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# **h5bench**

***Release 1.2***

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## GETTING STARTED

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h5bench is a suite of parallel I/O benchmarks or kernels representing I/O patterns that are commonly used in HDF5 applications on high performance computing systems. H5bench measures I/O performance from various aspects, including the I/O overhead, observed I/O rate, etc. You can check h5bench source-code at <https://github.com/hpc-io/h5bench>.

These are the benchmarks and kernels currently available in h5bench:

Benchmark	Name	SYNC	ASYNC	CACHE	LOG
h5bench write	h5bench_write				
h5bench read	h5bench_read				
Metadata Stress	h5bench_hdf5_iotest				
AMReX	h5bench_amrex				
Exerciser	h5bench_exerciser				
OpenPMD (write)	h5bench_openpmd_write				
OpenPMD (read)	h5bench_openpmd_read				
E3SM-IO	h5bench_e3sm				
MACSio	h5bench_macsisio				



## BUILD INSTRUCTIONS

### 1.1 Build with CMake (recommended)

First, clone the h5bench GitHub repository and ensure you are cloning the submodules:

```
git clone --recurse-submodules https://github.com/hpc-io/h5bench
```

If you are updating your h5bench, ensure you have the latest submodules that could be included in new releases:

```
git submodule update --init
```

#### 1.1.1 Dependency and environment variable settings

H5bench depends on MPI and parallel HDF5.

##### Use system provided by HDF5

For instance on the Cori system at NERSC:

```
module load cray-hdf5-parallel
```

You can also load any parallel HDF5 provided on your system, and you are good to go.

##### Use your own installed HDF5

Make sure to unload any system provided HDF5 version, and set an environment variable to specify the HDF5 install path:

```
export HDF5_HOME=/path/to/your/hdf5/installation
```

It should point to a path that contains the `include/`, `lib/`, and `bin/` subdirectories.

### 1.1.2 Compile with CMake

In the source directory of your cloned h5bench repository, run the following:

```
mkdir build
cd build

cmake ..

make
make install
```

**Warning:** If you plan on calling `make install` please notice that the default behavior of CMake is to install it system-wide. You can change the destination installation folder by passing the `-DCMAKE_INSTALL_PREFIX=<path>` to override with your defined installation directory.

By default, h5bench will only compile the base write and read benchmarks. To enable the additional benchmarks, you need to explicitly enable them before building h5bench. You can also enable all the benchmarks with `-DH5BENCH_ALL=ON`. Notice that some of them have additional dependencies.

Benchmark	Name	Build
h5bench write	h5bench_write	Always
h5bench read	h5bench_read	Always
Metadata Stress	h5bench_hdf5_iotest	<code>-DH5BENCH_METADATA=ON</code>
AMReX	h5bench_amrex	<code>-DH5BENCH_AMREX=ON</code>
Exerciser	h5bench_exerciser	<code>-DH5BENCH_EXERCISER=ON</code>
OpenPMD (write)	h5bench_openpmd_write	<code>-DH5BENCH_OPENPMD=ON</code>
OpenPMD (read)	h5bench_openpmd_read	<code>-DH5BENCH_OPENPMD=ON</code>
E3SM-IO	h5bench_e3sm	<code>-DH5BENCH_E3SM=ON</code>
MACSio	h5bench_macsio	<code>-DH5BENCH_MACSIO=ON</code>

**Warning:** If you want to specify the installation directory, you can pass `-DCMAKE_INSTALL_PREFIX` to `cmake`. If you are not installing it, make sure when you run h5bench, you update your environment variables to include the *build* directory. Otherwise, h5bench will not be able to find all the benchmarks.

### 1.1.3 Build with HDF5 ASYNC VOL connector support

To run `_async` benchmarks, you need the develop branch of **both** HDF5 and ASYNC-VOL. When building h5bench you need to specify the `-DWITH_ASYNC_VOL:BOOL=ON` option and have already compiled the VOL connector in the `$ASYNC_VOL` directory:

```
mkdir build
cd build

cmake .. -DWITH_ASYNC_VOL=ON -DCMAKE_C_FLAGS="-I/$ASYNC_VOL/src -L/$ASYNC_VOL/src"

make
make install
```



h5bench will automatically set the environment variables required to run the asynchronous versions, as long as you specify them in your JSON configuration file. However, if you run the benchmarks manually, you will need to set the following environment variables:

```
export HDF5_HOME="$YOUR_HDF5_DEVELOP_BRANCH_BUILD/hdf5"
export ASYNC_HOME="$YOUR_ASYNC_VOL/src"

export HDF5_VOL_CONNECTOR="async under_vol=0;under_info={}"
export HDF5_PLUGIN_PATH="$ASYNC_HOME"

# Linux
export LD_LIBRARY_PATH="$HDF5_HOME/lib:$ASYNC_HOME"
# MacOS
export DYLD_LIBRARY_PATH="$HDF5_HOME/lib:$ASYNC_HOME"
```

## 1.2 Build with Spack

You can also use Spack to install h5bench:

```
spack install h5bench
```

There are some variants available as described below:

CMakePackage: h5bench

Description:

A benchmark suite **for** measuring HDF5 performance.

Homepage: <https://github.com/hpc-io/h5bench>

Preferred version:

1.2 [git] <https://github.com/hpc-io/h5bench.git> at commit [866af6777573d20740d02acc47a9080de093e4ad](#)

Safe versions:

develop [git] <https://github.com/hpc-io/h5bench.git> on branch develop  
 1.2 [git] <https://github.com/hpc-io/h5bench.git> at commit [866af6777573d20740d02acc47a9080de093e4ad](#)  
 1.1 [git] <https://github.com/hpc-io/h5bench.git> at commit [1276530a128025b83a4d9e3814a98f92876bb5c4](#)  
 1.0 [git] <https://github.com/hpc-io/h5bench.git> at commit [9d3438c1bc66c5976279ef203bd11a8d48ade724](#)  
 latest [git] <https://github.com/hpc-io/h5bench.git> on branch master

Deprecated versions:

None

Variants:

Name [Default]	When	Allowed values	Description
=====	=====	=====	
→ =====			

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all [off]	@1.2:	on, off	Enables all_
↪h5bench benchmarks			
amrex [off]	@1.2:	on, off	Enables AMReX_
↪benchmark			
build_type [RelWithDebInfo]	--	Debug, Release, RelWithDebInfo, MinSizeRel	CMake build type
e3sm [off]	@1.2:	on, off	Enables E3SM_
↪benchmark			
exerciser [off]	@1.2:	on, off	Enables exerciser_
↪benchmark			
ipo [off]	--	on, off	CMake_
↪interprocedural optimization			
metadata [off]	@1.2:	on, off	Enables metadata_
↪benchmark			
openpmd [off]	@1.2:	on, off	Enables OpenPMD_
↪benchmark			
Build Dependencies:			
cmake hdf5 mpi parallel-netcdf			
Link Dependencies:			
hdf5 mpi parallel-netcdf			
Run Dependencies:			
None			

**Warning:** Current h5bench versions in Spack do not have support for the HDF5 VOL async/cache connectors yet.

## RUNNING H5BENCH

### 2.1 h5bench (recommended)

We provide a single script you can use to run the benchmarks available in the h5bench Benchmarking Suite. You can combine multiple benchmarks into a workflow with distinct configurations. If you prefer, you can also manually run each benchmark in h5bench. For more details, refer to the Manual Execution section.

```
usage: h5bench [-h] [-a] [-d] [-v] [-p PREFIX] [-f FILTER] [-V] setup
```

H5bench: a Parallel I/O Benchmark Suite **for** HDF5:

positional arguments:

    setup                            JSON file **with** the benchmarks to run

options:

-h, --help	Show this help message <b>and</b> exit
-a, --abort-on-failure	Stop h5bench <b>if</b> a benchmark failed
-d, --debug	Enable debug mode
-v, --validate-mode	Validated <b>if</b> the requested mode ( <b>async</b> /sync) was run
-p PREFIX, --prefix PREFIX	Prefix where <b>all</b> h5bench binaries were installed
-f FILTER, --filter FILTER	Execute only filtered benchmarks
-V, --version	Show program's <b>version number and exit</b>

You need to provide a JSON file with the configurations you want to run. If you're using h5bench, you should *not* call `mpirun`, `srun`, or any other parallel launcher on your own. Refer to the manual execution section if you want to follow that approach instead. The main script will handle setting and unsetting environment variables, launching the benchmarks with the provided configuration and HDF5 VOL connectors.

```
h5bench configuration.json
```

If you run it with the `--debug` option, h5bench will also print log messages `stdout`. The default behavior is to store it in a file.

**Warning:** Make sure you do not call `srun`, `mpirun`, etc directly but instead define that in the JSON configuration file. You should **always** call h5bench directly.

You can define a single `.json` file with a complete setup and a combination of kernels you want to run. You can filter which of those benchmarks h5bench should run by passing the `-filter` option when running. For instance, the following command will only run the `read`, and `openpmd` kernels defined in the `.json`. The remaining ones would be ignored.

```
h5bench --filter read,openpmd configuration.json
```

## 2.1.1 Configuration

The JSON configuration file has five main properties: `mpi`, `vol`, `file-system`, `directory`, `benchmarks`. All should be defined, even if empty.

### MPI

You can set the MPI launcher you want to use, e.g. `mpirun`, `mpiexec`, and `srun`, and provide the number of processes you want to use. For other methods or a fine grain control on the job configuration, you can define the configuration properties that h5bench will use to launch the experiments using the `command` property you provided. If the configuration option is defined, h5bench will ignore the `ranks` property.

```
"mpi": {  
  "command": "mpirun",  
  "ranks": "4",  
  "configuration": "-np 8 --oversubscribe"  
}
```

### VOL

You can use HDF5 VOL connectors (`async`, `cache`, etc) for `h5bench_write` and `h5bench_read`. Because some benchmarks inside h5bench do not have support for VOL connectors yet, you need to provide the necessary information in the configuration file to handle the VOL setup during runtime.

```
"vol": {  
  "library": "/vol-async/src:/hdf5-async-vol-register-install/lib:/argobots/install/  
↳lib:/hdf5-install/install:",  
  "path": "/vol-async/src",  
  "connector": "async under_vol=0;under_info={}"  
}
```

You should provide the absolute path for all the libraries required by the VOL connector using the `library` property, the path of the VOL connector, and the configuration in `connector`. The provided example depicts how to configure the HDF5 VOL async connector.

### Directory

h5bench will create a directory for the given execution workflow, where it will store all the generated files and logs. Additional options such as data striping for Lustre, if configured, will be applied to this directory.

```
"directory": "hdf5-output-directory"
```

## File System

You can use this property to configure some file system options. For now, you can use it for Lustre to define the striping count and size that should be applied to the directory that will store all the generated data from h5bench.

```
"file-system": {
  "lustre": {
    "stripe-size": "1M",
    "stripe-count": "4"
  }
}
```

## Benchmarks

You can specify which benchmarks h5bench should run using this property, their order, and configuration. You can choose between: write, write-unlimited, overwrite, append, read, metadata, exerciser, openpmd, amrex, e3sm.

For each pattern of h5bench, you should provide the file and the configuration:

```
{
  "benchmark": "write",
  "file": "test.h5",
  "configuration": {
    "MEM_PATTERN": "CONTIG",
    "FILE_PATTERN": "CONTIG",
    "NUM_PARTICLES": "16 M",
    "Timesteps": "5",
    "DELAYED_CLOSE_Timesteps": "2",
    "COLLECTIVE_DATA": "NO",
    "COLLECTIVE_METADATA": "NO",
    "EMULATED_COMPUTE_TIME_PER_Timestep": "1 s",
    "NUM_DIMS": "1",
    "DIM_1": "16777216",
    "DIM_2": "1",
    "DIM_3": "1",
    "MODE": "SYNC",
    "CSV_FILE": "output.csv"
  }
}
```

If you provide the same file name used for a previous write execution, it will read from that file. This way, you can configure a workflow with multiple interleaving files, e.g., write file-01, write file-02, read file-02, read file-01.

```
{
  "benchmark": "read": {
    "file": "test.h5",
    "configuration": {
      "MEM_PATTERN": "CONTIG",
      "FILE_PATTERN": "CONTIG",
      "NUM_PARTICLES": "16 M",
      "Timesteps": "5",
      "DELAYED_CLOSE_Timesteps": "2",
```

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```

    "COLLECTIVE_DATA": "NO",
    "COLLECTIVE_METADATA": "NO",
    "EMULATED_COMPUTE_TIME_PER_TIMESTEP": "1 s",
    "NUM_DIMS": "1",
    "DIM_1": "16777216",
    "DIM_2": "1",
    "DIM_3": "1",
    "MODE": "SYNC",
    "CSV_FILE": "output.csv"
  }
}

```

You can also provide the `align` settings for GPFS filesystem in the `benchmark` property configuration. Note, not in the `filesystem` property. This parameter is enabled only for `h5bench-write` and `h5bench-write-unlimited`.

```

{
  "benchmark": "write",
  "file": "test.h5",
  "configuration": {
    "MEM_PATTERN": "CONTIG",
    "FILE_PATTERN": "CONTIG",
    "TIMESTEPS": "5",
    "DELAYED_CLOSE_TIMESTEPS": "2",
    "COLLECTIVE_DATA": "YES",
    "COLLECTIVE_METADATA": "YES",
    "EMULATED_COMPUTE_TIME_PER_TIMESTEP": "1 s",
    "NUM_DIMS": "1",
    "DIM_1": "4194304",
    "DIM_2": "1",
    "DIM_3": "1",
    "MODE": "ASYNC",
    "CSV_FILE": "output.csv",
    "ALIGN": "YES",
    "ALIGN_THRESHOLD": "16777216",
    "ALIGN_LEN": "16777216"
  }
}

```

For the metadata stress benchmark, `file` and `configuration` properties must be defined:

```

{
  "benchmark": "metadata",
  "file": "hdf5_iotest.h5",
  "configuration": {
    "version": "0",
    "steps": "20",
    "arrays": "500",
    "rows": "100",
    "columns": "200",
    "process-rows": "2",
    "process-columns": "2",
    "scaling": "weak",

```

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```

    "dataset-rank": "4",
    "slowest-dimension": "step",
    "layout": "contiguous",
    "mpi-io": "independent",
    "csv-file": "hdf5_iotest.csv"
  }
}

```

For the **exerciser** benchmark, you need to provide the required runtime options in the JSON file inside the **configuration** property.

```

{
  "benchmark": "exerciser",
  "configuration": {
    "numdims": "2",
    "minels": "8 8",
    "nsizes": "3",
    "bufmult": "2 2",
    "dimranks": "8 4"
  }
}

```

You can find several samples of configuration file with all the options in the our [GitHub repository] (<https://github.com/hpc-io/h5bench/tree/master/samples>). You can also refer to this sample of a complete `configuration.json` file that defined the workflow of the execution of multiple benchmarks from h5bench Suite:

```

{
  "mpi": {
    "command": "mpirun",
    "ranks": "4",
    "configuration": "-np 8 --oversubscribe"
  },
  "vol": {
    "library": "/vol-async/src:/hdf5-async-vol-register-install/lib:/argobots/
↪install/lib:/hdf5-install/install:",
    "path": "/vol-async/src",
    "connector": "async under_vol=0;under_info={}"
  },
  "file-system": {
    "lustre": {
      "stripe-size": "1M",
      "stripe-count": "4"
    }
  },
  "directory": "full-teste",
  "benchmarks": [
    {
      "benchmark": "e3sm",
      "file": "coisa.h5",
      "configuration": {
        "k": "",
        "x": "blob",

```

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```

        "a": "hdf5",
        "r": "25",
        "o": "ON",
        "netcdf": "../..e3sm/datasets/f_case_866x72_16p.nc"
    }
},
{
    "benchmark": "write",
    "file": "test.h5",
    "configuration": {
        "MEM_PATTERN": "CONTIG",
        "FILE_PATTERN": "CONTIG",
        "NUM_PARTICLES": "16 M",
        "Timesteps": "5",
        "DELAYED_CLOSE_Timesteps": "2",
        "COLLECTIVE_DATA": "NO",
        "COLLECTIVE_METADATA": "NO",
        "EMULATED_COMPUTE_TIME_PER_Timestep": "1 s",
        "NUM_DIMS": "1",
        "DIM_1": "16777216",
        "DIM_2": "1",
        "DIM_3": "1",
        "MODE": "SYNC",
        "CSV_FILE": "output.csv"
    }
},
{
    "benchmark": "exerciser",
    "configuration": {
        "numdims": "2",
        "minels": "8 8",
        "nsizes": "3",
        "bufmult": "2 2",
        "dimranks": "8 4"
    }
},
{
    "benchmark": "exerciser",
    "configuration": {
        "numdims": "2",
        "minels": "8 8",
        "nsizes": "3",
        "bufmult": "2 2",
        "dimranks": "8 4"
    }
},
{
    "benchmark": "exerciser",
    "configuration": {
        "numdims": "2",
        "minels": "8 8",
        "nsizes": "3",

```

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```

        "bufmult": "2 2",
        "dimranks": "8 4"
    }
},
{
    "benchmark": "exerciser",
    "configuration": {
        "numdims": "2",
        "minels": "8 8",
        "nsizes": "3",
        "bufmult": "2 2",
        "dimranks": "8 4"
    }
},
{
    "benchmark": "exerciser",
    "configuration": {
        "numdims": "2",
        "minels": "8 8",
        "nsizes": "3",
        "bufmult": "2 2",
        "dimranks": "8 4"
    }
},
{
    "benchmark": "read",
    "file": "test.h5",
    "configuration": {
        "MEM_PATTERN": "CONTIG",
        "FILE_PATTERN": "CONTIG",
        "NUM_PARTICLES": "16 M",
        "Timesteps": "5",
        "DELAYED_CLOSE_TIMESTEPS": "2",
        "COLLECTIVE_DATA": "NO",
        "COLLECTIVE_METADATA": "NO",
        "EMULATED_COMPUTE_TIME_PER_TIMESTEP": "1 s",
        "NUM_DIMS": "1",
        "DIM_1": "16777216",
        "DIM_2": "1",
        "DIM_3": "1",
        "MODE": "SYNC",
        "CSV_FILE": "output.csv"
    }
},
{
    "benchmark": "write",
    "file": "test-two.h5",
    "configuration": {
        "MEM_PATTERN": "CONTIG",
        "FILE_PATTERN": "CONTIG",
        "NUM_PARTICLES": "2 M",
        "Timesteps": "20",

```

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```

        "DELAYED_CLOSE_TIMESTEPS": "2",
        "COLLECTIVE_DATA": "NO",
        "COLLECTIVE_METADATA": "NO",
        "EMULATED_COMPUTE_TIME_PER_TIMESTEP": "1 s",
        "NUM_DIMS": "1",
        "DIM_1": "16777216",
        "DIM_2": "1",
        "DIM_3": "1",
        "MODE": "SYNC",
        "CSV_FILE": "output.csv"
    },
    {
        "benchmark": "metadata",
        "file": "hdf5_iotest.h5",
        "configuration": {
            "version": "0",
            "steps": "20",
            "arrays": "500",
            "rows": "100",
            "columns": "200",
            "process-rows": "2",
            "process-columns": "2",
            "scaling": "weak",
            "dataset-rank": "4",
            "slowest-dimension": "step",
            "layout": "contiguous",
            "mpi-io": "independent",
            "csv-file": "hdf5_iotest.csv"
        }
    }
]
}

```

For a description of all the options available in each benchmark, please refer to their entries in the documentation.

When the `--debug` option is enabled, you can expect an output similar to:

```

2021-10-25 16:31:24,866 h5bench - INFO - Starting h5bench Suite
2021-10-25 16:31:24,889 h5bench - INFO - Lustre support detected
2021-10-25 16:31:24,889 h5bench - DEBUG - Lustre stripping configuration: lfs setstripe -
↳S 1M -c 4 full-teste
2021-10-25 16:31:24,903 h5bench - INFO - h5bench [write] - Starting
2021-10-25 16:31:24,903 h5bench - INFO - h5bench [write] - DIR: full-teste/504fc233/
2021-10-25 16:31:24,904 h5bench - INFO - Parallel setup: srun --cpu_bind=cores -n 4
2021-10-25 16:31:24,908 h5bench - INFO - srun --cpu_bind=cores -n 4 build/h5bench_write_
↳full-teste/504fc233/h5bench.cfg full-teste/test.h5
2021-10-25 16:31:41,670 h5bench - INFO - SUCCESS
2021-10-25 16:31:41,754 h5bench - INFO - Runtime: 16.8505464 seconds (elapsed time,
↳includes allocation wait time)
2021-10-25 16:31:41,755 h5bench - INFO - h5bench [write] - Complete
2021-10-25 16:31:41,755 h5bench - INFO - h5bench [exerciser] - Starting
2021-10-25 16:31:41,755 h5bench - INFO - h5bench [exerciser] - DIR: full-teste/247659d1/

```

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```

2021-10-25 16:31:41,755 h5bench - INFO - Parallel setup: srun --cpu_bind=cores -n 4
2021-10-25 16:31:41,756 h5bench - INFO - srun --cpu_bind=cores -n 4 build/h5bench_
↳exerciser --numdims 2 --minels 8 8 --nsizes 3 --bufmult 2 2 --dimranks 8 4
2021-10-25 16:31:49,174 h5bench - INFO - SUCCESS
2021-10-25 16:31:49,174 h5bench - INFO - Finishing h5bench Suite

```

## Cori (NERSC)

In Cori you need to load Python and its libraries for the main h5bench script to work. For manual execution of each benchmark that is not required.

```
module load python
```

In case you are running on Cori and the benchmark fails with an MPI message indicating no support for multiple threads:

```
Assertion `MPI_THREAD_MULTIPLE == mpi_thread_lvl_p rovided' failed.
```

Please, make sure you define the following:

```
export MPICH_MAX_THREAD_SAFETY="multiple"
```

## Sunspot (ALCF)

In Sunspot you need to export one additional environment variable related to ATS. ATS is the Address Translation Service support for using the IOMMU (Input-Output Memory Management Unit) for address translation. ATS is not supported on Intel processors at this time. The default is to NTA (NIC translation).

```
export FI_CXI_ATS=0
```

Otherwise you will encounter the following error:

```
libfabric:36807:1674015247::cxi:core:cxip_fc_notify_cb():4366<warn>
x1922c0s5b0n0: TXC (0x1081:5:0):: Fatal, unexpected event rc: 26
```

## 2.2 Manual Execution

If you prefer, you can execute each benchmark manually. In this scenario, you will be responsible for generating the input configuration file needed for each benchmark in the suite, ensuring it follows the pre-defined format unique for each one.

If you want to use HDF5 VOL connectors or tune the file system configuration, h5bench will *not* take care of that. Remember that not all benchmarks in the suite have support for VOL connectors yet.



## READ / WRITE

This set of benchmarks contains an I/O kernel developed based on a particle physics simulation's I/O pattern (VPIC-IO for writing data in a HDF5 file) and on a big data clustering algorithm (BDCATS-IO for reading the HDF5 file VPIC-IO wrote).

### 3.1 Configuration

You can configure the `h5bench_write` and `h5bench_read` benchmarks with the following options. Notice that if you use the `configuration.json` approach to define the runs for `h5bench`, we will automatically generate the final configuration file based on the options you provide in the JSON file. For standalone usage of this benchmark, you can check the input format at the end of this document and refer to its documentation.

Parameter	Description
MEM_PATTERN	Options: CONTIG, INTERLEAVED, and STRIDED
FILE_PATTERN	Options: CONTIG and STRIDED
TIMESTEPS	The number of iterations
EMULATED_COMPUTE_TIME_PER_TIMESTEP	Sleeps after each iteration to emulate computation
NUM_DIMS	The number of dimensions, valid values are 1, 2 and 3
DIM_1	The dimensionality of the source data
DIM_2	The dimensionality of the source data
DIM_3	The dimensionality of the source data

For `MEM_PATTERN`, `CONTIG` represents arrays of basic data types (i.e., int, float, double, etc.); `INTERLEAVED` represents an array of structure (AOS) where each array element is a C struct; and `STRIDED` represents a few elements in an array of basic data types that are separated by a constant stride. `STRIDED` is supported only for 1D arrays.

For `FILE_PATTERN`, `CONTIG` represents a HDF5 dataset of basic data types (i.e., int, float, double, etc.); `INTERLEAVED` represents a dataset of a compound datatype;

For `EMULATED_COMPUTE_TIME_PER_TIMESTEP`, you *must* provide the time unit (e.g. 10 s, 100 ms, or 5000us) to ensure correct behavior.

For `DIM_2` and `DIM_3` if **unused**, you should set both as 1. Notice that the total number of particles will be given by `DIM_1 * DIM_2 * DIM_3`. For example, `DIM_1=1024`, `DIM_2=256`, `DIM_3=1` is a valid setting for a 2D array and it will generate 262144 particles.

A set of sample configuration files can be found in the `samples/` directory in GitHub.

### 3.1.1 READ Settings (h5bench\_read)

Parameter	Description
READ_OPTION	Options: FULL, PARTIAL, and STRIDED

For the PARTIAL option, the benchmark will read only the first TO\_READ\_NUM\_PARTICLES particles.

### 3.1.2 Asynchronous Settings

Parameter	Description
MODE	Options: SYNC or ASYNC
IO_MEM_LIMIT	Memory threshold to determine when to execute I/O
DELAYED_CLOSE_TIMESTEPS	Groups and datasets will be closed later.

The IO\_MEM\_LIMIT parameter is optional. Its default value is 0 and it requires ASYNC, i.e., it only works in asynchronous mode. This is the memory threshold used to determine when to actually execute the I/O operations. The actual I/O operations (data read/write) will not be executed until the timesteps associated memory reaches the threshold, or the application run to the end.

For the ASYNC mode to work you **must** define the necessary HDF5 ASYNC-VOL connector. For more information about it, refer to its [documentation](#).

### 3.1.3 Compression Settings

Parameter	Description
COMPRESS	YES or NO (optional) enables parallel compression
CHUNK_DIM_1	Chunk dimension
CHUNK_DIM_2	Chunk dimension
CHUNK_DIM_3	Chunk dimension

Compression is only applicable for h5bench\_write. It has no effect for h5bench\_read. When enabled the chunk dimensions parameters (CHUNK\_DIM\_1, CHUNK\_DIM\_2, CHUNK\_DIM\_3) are required. The chunk dimension settings should be compatible with the data dimensions, i.e., they must have the same rank of dimensions, and chunk dimension size cannot be greater than data dimension size. Extra chunk dimensions have no effect and should be set to 1.

**Warning:** There is a known bug on HDF5 parallel compression that could cause the system run out of memory when the chunk amount is large (large number of particle and very small chunk sizes). On Cori Hasswell nodes, the setting of 16M particles per rank, 8 nodes (total 256 ranks), 64 \* 64 chunk size will crash the system by running out of memory, on single nodes the minimal chunk size is 4 \* 4.

### 3.1.4 Collective Operation Settings

Parameter	Description
COLLECTIVE_DATA	Enables collective operation (default is NO)
COLLECTIVE_METADATA	Enables collective HDF5 metadata (default is NO)

Both COLLECTIVE\_DATA and COLLECTIVE\_METADATA parameters are optional.

### 3.1.5 Subfiling Settings

Parameter	Description
SUBFILING	Enables HDF5 subfiling (default is NO)

**Attention:** In order to enable this option your HDF5 must have been compiled with support for the HDF5 Subfiling Virtual File Driver (VFD) which was introduced in the HDF5 1.14.0. For CMake you can use the `-DHDF5_ENABLE_PARALLEL=ON -DHDF5_ENABLE_SUBFILING_VFD=ON` and for autotools `--enable-parallel --enable-subfiling-vfd=yes`. Without this support, this parameter has no effect.

### 3.1.6 CSV Settings

Performance results will be written to this file and standard output once a file name is provided.

Parameter	Description
CSV_FILE	CSV file name to store benchmark results

## 3.2 Supported Patterns

**Attention:** Not every pattern combination is covered by the benchmark. Supported benchmark parameter settings are listed below.

### 3.2.1 Supported Write Patterns (h5bench\_write)

The I/O patterns include array of structures (AOS) and structure of arrays (SOA) in memory as well as in file. The array dimensions are 1D, 2D, and 3D for the write benchmark. This defines the write access pattern, including CONTIG (contiguous), INTERLEAVED and STRIDED for the source (the data layout in the memory) and the destination (the data layout in the resulting file). For example, MEM\_PATTERN=CONTIG and FILE\_PATTERN=INTERLEAVED is a write pattern where the in-memory data layout is contiguous (see the implementation of `prepare_data_contig_2D()` for details) and file data layout is interleaved by due to its compound data structure (see the implementation of `data_write_contig_to_interleaved()` for details).

- 4 patterns for both 1D and 2D array write (NUM\_DIMS=1 or NUM\_DIMS=2)

```
'MEM_PATTERN': 'CONTIG'
'FILE_PATTERN': 'CONTIG'
```

```
'MEM_PATTERN': 'CONTIG'
'FILE_PATTERN': 'INTERLEAVED'
```

```
'MEM_PATTERN': 'INTERLEAVED'
'FILE_PATTERN': 'CONTIG'
```

```
'MEM_PATTERN': 'INTERLEAVED'
'FILE_PATTERN': 'INTERLEAVED'
```

- 1 pattern for 3D array (NUM\_DIMS=3)

```
'MEM_PATTERN': 'CONTIG'
'FILE_PATTERN': 'CONTIG'
```

- 1 strided pattern for 1D array (NUM\_DIMS=1)

```
'MEM_PATTERN': 'CONTIG'
'FILE_PATTERN': 'STRIDED'
```

### 3.2.2 Supported Read Patterns (h5bench\_read)

- 1 pattern for 1D, 2D and 3D read (NUM\_DIMS=1 or NUM\_DIMS=2)

Contiguously read through the whole data file:

```
'MEM_PATTERN': 'CONTIG'
'FILE_PATTERN': 'CONTIG'
'READ_OPTION': 'FULL'
```

- 2 patterns for 1D read

Contiguously read the first TO\_READ\_NUM\_PARTICLES elements:

```
'MEM_PATTERN': 'CONTIG'
'FILE_PATTERN': 'CONTIG'
'READ_OPTION': 'PARTIAL'
```

```
'MEM_PATTERN': 'CONTIG'
'FILE_PATTERN': 'STRIDED'
'READ_OPTION': 'STRIDED'
```

## 3.3 Understanding the Output

The metadata and raw data operations are timed separately, and the overserved time and I/O rate are based on the total time.

Sample output of h5bench\_write:

```
===== Performance results =====
Total emulated compute time 4000 ms
Total write size = 2560 MB
```

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```

Data preparation time = 739 ms
Raw write time = 1.012 sec
Metadata time = 284.990 ms
H5Fcreate() takes 4.009 ms
H5Fflush() takes 14.575 ms
H5Fclose() takes 4.290 ms
Observed completion time = 6.138 sec
Raw write rate = 2528.860 MB/sec
Observed write rate = 1197.592 MB/sec

```

Sample output of h5bench\_read:

```

===== Performance results =====
Total emulated compute time = 4 sec
Total read size = 2560 MB
Metadata time = 17.523 ms
Raw read time = 1.201 sec
Observed read completion time = 5.088 sec
Raw read rate = 2132.200 MB/sec
Observed read rate = 2353.605225 MB/sec

```

## 3.4 Known Issues

**Warning:** In Cori/NERSC or similar platforms that use Cray-MPICH library, if you encounter a failed assertion regarding support for MPI\_THREAD\_MULTIPLE you should define the following environment variable:

```
export MPICH_MAX_THREAD_SAFETY="multiple"
```

**Warning:** If you're trying to run the benchmark with the HDF5 VOL ASYNC connector in MacOS and are getting segmentation fault (from ABT\_thread\_create), please try to set the following environment variable:

```
export ABT_THREAD_STACKSIZE=1000000
```



## METADATA STRESS

The Metadata Stress benchmark (`h5bench_hdf5_iotest`) is a simple I/O performance tester for HDF5. Its purpose is to assess the performance variability of a set of logically equivalent HDF5 representations of a common pattern. The test repeatedly writes (and reads) in parallel a set of 2D array variables in a tiled fashion, over a set of time steps. For more information referer to HDF Group [GitHub repository](#). We modified this benchmark slightly so to be able to specify the config file location, everything else remains untouched.

### 4.1 Configuration

You can configure the Metadata Stress test with the following options. Notice that if you use the `configuration.json` approach to define the runs for `h5bench`, we will automatically generate the final configuration file based on the options you provide in the JSON file. For standalone usage of this benchmark, you can check the input format at the end of this document and refer to its documentation.

Parameter	Description
<code>steps</code>	Number of steps
<code>arrays</code>	Number of arrays
<code>rows</code>	Total number of array rows for strong scaling. Number of array rows per block for weak scaling.
<code>columns</code>	Total number of array columns for strong scaling. Number of array columns per block for weak scaling.
<code>process-rows</code>	Number of MPI-process rows: <code>rows % proc-rows == 0</code> for strong scaling
<code>process-columns</code>	Number of MPI-process columns: <code>columns % proc-columns == 0</code> for strong scaling
<code>scaling</code>	Scaling ([weak, strong])
<code>dataset-rank</code>	Rank of the dataset(s) in the file ([2, 3, 4])
<code>slowest-dimension</code>	Slowest changing dimension ([step, array])
<code>layout</code>	HDF5 dataset layout ([contiguous, chunked])
<code>mpi-io</code>	MPI I/O mode ([independent, collective])
<code>hdf5-file</code>	HDF5 output file name
<code>csv-file</code>	CSV results file name

### 4.1.1 JSON Configuration (recommended)

To run an instance of Metadata Stress Test benchmark you need to include the following in the `benchmarks` property of your `configuration.json` file:

```
{
  "benchmark": "metadata",
  "file": "hdf5_iotest.h5",
  "configuration": {
    "version": "0",
    "steps": "20",
    "arrays": "500",
    "rows": "100",
    "columns": "200",
    "process-rows": "2",
    "process-columns": "2",
    "scaling": "weak",
    "dataset-rank": "4",
    "slowest-dimension": "step",
    "layout": "contiguous",
    "mpi-io": "independent",
    "csv-file": "hdf5_iotest.csv"
  }
}
```

### 4.1.2 Standalone Configuration

For standalone usage of this benchmark, this is the observed input configuration you should provide to the `h5bench_hdf5_iotest` executable.

```
[DEFAULT]
version = 0
steps = 20
arrays = 500
rows = 100
columns = 200
process-rows = 1
process-columns = 1
scaling = weak
dataset-rank = 4
slowest-dimension = step
layout = contiguous
mpi-io = independent
hdf5-file = hdf5_iotest.h5
csv-file = hdf5_iotest.csv
```

## AMREX

AMReX is a software framework for massively parallel, block-structured adaptive mesh refinement (AMR) applications.

You can find more information in AMReX-Codes [GitHub repository](#).

### 5.1 Configuration

You can configure the AMReX HDF5 benchmark with the following options. Notice that if you use the `configuration.json` approach to define the runs for `h5bench`, we will automatically generate the final configuration file based on the options you provide in the JSON file. For standalone usage of this benchmark, you can check the input format at the end of this document and refer to its documentation.

Parameter	Description
<code>ncells</code>	Domain size
<code>max_grid_size</code>	The maximum allowable size of each subdomain (used for parallel decomposal)
<code>nlevs</code>	Number of levels
<code>ncomp</code>	Number of components in the multifabs
<code>nppc</code>	Number of particles per cell
<code>nplotfile</code>	Number of plot files to write
<code>nparticlefile</code>	Number of particle files to write
<code>sleeptime</code>	Time to sleep before each write
<code>restart_check</code>	Whether to check the correctness of checkpoint/restart
<code>grids_from_file</code>	Enable AMReX to read grids from file
<code>ref_ratio_file</code>	Refinement ratios for different AMReX refinement levels
<code>hdf5compression</code>	Define the HDF5 compression algorithm to use

#### 5.1.1 JSON Configuration (recommended)

To run an instance of AMReX HDF5 benchmark you need to include the following in the `benchmarks` property of your `configuration.json` file:

```
{  
  "benchmark": "amrex",  
  "file": "amrex.h5",  
  "configuration": {  
    "ncells": "64",  
    "max_grid_size": "8",  
  }  
}
```

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```

        "nlevs": "1",
        "ncomp": "6",
        "nppc": "2",
        "nplotfile": "2",
        "nparticlefile": "2",
        "sleeptime": "2",
        "restart_check": "1",
        "hdf5compression": "ZFP_ACCURACY#0.001",
        "mode": "SYNC"
    }
}

```

To read grids from file you need to set: `grids_from_file`, `nlevels`, and `ref_ratio_file`.

```

{
  "benchmark": "amrex",
  "file": "amrex.h5",
  "configuration": {
    "ncells": "64",
    "max_grid_size": "8",
    "nlevs": "1",
    "ncomp": "6",
    "nppc": "2",
    "nplotfile": "2",
    "nparticlefile": "2",
    "sleeptime": "2",
    "restart_check": "1",
    "hdf5compression": "ZFP_ACCURACY#0.001",
    "mode": "SYNC",
    "nlevs": "3",
    "grids_from_file": "1",
    "ref_ratio_file": "4 2"
  }
}

```

### 5.1.2 HDF5 ASYNC VOL Connector

AMReX supports the [HDF5 ASYNC VOL connector](#). To enable it, you should specify in the `vol` property of your `configuration.json` file: the required library paths, the VOL ASYNC source path, and the connector setup.

```

"vol": {
  "library": "/vol-async/src:/hdf5-async-vol-register-install/lib:/argobots/install/
↳ lib:/hdf5-install/install:",
  "path": "/vol-async/src",
  "connector": "async under_vol=0;under_info={}"
}

```

### 5.1.3 Standalone Configuration

For standalone usage of this benchmark, this is the observed input configuration you should provide to the `h5bench_amrex` executable.

```
ncells = 64
max_grid_size = 8
nlevs = 1
ncomp = 6
nppc = 2
nplotfile = 2
nparticlefile = 2
sleeptime = 2
restart_check = 1

# Uncomment to read grids from file
# nlevs = 3
# grids_from_file = 1
# ref_ratio_file = 4 2

# Uncomment to enable compression
# hdf5compression=ZFP_ACCURACY#0.001

directory = .
```





## OPENPMD

OpenPMD is an open meta-data schema that provides meaning and self-description for data sets in science and engineering.

The openPMD-api library provides a reference API for openPMD data handling. In the h5bench Benchmarking Suite we provide support for the write and read parallel benchmarks with HDF5 backend. You can find more information in [OpenPMD documentation](#).

### 6.1 Configuration

You can configure the openPMD write HDF5 benchmark with the following options. Notice that if you use the `configuration.json` approach to define the runs for h5bench, we will automatically generate the final configuration file based on the options you provide in the JSON file. For standalone usage of this benchmark, you can check the input format at the end of this document and refer to its documentation.

Parameter	Description
<code>operation</code>	Operation: write or read
<code>fileLocation</code>	Directory where the file will be written to or read from

When running with the `write` operation, you have to define the following options:

`dim` Number of dimensions (1, 2, or 3) `balanced` Should it use a balanced load? (`true` or `false`) `ratio` Particle to mesh ratio `steps` Number of iteration steps `minBlock` Meshes are viewed as grid of mini blocks `grid` Grid based on the mini block

When running with the `read` operation, you have to define the pattern:

`pattern` Read access pattern

The `minBlock` and `grid` parameters must include the values for each of the `dim` dimensions. For example, if `"dim": "3"` (for a 3D mesh) `minBlock` should contain three values, one for each dimension `"16 32 32"` and `grid` (which is based on the mini block) should also contain three values, one for each dimension `"32 32 16"`.

For the `pattern` attribute for read you can choose:

- `m`: metadata only
- `sx`: slice of the 'rho' mesh in the x-axis (eg. `x=0`)
- `sy`: slice of the 'rho' mesh in the y-axis (eg. `y=0`)
- `sz`: slice of the 'rho' mesh in the z-axis (eg. `z=0`)
- `fx`: slice of the 3D magnetic field in the x-axis (eg. `x=0`)
- `fy`: slice of the 3D magnetic field in the y-axis (eg. `y=0`)

- fz: slice of the 3D magnetic field in the z-axis (eg. z=0)

### 6.1.1 JSON Configuration (recommended)

To run an instance of openPMD HDF5 benchmark you need to include the following in the `benchmarks` property of your `configuration.json` file:

```
{
  "benchmark": "openpmd",
  "configuration": {
    "operation": "write",
    "dim": "3",
    "balanced": "true",
    "ratio": "1",
    "steps": "1",
    "minBlock": "8 16 16",
    "grid": "16 16 8"
  }
},
{
  "benchmark": "openpmd",
  "configuration": {
    "operation": "read",
    "dim": "3",
    "balanced": "true",
    "ratio": "1",
    "steps": "1",
    "minBlock": "8 16 16",
    "grid": "16 16 8"
  }
}
```

### 6.1.2 Standalone Configuration

For standalone usage of this benchmark, this is the observed input configuration you should provide to the `h5bench_openpmd_write`.

```
dim=3
balanced=true
ratio=1
steps=10
minBlock=16 32 32
grid=32 32 16
```

For the `h5bench_openpmd_read`, you need to provide two arguments: the file prefix and the pattern.

## EXERCISER

**Attention:** For more-detailed instructions of how to build and run the exerciser code on specific machines (at ALCF), see the `Exerciser/BGQ/VESTA_XL/README.md` and `Exerciser/BGQ/THETA/README.md` directories of this repository. Those README files also include instructions for building the CCIO and develop versions of HDF5 for use with this benchmark.

The **HDF5 Exerciser Benchmark** creates an HDF5 use case with some ideas/code borrowed from other benchmarks (namely IOR, VPICIO and FLASHIO). Currently, the algorithm does the following in parallel over all MPI ranks:

- For each rank, a local data buffer (with dimensions given by `numdims` is initialized with `minNEls` double-precision elements in each dimension
- If the `derivedtype` flag is used, a second local dataset is also specified with a derived data type assigned to each element
- For a given number of iterations (hardcoded as `NUM_ITERATIONS`):
  - Open a file, create a top group, set the MPI-IO transfer property, and (optionally) add a simple attribute string to the top group
  - Create memory and file dataspace with hyperslab selections for simple rank-ordered offsets into the file. The `rshift` option can be used to specify the number of rank positions to shift the write position in the file (the read will be shifted twice this amount to avoid client-side caching effects)
  - Write the data and close the file
  - Open the file, read in the data, and check correctness (if dataset is small enough)
  - Close the dataset (but not the file)
  - If the second (derived-type) data set is specified: (1) create a derived type, (2) open a new data set with the same number of elements and dimension, (3) write the data and (4) close everything
- Each dimension of `curNEls` is then multiplied by each dimension of `bufMult`, and the previous steps (the loop over `NUM_ITERATIONS`) are repeated. This outer loop over local buffer sizes is repeated a total of `nsizes` times

## 7.1 Configuration

You can configure the `h5bench_write` and `h5bench_read` benchmarks with the following options. Notice that if you use the `configuration.json` approach to define the runs for `h5bench`, we will automatically generate the final configuration file based on the options you provide in the JSON file. For standalone usage of this benchmark, you can check the input format at the end of this document and refer to its documentation.

### 7.1.1 Required

Parameter	Description
<code>numdims &lt;x&gt;</code>	Dimension of the datasets to write to the HDF5 file
<code>minels &lt;x&gt; ... &lt;x&gt;</code>	Min number of double elements to write in each dim of the dataset (one value for each dimension)

### 7.1.2 Optional

Parameter	Description
<code>nsizes &lt;x&gt;</code>	How many buffer sizes to use (Code will start with <code>minbuf</code> and loop through <code>nsizes</code> iterations, with the buffer size multiplied by <code>bufmult</code> in each dim, for each iteration)
<code>bufmult &lt;x&gt; ... &lt;x&gt;</code>	Constant, for each dimension, used to multiply the buffer [default: 2 2 ... ]
<code>metacoll</code>	Whether to set meta data collective usage [default: False]
<code>derivedtype</code>	Whether to create a second data set containing a derived type [default: False]
<code>addattr</code>	Whether to add attributes to group 1 [default: False]
<code>indepio</code>	Whether to use independant I/O (not MPI-IO) [default: False]
<code>keepfile</code>	Whether to keep the file around after the program ends for further analysis, otherwise deletes it [default: False]
<code>usechunked</code>	Whether to chunk the data when reading/writing [default: False]
<code>maxcheck &lt;x&gt;</code>	Maximum buffer size (in bytes) to validate. Note that all buffers will be validated if this option is not set by this command-line argument [default: Inf]
<code>memblock &lt;x&gt;</code>	Define the block size to use in the local memory buffer (local buffer is always 1D for now, Note: This currently applies to the 'double' dataset only) [default: local buffer size]
<code>memstride &lt;x&gt;</code>	Define the stride of the local memory buffer (local buffer is always 1D for now, Note: This currently applies to the 'double' dataset only) [default: local buffer size]
<code>fileblocks &lt;x&gt; ...&lt;x&gt;</code>	Block sizes to use in the file for each dataset dimension (Note: This currently applies to the 'double' dataset only) [default: 1 ... 1]
<code>filestrides &lt;x&gt; ...&lt;x&gt;</code>	Stride dist. to use in the file for each dataset dimension (Note: This currently applies to the 'double' dataset only) [default: 1 ... 1]

The exerciser also allows the MPI decomposition to be explicitly defined:

Parameter	Description
<code>dimranks &lt;x&gt; ...&lt;x&gt;</code>	MPI-rank division in each dimension. Note that, if not set, decomposition will be in 1st dimension only.

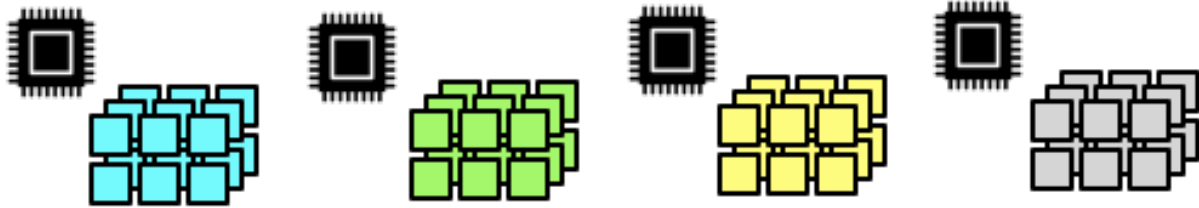
## 7.2 Exerciser Basics

In the simplest case, the Exerciser code will simply write and then read an n-dimensional double-precision dataset in parallel (with all the necessary HDF5 steps in between). At a minimum, the user must specify the number of dimensions to use for this dataset (using the `numdims` flag), and the size of each dimension (using the `minels` flag). By default, the maximum number of dimensions allowed by the code is set by `MAX_DIM` (currently 4, but can be modified easily). Note that the user is specifying the number of elements to use in each dimension with `minels`. Therefore, the local buffer size is the product of the dimension sizes and `sizeof(double)` (and the global dataset in the file is a product of the total MPI ranks and the local buffer size). As illustrated in Fig. 1, the mapping of ranks to hyper-slabs in the global dataset can be specified with the `dimranks` flag (here, Example 1 is the default decomposition, while Example 2 corresponds to: `"dimranks": "2 2"`). This flag simply allows the user to list the number of spatial decompositions in each dimension of the global dataset, and requires that the product of the input to be equal to the total number of MPI ranks.

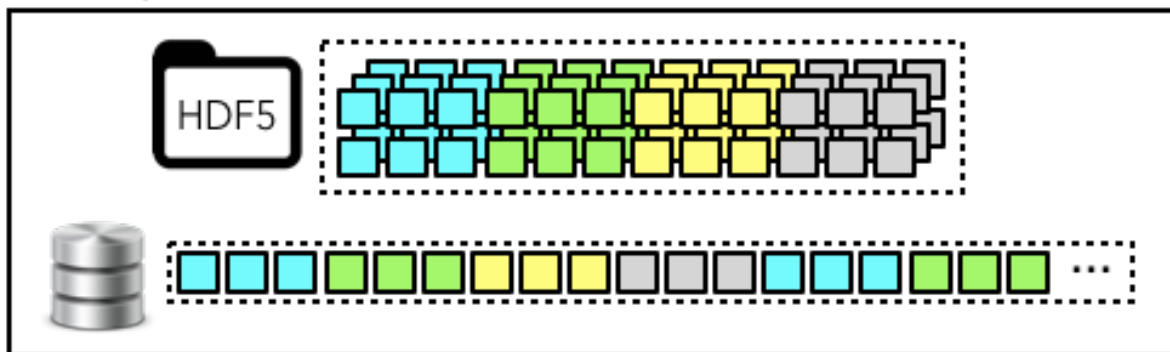
---

**Note:** Authors:

- Richard J. Zamora ([rzamora@anl.gov](mailto:rzamora@anl.gov))
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  - Venkatram Vishwanath ([venkat@anl.gov](mailto:venkat@anl.gov))
-



Example 1:



Example 2:

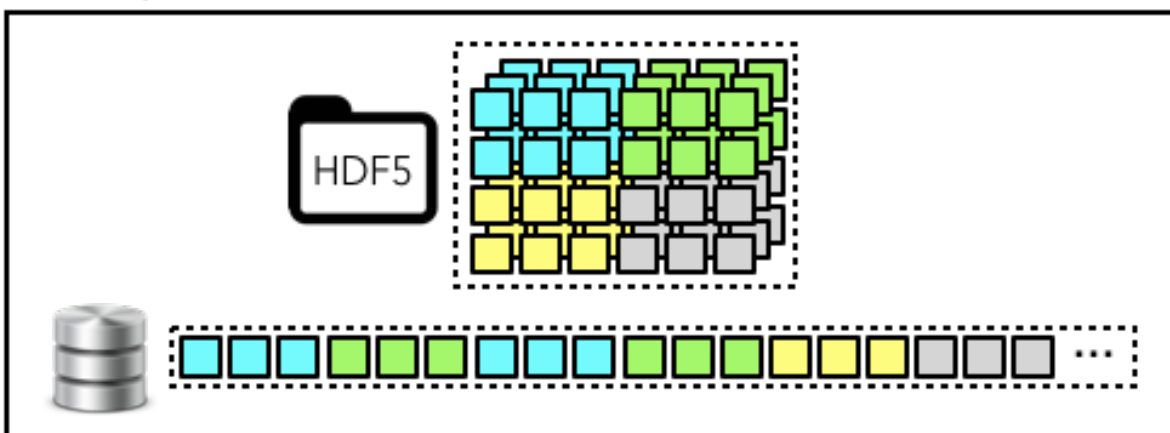


Fig. 1: Fig. 1 - Illustration of different local-to-global dataset mapping options.

**E3SM**

E3SM-IO is the parallel I/O kernel from the E3SM climate simulation model. It makes use of PIO library which is built on top of PnetCDF.

This benchmark currently has two cases from E3SM, namely F and G cases. The F case uses three unique data decomposition patterns shared by 388 2D and 3D variables (2 sharing Decomposition 1, 323 sharing Decomposition 2, and 63 sharing Decomposition 3). The G case uses 6 data decompositions shared by 52 variables (6 sharing Decomposition 1, 2 sharing Decomposition 2, 25 sharing Decomposition 3, 2 sharing Decomposition 4, 2 sharing Decomposition 5, and 4 sharing Decomposition 6).

You can find more information in Parallel-NetCDF [GitHub repository](#).

## 8.1 Configuration

You can configure the E3SM-IO benchmark with the following options. Notice that if you use the `configuration.json` approach to define the runs for `h5bench`, we will automatically generate the final configuration file based on the options you provide in the JSON file. For standalone usage of this benchmark, you can refer to E3SM-IO repository.

Parameter	Description
<code>k</code>	Keep the output files when program exits
<code>x</code>	I/O strategy to write ( <code>canonical</code> , <code>log</code> , and <code>blob</code> )
<code>a</code>	I/O library name to perform write operation ( <code>hdf5</code> , <code>hdf5_log</code> , <code>hdf5_md</code> )
<code>r</code>	Number of records/time steps for F case h1 file
<code>o</code>	Enable write performance evaluation
<code>netcdf</code>	Define the HDF5 compression algorithm to use

**Warning:** `h5bench` temporarily only supports `-x blob` and `-a hdf5`. If you set other options, they will be overwritten to the supported version.

### 8.1.1 JSON Configuration (recommended)

To run an instance of AMReX HDF5 benchmark you need to include the following in the `benchmarks` property of your `configuration.json` file:

```
{
  "benchmark": "e3sm",
  "file": "coisa.h5",
  "configuration": {
    "k": "",
    "x": "blob",
    "a": "hdf5",
    "r": "25",
    "o": "ON",
    "netcdf": "../../../e3sm/datasets/f_case_866x72_16p.nc"
  }
}
```



## MACSIO

MACSio (Multi-purpose, Application-Centric, Scalable I/O Proxy Application) is being developed to fill a long existing void in co-design proxy applications that allow for I/O performance testing and evaluation of tradeoffs in data models, I/O library interfaces and parallel I/O paradigms for multi-physics, HPC applications.

MACSio in h5bench only supports the HDF5 interface. You need to have the [json-cwx](#) dependency library installed prior to compiling it in h5bench.

You can find more information in MACSio [GitHub repository](#).

### 9.1 Configuration

You can configure the MACSio benchmark with the following options. Notice that if you use the `configuration.json` approach to define the runs for h5bench, we will automatically generate the final configuration file based on the options you provide in the JSON file. For standalone usage of this benchmark, you can refer to MACSio repository.

Parameter	Description
<code>parallel_file_mode</code>	Defines the parallel file mode
<code>part_size</code>	Defines the request sized (default per-proc request size is 80,0000 bytes (10K doubles))
<code>avg_num_parts</code>	Number of parts per process (default is 2)
<code>num_dumps</code>	Number of dumps (default is 10)

For the `parallel_file_mode` you can use Multiple Independent File (MIF) or Single Shared File (SIF) and specify the number of HDF5 files right after this parameter (see example below).

#### 9.1.1 JSON Configuration (recommended)

To run an instance of MACSio HDF5 benchmark you need to include the following in the `benchmarks` property of your `configuration.json` file:

```
{
  "benchmark": "macsio",
  "file": "test.h5",
  "configuration": {
    "parallel_file_mode": "MIF 8",
    "part_size": "1M"
  }
}
```



## WAYS TO CONTRIBUTE

We appreciate your interest in **h5bench**, and thank you for taking the time to contribute!

We have compiled a set of instructions to help us make h5bench even better.

### 10.1 Reporting bugs

You can open a new issue using our GitHub [issue tracker](#). If you run into an issue, please search first to ensure the issue has not been reported before. Open a new issue only if you have not found anything similar to your issue. Please, try to provide as much information as possible to reproduce your bug quickly.

### 10.2 Suggesting enhancements

You can use our GitHub [issue tracker](#) to describe your proposed feature. Please, provide the necessary context, covering why it is needed and what problem does it solve.

### 10.3 Adding new benchmarks

We provide a set of instructions on how to add new benchmarks to the h5bench Benchmarking Suite. However, please notice that you might require some changes depending on how your benchmark work. You can contribute in two ways:

- **Adding existing benchmarks as submodules:** We plan to support only the version included in the original PR, based on its commit hash. Updates on the submodule require the contributor's help to ensure we can smoothly upgrade the available version without breaking existing features (both in the benchmark and in h5bench).
- **Adding newly developed benchmarks:** The community may perform maintenance, requiring you to provide comprehensive documentation (in code and usage) and examples to understand the benchmark module.

#### 10.3.1 Example

To illustrate how you can add a new benchmark using the submodule approach we will use AMReX:

1. You need to include the AMReX repository as a submodule:

```
git submodule add https://github.com/AMReX-Codes/amrex amrex
```

2. For this benchmark, we need some libraries to be compiled and available as well, so we will need to modify our `CMakeLists.txt`, so it builds that subdirectory:

```
set(AMReX_HDF5 YES)
set(AMReX_PARTICLES YES)
set(AMReX_MPI_THREAD_MULTIPLE YES)
add_subdirectory(amrex)
```

- AMReX comes with several other benchmarks. Still, since we are only interested in the HDF5 one, we will only compile that code. For that, we will need to add the following to our `CMakeLists.txt`. This is based on how that benchmark is normally compiled within AMReX.

```
set(amrex_src amrex/Tests/HDF5Benchmark/main.cpp)
add_executable(h5bench_amrex ${amrex_src})
```

- Be sure to follow the convention of naming the executable as `h5bench_` plus the benchmark name, e.g. `h5bench_amrex`.
- If you are going to provide support for the HDF5 async VOL connector with explicit implementation (which require changes in the original code), make sure you link the required libraries (`asynchdf5` and `h5async`):

```
if(WITH_ASYNC_VOL)
    set(AMREX_USE_HDF5_ASYNC YES)
    target_link_libraries(h5bench_amrex hdf5 z m amrex asynchdf5 h5async MPI::MPI_C)
else()
    target_link_libraries(h5bench_amrex hdf5 z m amrex MPI::MPI_C)
endif()
```

- The last step is to update the `h5bench` Python-based script to handle the new benchmark. On the top of the file, add the path of your benchmark:

```
H5BENCH_AMREX = 'h5bench_amrex'
```

Update the `run()` function that iterates over the `benchmarks` property list defined by the user in the `configuration.json` file to accept the new benchmark name:

```
elif name == 'amrex':
    self.run_amrex(id, benchmark[name], setup['vol'])
```

You then need to define the `run_` function for the benchmark you're adding. The most important part is translating the configuration defined in the `configuration.json` file into a format accepted by your benchmark (e.g., a file, a JSON, command line). For AMReX, it requires an `amrex.ini` file with key-value configurations defined in the format `key = value`, one per line:

```
# Create the configuration file for this benchmark
with open(configuration_file, 'w+') as f:
    for key in configuration:
        f.write('{} = {}\n'.format(key, configuration[key]))

    f.write('directory = {}\n'.format(file))
```

If you plan to support the HDF5 async VOL connector, make sure you can `enable_vol()` and `disable_vol()` at the beginning and end of this `run_` function.

Here you can check an example of the complete `run_amrex` function:

```
def run_amrex(self, id, setup, vol):
    """Run the AMReX benchmark."""
```

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```

self.enable_vol(vol)

try:
    start = time.time()

    file = '{}/{}'.format(self.directory, setup['file'])
    configuration = setup['configuration']

    configuration_file = '{}/{}amrex.ini'.format(self.directory, id)

    # Create the configuration file for this benchmark
    with open(configuration_file, 'w+') as f:
        for key in configuration:
            f.write('{} = {}\n'.format(key, configuration[key]))

        f.write('directory = {}\n'.format(file))

    command = '{} {} {}'.format(
        self.mpi,
        self.H5BENCH_AMREX,
        configuration_file
    )

    self.logger.info(command)

    # Make sure the command line is in the correct format
    arguments = shlex.split(command)

    stdout_file_name = 'stdout'
    stderr_file_name = 'stderr'

    with open(stdout_file_name, mode='w') as stdout_file, open(stderr_file_name,
↪mode='w') as stderr_file:
        s = subprocess.Popen(arguments, stdout=stdout_file, stderr=stderr_file,
↪env=self.vol_environment)
        sOutput, sError = s.communicate()

        if s.returncode == 0:
            self.logger.info('SUCCESS')
        else:
            self.logger.error('Return: %s (check %s for detailed log)', s.returncode,
↪ stderr_file_name)

            if self.abort:
                self.logger.critical('h5bench execution aborted upon first error')

                exit(-1)

    end = time.time()

    self.logger.info('Runtime: {:.7f} seconds (elapsed time, includes allocation,
↪wait time)'.format(end - start))

```

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```
except Exception as e:
    self.logger.error('Unable to run the benchmark: %s', e)

self.disable_vol(vol)
```

7. Make sure you provide some sample JSON configuration files in the `configurations` directory.

Please, feel free to reach us if you have questions!

## 10.4 Testing

h5bench constantly receives updates and improvements. If you can run the latest version, please consider helping us by reporting your findings, including bugs and performance regressions. Running the benchmarks contained in the h5bench Benchmarking Suite with different configurations and platforms helps us a lot in making it more robust by quickly identifying and solving issues.

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