support CPU execution of ML algorithms
	- Conventional wisdom has been that only DNNs are a good match for GPUs because of high computational requirements

Main components of the cuML library

- Main components
	- Python layer
		- Provides a Scikit-learn like interface
		- Hides the complexities of the CUDA/C/C++ layer
	- Primitives and cuML algorithms built on top of CUDA
		- ML Algorithms
		- Primitives
			- Reusable building blocks for building machine learning algorithms
			- Common for different machine learning algorithms
			- Used to build different machine learning algorithms
	- Communication Support in cuML:
		- Point-to-point communication: Dask
		- Collective communication: NVIDIA Collective Communications Library (NCCL)

Accelerating cuML with MVAPICH2-GDR

- Utilize MVAPICH2-GDR (with mpi4py) as communication backend during the training phase (the fit() function) in the Multi-node Multi-GPU (MNMG) setting over cluster of GPUs
- Communication primitives:
	- Allreduce
	- Reduce
	- Broadcast
- Exploit optimized collectives

MPI4cuML 0.5 release

(http://hidl.cse.ohio-state.edu)

MPI4cuML 0.5 Release - MPI-Driven ML Training

- cuML is a distributed machine learning training framework with a focus on GPU acceleration and distributed computing. MVAPICH2-GDR provides many features to augment distributed training with cuML on GPUs
- (NEW) Based on cuML 22.02.00
	- Include ready-to-use examples for KMeans, Linear Regression, Nearest Neighbors, and tSVD
- (NEW) MVAPICH2 support for RAFT 22.02.00
	- (NEW) Enabled cuML's communication engine, RAFT, to use MVAPICH2-GDR backend for Python and C++ cuML applications
	- KMeans, PCA, tSVD, RF, LinearModels
	- Added switch between available communication backends (MVAPICH2 and NCCL)
- Built on top of mpi4py over the MVAPICH2-GDR library
- Tested with
	- Mellanox InfiniBand adapters (FDR and HDR)
	- (NEW) NVIDIA GPU A100, V100 and, P100
	- Various x86-based multi-core platforms (AMD and Intel)
- <http://hidl.cse.ohio-state.edu/>

K-Means

M. Ghazimirsaeed , Q. Anthony , A. Shafi , H. Subramoni , and D. K. Panda, Accelerating GPU-based Machine Learning in Python using MPI Library: A Case Study with MVAPICH2-GDR, MLHPC Workshop, Nov 2020

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Understanding the Deep Neural Network Concepts

• Example of a 3-layer Deep Neural Network (DNN) – (input layer is not counted)

Courtesy: <http://cs231n.github.io/neural-networks-1/>

Essential Concepts: Learning Rate (α)

Courtesy: <https://www.jeremyjordan.me/nn-learning-rate/>

Essential Concepts: Batch Size

- Batched Gradient Descent
	- Batch Size = **N**
- Stochastic Gradient Descent
	- $-$ Batch Size $= 1$
- Mini-batch Gradient Descent
	- Somewhere in the middle
	- Common:
		- **Batch Size** = 64, 128, 256, etc.
- Finding the optimal batch size will yield the fastest learning.

Courtesy: <https://www.jeremyjordan.me/gradient-descent/>

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The Need for Parallel and Distributed Training

- Why do we need Parallel Training?
- Larger and Deeper models are being proposed
	- **Language Models: RNNs -> Transformers -> BERT – GPT – LLaMA**
	- **Vision Models: AlexNet** -> **ResNet** -> **NASNet – AmoebaNet Vision Transformers**
	- DNNs require a lot of memory and a lot of computation
	- Larger models cannot fit a GPU's memory
- Single GPU training cannot keep up with ever-larger models
- Community has moved to multi-GPU training
- Multi-GPU in one node is good but there is a limit to Scale-up (8-16 GPUs)
- **Multi-node (Distributed or Parallel) Training is necessary!!**

Parallelization Strategies

- Some parallelization strategies..
	- Data Parallelism or Model Parallelism
	- Hybrid Parallelism

Machine 2

Machine 4

Hybrid (Model and Data) Parallelism

Data Parallelism

Machine 1

Machine 3

Courtesy: <http://engineering.skymind.io/distributed-deep-learning-part-1-an-introduction-to-distributed-training-of-neural-networks>

 $\frac{1}{1}$ GPU

GPU

GPU3

Machine 1

Data Parallelism and MPI Collectives

- Step1: Data Propagation
	- Distribute the Data among GPUs
- Step2: Forward Backward Pass
	- Perform forward pass and calculate the prediction
	- Calculate Error by comparing prediction with actual output
	- Perform backward pass and calculate gradients
- Step3: Gradient Aggregation
	- Call MPI_Allreduce to reduce the local gradients
	- Update parameters locally using global gradients

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MVAPICH2 (MPI)-driven Infrastructure for ML/DL Training

More details available from: http://hidl.cse.ohio-state.edu

HiDL Software Stack Release v1.0

- Based on Horovod
-
- Optimized support for MPI controller in deep learning workloads
Efficient large-message collectives (e.g. Allreduce) on various • Efficient large-message collectives (e.g. Allreduce) on various CPUs and GPUs
- GPU-Direct algorithms for collective operations (including those commonly used for data- and model-parallelism, e.g. Allgather and Alltoall)
- Support for fork safety
- Exploits efficient large message collectives
- Compatible with
	- Mellanox InfiniBand adapters (EDR, FDR, HDR)
	- Various x86-based multi-core CPUs (AMD and Intel)
	- NVIDIA A100, V100, P100, Quadro RTX 5000 GPUs
	- CUDA [9.x, 10.x, 11.x] and cuDNN [7.5.x, 7.6.x, 8.0.x, 8.2.x, 8.4.x]
	- AMD MI100 GPUs
	- ROCm [5.1.x]

Install Horovod with MVAPICH2-X and MVAPICH2-GDR

Command to install Horovod for CPU

\$ HOROVOD_WITH_MPI=1 pip install --no-cache-dir horovod

Command to install Horovod for GPU

\$ HOROVOD_GPU_ALLREDUCE=MPI HOROVOD_CUDA_HOME=/opt/cuda/11.3 HOROVOD_WITH_MPI=1 pip install --no-cache-dir horovod

Run PyTorch on a single GPU

+ python pytorch_synthetic_benchmark.py --batch-size 64 --num-iters=5

. Model: resnet50 Batch size: 64 Number of GPUs: 1 Running warmup... Running benchmark... Iter #0: 333.9 img/sec per GPU Iter #1: 334.2 img/sec per GPU Iter #2: 333.9 img/sec per GPU Iter #3: 333.8 img/sec per GPU Iter #4: 333.9 img/sec per GPU Img/sec per GPU: 334.0 +-0.2 --- Total img/sec on 1 GPU(s): 334.0 +-0.2 --- V100

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Run PyTorch on two nodes with 1 GPU/node (using MVAPICH2- GDR)

+ mpirun_rsh -np 2 gpu11 gpu12 MV2_USE_CUDA=1 MV2_CPU_BINDING_POLICY=hybrid MV2 HYBRID BINDING POLICY=spread MV2 USE_RDMA_CM=0 MV2 GPUDIRECT GDRCOPY LIB=/opt/gdrcopy2.0/lib64/libgdrapi.so LD PRELOAD=mv2-gdr/lib/libmpi.so python pytorch synthetic benchmark.py --batch-size 64 --num-iters=5

. Model: resnet50 Batch size: 64 Number of GPUs: 2 Running warmup... Running benchmark... Iter #0: 317.0 img/sec per GPU Iter #1: 314.9 img/sec per GPU Iter #2: 315.4 img/sec per GPU Iter #3: 318.0 img/sec per GPU Iter #4: 316.7 img/sec per GPU Img/sec per GPU: 316.4 +-2.2 --- Total img/sec on 2 GPU(s): 632.8 +-4.3 --- V100 **~1.89X on 2 GPUs**

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Solutions and Case Studies: Exploiting HPC for DL

- **Data Parallelism**
- Model-Parallelism

Distributed TensorFlow on ORNL Summit (1,536 GPUs)

- ResNet-50 Training using TensorFlow benchmark on SUMMIT -- 1536 Volta GPUs!
- 1,281,167 (1.2 mil.) images

Image per second

Image per second

Thousands

Thousands

- $Time/epoch = 3 seconds$
- Total Time (90 epochs) = 3 x 90 = 270 seconds = **4.5 minutes!**

0 50 100 150 200 250 300 350 400 450 1 2 4 6 12 24 48 96 192 384 768 1536 Number of GPUs MVAPICH2-GDR 2.3.4 **MVAPICH2-GDR 2.3.4** *MVAPICH2-GDR reaching ~0.42 million images per second for ImageNet-1k! ImageNet-1k has 1.2 million images*

Platform: The Summit Supercomputer (#2 on Top500.org) – 6 NVIDIA Volta GPUs per node connected with NVLink, CUDA 10.1

Distributed TensorFlow on TACC Frontera (2048 CPU nodes)

- Scaled TensorFlow to 2048 nodes on Frontera using MVAPICH2 and IntelMPI
- MVAPICH2 delivers close to the ideal performance for DNN training
- Report a peak of 260,000 images/sec on 2048 nodes
- On 2048 nodes, ResNet-50 can be trained in 7 minutes!

AccDP: Exploiting Data Parallelism

ResNet18 training throughput comparison between regular training and AccDP (proposed design) for different DNN models on up to 8 nodes 2 GPUs per node (16 GPUs) with 4 MPS clients per GPU

Multi node with ResNet18 Multi node with ShuffleNet

• ShuffleNet training throughput comparison between regular training and AccDP (proposed design) for different DNN models on up to 8 nodes 2 GPUs per node (16 GPUs) with 4 MPS clients per GPU.

N. Alnaasan, A. Jain, A. Shafi, H. Subramoni, and DK Panda, "AccDP: Accelerated Data-Parallel Distributed DNN Training for Modern GPU-Based HPC Clusters", HiPC'22.

Solutions and Case Studies: Exploiting HPC for DL

- **Data Parallelism**
- **Model-Parallelism**

Exploiting Model Parallelism in AI-Driven Digital Pathology

- Pathology whole slide image (WSI)
	- $-$ Each WSI = 100,000 x 100,000 pixels
	- Can not fit in a single GPU memory
	- Tiles are extracted to make training possible
- Two main problems with tiles
	- Restricted tile size because of GPU memory limitation
	- Smaller tiles loose structural information
- Reduced training time significantly
	- **GEMS-Basic: 7.25 hours (1 node, 4 GPUs)**
	- **GEMS-MAST: 6.28 hours (1 node, 4 GPUs)**
	- **GEMS-MASTER: 4.21 hours (1 node, 4 GPUs)**
	- **GEMS-Hybrid: 27 mins (32 nodes, 128 GPUs)**
	- **Overall 15x reduction in training time!!!!**

A. Jain, A. Awan, A. Aljuhani, J. Hashmi, Q. Anthony, H. Subramoni, D. K. Panda, R. Machiraju, and A. Parwani, "GEMS: GPU Enabled Memory Aware Model Parallelism System for Distributed DNN Training", Supercomputing (SC '20).

Courtesy: [https://blog.kitware.com/digital-slide](https://blog.kitware.com/digital-slide-archive-large-image-and-histomicstk-open-source-informatics-tools-for-management-visualization-and-analysis-of-digital-histopathology-data/)[archive-large-image-and-histomicstk-open-source](https://blog.kitware.com/digital-slide-archive-large-image-and-histomicstk-open-source-informatics-tools-for-management-visualization-and-analysis-of-digital-histopathology-data/)[informatics-tools-for-management-visualization-and](https://blog.kitware.com/digital-slide-archive-large-image-and-histomicstk-open-source-informatics-tools-for-management-visualization-and-analysis-of-digital-histopathology-data/)[analysis-of-digital-histopathology-data/](https://blog.kitware.com/digital-slide-archive-large-image-and-histomicstk-open-source-informatics-tools-for-management-visualization-and-analysis-of-digital-histopathology-data/)

Scaling ResNet110 v2 on 1024×1024 image tiles using histopathology data

MPI4DL v0.6

[MPI4DL v0.6](https://github.com/OSU-Nowlab/MPI4DL) is a distributed, accelerated and memory efficient training framework for very high-resolution images that integrates Spatial Parallelism, Bidirectional Parallelism, Layer Parallelism, and Pipeline Parallelism.

Features:

- Based on PyTorch
- Support for training very high-resolution images
- Distributed training support for:
	- Model Parallelism
- Spatial Parallelism for High Resolution Images
- Layer Parallelism (LP)

- Pipeline Parallelism (PP)

- Spatial and Layer Parallelism (SP+LP) - Spatial and Pipeline Parallelism (SP+PP)

- Memory Efficient Bidirectional Parallelism (GEMS)
	- Bidirectional and Layer Parallelism (GEMS+LP)
	- Bidirectional and Pipeline Parallelism (GEMS+PP)
	- Spatial, Bidirectional and Layer Parallelism (SP+GEMS+LP)
	- Spatial, Bidirectional and Pipeline Parallelism (SP+GEMS+PP)

- Support for different image sizes and custom datasets.
- Exploits collective features of MVAPICH2-GDR

Throughput comparison of Pipeline Parallelism and Pipeline + Spatial Parallelism techniques for AmoebaNet on 1024 * 1024 and 2048 * 2048 image sizes.

Throughput comparison of Spatial Parallelism and Spatial + Bidirectional Parallelism for AmoebaNet and ResNet with the following configurations: 5 model splits,4 spatial parts, and 2 model replicas for Bidirectional Parallelism.

ParaInfer-X v1.0: a Temporal Fusion framework for LLM inference

ParaInfer-X is a collection of parallel inference techniques that can facilitate the deployment of emerging AI models on edge devices and HPC clusters. In v1.0, we proposes a temporal fusion framework, named Flover, to smartly batch multiple requests during LLM generation, which is also known as temporal fusion/in-flight batching. **Features:**

- Based on Faster Transformer
- Support for inference of various large language models:
	- GPT-J 6B, LlaMA 7B, 13B**,** 33B**,** 65B
	- Support for persistent model inference stream
	- Support for temporal fusion/in-flight batching of multiple requests
	- Support for multiple GPU tensor parallelism
	- Support for asynchronous memory reordering for evicting finished requests

(a). Parallel inference on different models. We measure the overall time spent on parallel inference $1, 2, 4, 8, 16, 32$ requests.

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Conclusion

- Exponential growth in Machine Learning and Deep Learning frameworks
- Provided an overview of issues, challenges, and opportunities for designing efficient communication runtimes
	- Efficient, scalable, and hierarchical designs are crucial for ML and DL frameworks
	- Co-design of communication runtimes and ML and DL frameworks will be essential
- Presented use-cases to demonstrate the complex interaction between DL and ML middleware with the underling HPC technologies and middleware
- Need collaborative efforts to achieve the full potential

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Network-Based Computing Laboratory http://nowlab.cse.ohio-state.edu/

The High-Performance MPI/PGAS Project http://mvapich.cse.ohio-state.edu/

High-Performance **Big Data**

The High-Performance Big Data Project http://hibd.cse.ohio-state.edu/

The High-Performance Deep Learning Project http://hidl.cse.ohio-state.edu/