

# **MPI4Dask 0.3 User Guide**

HIGH-PERFORMANCE BIG DATA TEAM  
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# 1 Overview of the MPI4Dask Project

MPI4Dask is a custom version of the Dask Distributed library that contains high-performance MPI communication backend for Dask. The MPI backend in MPI4Dask uses mpi4py over an MVAPICH2 library and targets modern HPC clusters. The current version of MPI4Dask supports both CPU and GPU-based Dask applications. MPI4Dask embodies enhancements to the Dask Distributed package.

MPI4Dask provides point-to-point asynchronous I/O communication *coroutines*, which are non-blocking concurrent operations defined using the `async/await` keywords from the Python's `asyncio` framework. If there are any questions, comments or feedback regarding this software package, please post them to ([mvapich-discuss@lists.osu.edu](mailto:mvapich-discuss@lists.osu.edu)).

## 2 Features

High-level features of MPI4Dask are listed below. New features and enhancements are marked as **(NEW)**.

- **(NEW)** Based on Dask Distributed 2022.8.1
- Compliant with user-level Dask APIs and packages
- Support for MPI-based communication for GPU-based Dask applications
- Implements point-to-point communication co-routines
- Efficient chunking mechanism implemented for large messages
- Support for MPI-based communication for CPU-based Dask applications
- Built on top of mpi4py over MVAPICH2, MVAPICH2-X, and MVAPICH2-GDR libraries
- Supports starting execution of Dask programs using Dask-MPI
- Tested with
  - CPU-based Dask applications using numPy and Pandas data frames
  - GPU-based Dask applications using cuPy and cuDF
  - Mellanox InfiniBand adapters (FDR and EDR)
  - Various multi-core platforms
  - **(NEW)** Scaling out to hundreds of nodes
  - **(NEW)** NVIDIA V100 and A100 GPUs

## 3 Setup Instructions

### 3.1 Installation Pre-requisites

#### 3.1.1 Miniconda

1. Install [Miniconda](#), which is a free minimal installer for the conda package manager. The following commands download and install latest version of Miniconda3 for Python 3.x:

```
$ wget https://repo.anaconda.com/miniconda/Miniconda3-latest-Linux-x86_64.sh
$ chmod a+x Miniconda3-latest-Linux-x86_64.sh
$ ./Miniconda3-latest-Linux-x86_64.sh
```

1. Create a directory for MPI4Dask.

```
$ mkdir root-directory/mpi4dask
```

#### 3.1.2 Modules/Libraries for GPU-based Dask Applications

1. Make sure that relevant/required modules are loaded in the environment. One of the required module is the CUDA toolkit (11.2 in this case). A list of all installed modules on the system can be seen as follows:

```
$ module list
Currently Loaded Modulefiles:
  1) cuda/11.2
```

2. Install pre-requisite libraries including build tools, Python (version 3.8 here), Cython, CUDA toolkit (version 11.2 here), Dask packages, RAPIDS libraries, and array packages. Users may install specific versions of these packages. To avoid potential conda environment compatibility check issues, we recommend using the following commands to install:

```
$ cd root-directory/mpi4dask
$ git clone https://github.com/aamirshafi/distributed.git
$ cd distributed
All required packages and configuration are included in mpi4dask_cuda.yml.
Default env name is mpi4dask_cuda, which can be modified in the .yml 'name' space.
$ conda env create -f mpi4dask_cuda.yml
```

#### 3.1.3 Modules/Libraries for CPU-based Dask Applications

1. Install pre-requisite libraries including build tools, Python (version 3.8 here), Cython, Dask packages, and array packages. Users may install specific versions of these packages. Also, these packages are installed from three conda channels, including conda-forge, rapidsai, and nvidia, as follows:

```
$ cd root-directory/mpi4dask
$ git clone https://github.com/aamirshafi/distributed.git
$ cd distributed
All required packages and configuration are included in mpi4dask.yml.
Default env name is mpi4dask, which can be modified in the .yml 'name' space.
$ conda env create -f mpi4dask.yml
```

### 3.1.4 Install the MVAPICH2 Library

1. Download and install an appropriate flavor of the MVAPICH2 library from <http://mvapich.cse.ohio-state.edu>. For running GPU-based Dask applications, the MVAPICH2-GDR library is recommended. For CPU-based Dask applications, either the MVAPICH2 or the MVAPICH2-X library can be used. The MVAPICH2-X software provides advanced features and better performance and, hence, is recommended for CPU-based Dask applications. Please become familiar with advanced configurations of the MVAPICH2 software—this is important to get maximum performance from the hardware system.

MVAPICH2 user-guides can be seen here:

- MVAPICH2: <https://mvapich.cse.ohio-state.edu/static/media/mvapich/mvapich2-quickstart.pdf>
- MVAPICH2-GDR: <https://mvapich.cse.ohio-state.edu/userguide/gdr/>
- MVAPICH2-X: <https://mvapich.cse.ohio-state.edu/static/media/mvapich/mvapich2-x-userguide.pdf>

2. This guide assumes that an environment variable \$MPILIB points to the installation directory of the MVAPICH2 software.

```
$ export MPILIB=/path/to/MVAPICH2-GDR/install/directory
```

3. It is important to make sure that the MVAPICH2-GDR is the MPI library used with MPI4Dask. Other installations of MPI libraries can result in errors. In order to avoid such situations, it is important to make sure that the correct version of the MVAPICH2-GDR library is available on the \$PATH variable. This can be checked as follows:

```
$ $MPILIB/bin/mpiname -a
```

If the \$PATH is set correctly, the output of the following command should match the previous command output.

```
$ mpiname -a
```

### 3.1.5 Install the mpi4py Library

1. Download and install the mpi4py library. Details about mpi4py can be seen at <https://mpi4py.readthedocs.io>. Download the software as follows:

```
$ cd root-directory/mpi4dask  
  
$ git clone https://github.com/mpi4py/mpi4py.git  
  
$ cd mpi4py
```

2. As part of the installation, the configuration file (`mpi.cfg`) in the root directory of `mpi4py` software needs to be updated to point to the `MVAPICH2` library. It is recommended to make a new section—in the `mpi.cfg` file—for `MVAPICH2` as follows:

```
# MVAPICH2  
# -----  
[MVAPICH2]  
mpi_dir = /path/to/MVAPICH2-GDR/install/directory  
mpicc = %(mpi_dir)s/bin/mpicc  
mpicxx = %(mpi_dir)s/bin/mpicxx  
include_dirs = %(mpi_dir)s/include  
library_dirs = %(mpi_dir)s/lib64  
runtime_library_dirs = %(library_dirs)s
```

3. After updating the `mpi.cfg` configuration file, install `mpi4py` with the following commands:

```
$ python setup.py build --mpi=MVAPICH2  
  
$ pip install .
```

### 3.1.6 Install the `dask-mpi` library

Download and Install `dask-mpi` package from the github repository. `MPI4Dask` relies on `Dask-MPI` to start execution of `Dask` scheduler, client, and workers. Details about `Dask-MPI` package can be seen at <http://mpi.dask.org>. It is recommended not to install `Dask-MPI` through the `conda` package manager as it installs some conflicting packages as dependencies.

```
$ cd root-directory/mpi4dask  
  
$ git clone https://github.com/dask/dask-mpi.git  
  
$ cd dask-mpi  
  
$ python setup.py build  
  
$ pip install .
```

## 3.2 Install `MPI4Dask` distributed

As part of this step, an enhanced version of the `Dask Distributed` library with support for `MPI` communication-backend is installed using the following commands:

1. Download the integrated `MPI4Dask` distributed:

```
$ cd root-directory/mpi4dask
$ git clone https://github.com/aamirshafi/distributed.git
```

2. Change directory to the custom Dask Distributed library:

```
$ cd distributed
```

3. Install the custom Dask Distributed package with the following commands:

```
$ python setup.py build
$ pip install .
```

4. At this time, MPI4Dask with all its dependencies should be installed. Verify this by running the following commands:

```
$ conda list
$ conda list | grep distributed
$ conda list | grep dask
```

## 4 Basic Usage Instructions

The MPI4Dask package contains two sample applications that can be used to test and benchmark performance of MPI4Dask. These applications include: 1) Sum of cuPy Array and its Transpose, and 2) cuDF Merge.

Notice that we highly recommend users always do `import mpi4py` before `import dask` or `import distributed`. This is due to the `Dask-distributed` will interfere with the initialization of MPI communication after a certain scale. An example is:

```
import mpi4py
from mpi4py import MPI
import mpi4py.rc
mpi4py.rc.threads = True
from dask_mpi import initialize
import dask
from dask.distributed import *
```

And specifically, when users call `initialize()` function from `dask-mpi`, please do the following to make sure that MPI Comm world is created before initializing dask scheduler/workers.

```
initialize(
    interface='ib0',
    protocol='mpi',
    nthreads=8, # user can specify this accordingly
    comm=MPI.COMM_WORLD, # MPI comm has to be created first
    local_directory=os.environ.get("TMPDIR", None)
)
```

## 4.1 Sum of cuPy Array and its Tranpose

This benchmark creates a cuPy array and distributes its chunks across Dask workers. The benchmark adds these distributed chunks to their transpose, forcing the GPU data to move around over the network. This benchmark is originally taken from the dask-cuda package repository [https://github.com/rapidsai/dask-cuda/tree/branch-0.18/dask\\_cuda/benchmarks](https://github.com/rapidsai/dask-cuda/tree/branch-0.18/dask_cuda/benchmarks).

Write the hosts files. This file contains names of the compute nodes, with GPUs, that will execute the parallel Dask program.

```
$ cd root-directory/mpi4dask/distributed/dask_apps

$ vim hosts
.. write name of the compute nodes ..

$ cat hosts
machine1
machine2
machine3
machine4
```

Note that users need to customize the hosts file as per their environment. This benchmark can be executed as follows:

```
$ LD_PRELOAD=$MPILIB/lib/libmpi.so $MPILIB/bin/mpirun_rsh -export-all -np 4 -hostfile
  hosts MV2_USE_CUDA=1 MV2_USE_GDRCOPY=1
  MV2_GPUDIRECT_GDRCOPY_LIB=/opt/gdrcopy2.0/lib64/libgdrapi.so
  MV2_CPU_BINDING_LEVEL=SOCKET MV2_CPU_BINDING_POLICY=SCATTER python cupy_sum_mpi.py
```

This command is starting the execution of the Dask application using MPI—this is enabled by the Dask-MPI package. A total of 4 MPI processes are started with process 0 and 1 assuming the role of Dask scheduler and client respectively. All other processes with ranks greater than 1 becoming Dask workers. Here processes 2 and 3 are Dask workers. The LD\_PRELOAD environment variable is required for Python applications to run correctly with the MVAPICH2-GDR library.

Details on various options, passed to the MVAPICH2-GDR library, are as follows:

- `-np 4` — specifies the total number of parallel processes started by the MPI library
- `-hostfile hosts` — specifies the names of machines where to start parallel processes in the “hosts” file
- `MV2_USE_CUDA=1` — ensures that MVAPICH2-GDR library supports GPU-to-GPU communication
- `MV2_USE_GDRCOPY=1` — turns on the GDRCopy protocol
- `MV2_GPUDIRECT_GDRCOPY_LIB` — specifies path to the GDRCopy dynamic library
- `MV2_CPU_BINDING_LEVEL=SOCKET MV2_CPU_BINDING_POLICY=SCATTER` — selects a socket-level scatter process binding policy

If users are deploying MPI4Dask on nodes that each have multiple GPUs, please specify the CPU mapping explicitly to improve the performance and avoid potential issues. The standard procedure is as follows:



### 1. Check the Nvidia device topology.

```
$ nvidia-smi topo -m
$
    GPU0 GPU1 mlx5_0 mlx5_1 CPU Affinity NUMA Affinity
GPU0 X SYS NODE SYS 0-63 0
GPU1 SYS X SYS NODE 64-127 1
mlx5_0 NODE SYS X SYS
mlx5_1 SYS NODE SYS X

Legend:

X = Self
SYS = Connection traversing PCIe as well as the SMP interconnect between NUMA
    nodes (e.g., QPI/UPI)
NODE = Connection traversing PCIe as well as the interconnect between PCIe Host
    Bridges within a NUMA node
PHB = Connection traversing PCIe as well as a PCIe Host Bridge (typically the
    CPU)
PXB = Connection traversing multiple PCIe bridges (without traversing the PCIe
    Host Bridge)
PIX = Connection traversing at most a single PCIe bridge
NV# = Connection traversing a bonded set of # NVLinks
```

2. For example, the above info is about a node with 2 Nvidia A100 GPUs, with 2 AMD EPYC 7713 64-Core Processors (w. 128 HyperThreading). Check the 'CPU Affinity' column, 'GPU0' is mapped to Core 0-63, 'GPU1' is mapped to Core 64-127.

3. If users launch two dask-workers (two MPI processes) on this node, first, remove two parameters from the above command

```
MV2_CPU_BINDING_LEVEL=SOCKET MV2_CPU_BINDING_POLICY=SCATTER
```

4. And specify three MPI runtime parameters. Here is an example where each dask-worker uses 8 threads.

```
$ OMP_NUM_THREADS=8 MV2_THREADS_PER_PROCESS=8
    MV2_CPU_MAPPING=0,1,2,3,4,5,6,7:64,65,66,67,68,69,70,71
```

5. Therefore, the overall command to run will be:

```
$ LD_PRELOAD=$MPILIB/lib/libmpi.so $MPILIB/bin/mpirun_rsh -export-all -np 4
    -hostfile hosts MV2_USE_CUDA=1 MV2_USE_GDRCOPY=1
    MV2_GPUDIRECT_GDRCOPY_LIB=/opt/gdrcopy2.0/lib64/libgdrapi.so
    OMP_NUM_THREADS=8 MV2_THREADS_PER_PROCESS=8
    MV2_CPU_MAPPING=0,1,2,3,4,5,6,7:64,65,66,67,68,69,70,71 python cupy_sum_mpi.py
```

This is not an exhaustive list of options that can be passed to the MVAPICH2-GDR library. Details on these—and other flags for the MVAPICH2-GDR library—can be seen at: <http://mvapich.cse.ohio-state.edu/userguide/gdr/>. As mentioned earlier, it is important to configure the MVAPICH2-GDR installation correctly to ensure optimal performance.

The `cupy_sum_mpi.py` script has some configuration parameters that can be modified, in the source-code, by the users. These include:

- `DASK_PROTOCOL = 'mpi'` — specifies the protocol used by the Dask Distributed library. Options include 'mpi', 'tcp', 'ucx'
- `DASK_INTERFACE = 'ib0'` — specifies the interface used by the Dask Distributed library
- `GPUS_PER_NODE = 1` — specifies the number of GPUs in the system
- `THREADS_PER_NODE = 8` — specifies the number of threads (cores) in the system

## 4.2 cuDF Merge

cuDF DataFrames are table-like data-structure that are stored in the GPU memory. As part of this application, the merge operation is carried out for multiple cuDF data frames. This benchmark is originally taken from the dask-cuda package repository [https://github.com/rapidsai/dask-cuda/tree/branch-0.18/dask\\_cuda/benchmarks](https://github.com/rapidsai/dask-cuda/tree/branch-0.18/dask_cuda/benchmarks). This application can be executed as follows:

```
$ cd root-directory/mpi4dask/distributed/dask_apps

$ LD_PRELOAD=$MPILIB/lib/libmpi.so $MPILIB/bin/mpirun_rsh -export-all -np 4 -hostfile
  hosts MV2_USE_CUDA=1 MV2_USE_GDRCOPY=1
  MV2_GPU_DIRECT_GDRCOPY_LIB=/opt/gdrcopy2.0/lib64/libgdrapi.so
  MV2_CPU_BINDING_LEVEL=SOCKET MV2_CPU_BINDING_POLICY=SCATTER python
  cudf_merge_mpi.py
```

The `cudf_merge_mpi.py` script accepts some runtime arguments. Some of the important ones include:

- `--type gpu` — specifies that the data frame is GPU-based. Currently only GPU has been tested
- `--protocol mpi` — specifies the communication backend used by Dask Distributed. Options include 'mpi', 'tcp', 'ucx'
- `--runs 20` — specifies the number of repetitions for the experiment
- `--chunk-size 1.000.000.00` — specifies the chunk size for the data frame

In addition, the `cudf_merge_mpi.py` script has some configuration parameters that can be modified by the users. These include:

- `DASK_INTERFACE = 'ib0'` — specifies the interface used by the Dask Distributed library
- `GPUS_PER_NODE = 1` — specifies the number of GPUs in the system
- `THREADS_PER_NODE = 8` — specifies the number of threads (cores) in the system

If users are deploying MPI4Dask on nodes that each have multiple GPUs, please specify the CPU mapping explicitly to improve the performance and avoid potential issues. Please follow the same standard procedure as in §4.1.

## **5 Troubleshooting with MPI4Dask**

If you are experiencing any problems with MPI4Dask, please feel free to contact us by sending an email to ([mvapich-discuss@lists.osu.edu](mailto:mvapich-discuss@lists.osu.edu)).