





# 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:** <u>https://hackernoon.com/difference-between-artificial-intelligence-machine-learning-and-deep-learning-1pcv3zeg</u>, <u>https://blog.dataiku.com/ai-vs.-machine-learning-vs.-deep-learning</u>, https://en.wikipedia.org/wiki/Machine\_learning

Feature extraction + Classification

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Input

Output

#### **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/

#### NVIDIA's NCCL **Network Based Computing Laboratory**

# **Support for Parallel and Distributed Execution**

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









- Execution is supposed on multiple nodes using Dask (<u>http://dask.org</u>) and

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# **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



#### 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)
- <a href="http://hidl.cse.ohio-state.edu/">http://hidl.cse.ohio-state.edu/</a>

#### 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:** <u>http://cs231n.github.io/neural-networks-1/</u>

#### **Essential Concepts: Learning Rate (\alpha)**



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

#### **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:
    - **<u>Batch Size</u>** = 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: <a href="http://engineering.skymind.io/distributed-deep-learning-part-1-an-introduction-to-distributed-training-of-neural-networks">http://engineering.skymind.io/distributed-deep-learning-part-1-an-introduction-to-distributed-training-of-neural-networks</a>

GPU

GPU 3

Machine 1

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#### **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 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

V100 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):  $334.0 \pm 0.2$ 

# 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

V100 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 + -4.3~1.89X on 2 GPUs

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



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

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#### **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**

#### Multi node with ResNet18

 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 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**

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# **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: <u>https://blog.kitware.com/digital-slide-</u> <u>archive-large-image-and-histomicstk-open-source-</u> <u>informatics-tools-for-management-visualization-and-</u> <u>analysis-of-digital-histopathology-data/</u>



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

#### **MPI4DL v0.6**

<u>MPI4DL v0.6</u> 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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The High-Performance MPI/PGAS Project http://mvapich.cse.ohio-state.edu/



High-Performance Big Data

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



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