

# UNIVERSITY OF A CORUÑA Computer Architecture Group

# MapReduce Evaluator 1.1: User Guide

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## Contents

1	Overview	3			
2	Features         2.1       User-defined parameters         2.2       Galaxie	<b>3</b> 3			
	2.2 Solutions and benchmarks	4			
3	Example	<b>5</b>			
	3.1 Configuration of the experiments	5			
	3.2 Execution	8			
	3.3 Results	8			
	3.3.1 Log & configuration $\ldots$	8			
	3.3.2 Performance	8			
	3.3.3 Resource Utilization	8			
Α	About Open Grid Scheduler/Grid Engine	10			
в	B About Environment Modules				
С	C System Requirements				

#### 1 Overview

The MapReduce Evaluator 1.1 (MREv) is an evaluation tool to extract valuable information about the performance, scalability and resource efficiency of HPC-oriented MapReduce solutions. It allows to compare several frameworks by means of different workloads, including micro-benchmarks and real-world applications.

MREv uses multiple user-defined parameters to unify of the configuration the solutions, ensuring a fair comparison between them. In each experiment, the user can select the workloads and the solutions to be run. Several cluster sizes can be used in order to check the scalability of the frameworks and ensure an optimal use of the nodes of the system. The user can also determine the number of times each workload is executed in order to obtain statistical information.

#### 2 Features

This section discusses the user-defined parameters, solutions and benchmarks currently available in MREv. It also describes the main behaviour of the execution of the experiments.

#### 2.1 User-defined parameters

Table 1 shows the configuration parameters currently available in MREv. They are classified depending on whether they affect the configuration of the solutions, the experiments to run or the input data sets of the workloads.

System dependent	Experiment dependent	Workload dependent
Mappers per node	Benchmarks	RandomWriter/Teragen data size
Reducers per node	Cluster sizes	TestDFSIO: number of files
Java heap size (slave)	Solutions	TestDFSIO: file size
Java heap size (master)	Number of executions	PageRank: number of iterations
HDFS block size		Bayes: input directory
HDFS replication factor		Command: workload command
Temporary directory		Command: data generation command
SSD optimization		Benchmark timeout
GbE interface		
IPoIB interface		
Cores per node		
Memory per node		
Hadoop specific paramete	ers	Mahout specific parameters
io.sort.factor		Mahout heap size
io.sort.mb		
shuffle.parallelcopies		
MREv parameters		
Enable plots		
Enable stat records		
Default timeout		
MREv out directory		

Table 1: Available MREv configuration parameters

#### $\mathbf{2.2}$ Solutions and benchmarks

Table 2 and 3 show the frameworks, and the benchmarks, respectively, currently supported by MREv. All benchmark are available for Hadoop-based solutions, although DataMPI only supports WordCount, Sort and TeraSort, provided with the distribution. Benchmarks for Spark are still under development.

Solution	Release Date	Interconnect
Hadoop-1.2.1-GbE	01/08/2013	Gigabit Ethernet
Hadoop-1.2.1-IPoIB	01/08/2013	InfiniBand (IPoIB)
Hadoop-1.2.1-UDA	07/06/2013	InfiniBand (RDMA & IPoIB)
Hadoop-2.5.1-GbE	12/09/2014	Gigabit Ethernet
Hadoop-2.5.1-IPoIB	12/09/2014	InfiniBand (IPoIB)
Hadoop-2.5.1-UDA	03/09/2014	InfiniBand (RDMA & IPoIB) <sup>a</sup>
Hadoop-2.7.1-GbE	06/07/2015	Gigabit Ethernet
Hadoop-2.7.1-IPoIB	06/07/2015	InfiniBand (IPoIB)
Hadoop-2.7.1-UDA	06/07/2015	InfiniBand (RDMA & IPoIB) <sup>a</sup>
RDMA-Hadoop-0.9.9-GbE	31/03/2014	InfiniBand (RDMA & Gigabit Ethernet)
RDMA-Hadoop-0.9.9-IPoIB	31/03/2014	InfiniBand (RDMA & IPoIB)
RDMA-Hadoop-2-0.9.7-GbE	26/05/2015	InfiniBand (RDMA & Gigabit Ethernet)
RDMA-Hadoop-2-0.9.7-IPoIB	26/05/2015	InfiniBand (RDMA & IPoIB)
$DataMPI-0.6.0-HDFS-GbE^{b}$	16/04/2014	InfiniBand (RDMA) & Gigabit Ethernet
$DataMPI-0.6.0-HDFS-IPoIB^{b}$	16/04/2014	InfiniBand (RDMA & IPoIB)
Spark-1.4.1-YARN-GbE	26/07/2015	Gigabit Ethernet
Spark-1.4.1-YARN-IPoIB	26/07/2015	InfiniBand (IPoIB)

Table 2: Supported frameworks

 $^{\rm a}$  Mellanox UDA uses RDMA communications in the shuffle phase and IPoIB for HDFS  $^{\rm b}$  DataMPI uses HDFS (Hadoop 1.2.1) as distributed filesystem configured with IPoIB

Micro-benchmarks						
TestDFSIO	Tests the read and write throughput of HDFS by generating a large number					
	of tasks performing reads and writes simultaneously					
Wordcount	Counts the number of times each word appears in the input text data set,					
	which is set up using the RandomTextWriter data generator					
Sort	Sorts the input text data set, also generated by RandomTextWriter					
TeraSort	Sorts 100B-sized $\langle key, value \rangle$ tuples, generated by the TeraGen data					
	generator. Each key is 10B-sized and each value is 90B-sized					
Real-world A	Applications					
PageRank	Ranks websites by counting the number and quality of the links to each one.					
	Developed by Google, it is used to obtain Google search results					
Bayes	Performs a classification algorithm, based on Bayes' Theorem					
Other						
Command	Executes a user-defined action, which can be set in the MREv configuration.					
	If it is not specified, MREv opens an interactive shell, which is useful for					
	debugging and interacting with other Hadoop components like HDFS.					

 Table 3: Supported benchmarks

### 3 Example

This Section describes a practical use case of MREv, including the configuration of user-defined parameters, the execution of the experiments and some comments about the provided results.

#### 3.1 Configuration of the experiments

The configuration of a experiment affects the following files:

- mrev-conf.sh
- $\bullet\,$  system-conf.sh
- $\bullet \ {\rm experiment-conf.sh}$
- solutions.lst
- benchmarks.lst
- $\bullet$  cluster\_sizes.lst

The environment variables and the configuration files are explained below, including the default values of the parameters.

**Environment variables** There are two environment variables that MREv uses to know where to find the configuration of the experiments. First, the **EXP\_DIR** variable determines the directory that contains the configuration files mentioned above. If this variable is not set, the value taken by default is **\$MREv\_HOME/experiment**. Second, the **HOSTFILE** variable contains the file that lists the compute nodes. The first line of the file will be the master, and the remaining lines will be the slaves. If this variable is not set, the value taken by default is **\$EXP\_DIR/hostfile**.

```
export EXP_DIR=$MREv_HOME/experiment
export HOSTFILE=$EXP_DIR/hostfile
```

**mrev-con.sh** This file contains some parameters that configure the behaviour of MREv along the experiments. First, the ENABLE\_PLOT parameter determines if MREv will generate performance graphs of the execution of the workloads. Similarly, ENABLE\_STAT is set to perform the stat recording when executing the workloads, and generating the graphs afterwards. The DEFAULT\_TIMEOUT parameter define the maximum execution time for a workload. Those which run above this limit will be killed, and the execution of the solution will be finished. Of course, this MREv feature can be disabled by setting DEFAULT\_TIMEOUT to 0. Finally, the directory where MREv will write the results of the experiment can also be configured by using the OUT\_DIR variable. If this variable is not set, the value taken by default is \$PWD/MREv\_OUT.

#!/bin/sh

export ENABLE\_PLOT="true"
export ENABLE\_STAT="true"
export DEFAULT\_TIMEOUT="8h"
export OUT\_DIR=\$PWD/\${METHOD\_NAME}\_OUT

**system-conf.sh** This file contains the parameters related to the system where MREv is being run. Some of them are automatically detected from the system, but can also be tuned by the user in order to maximize the leveraging of the system resources.

#!/bin/sh

export TMP\_DIR=/tmp/\$USER # Directory used to store local data in each node export GBE\_INTERFACE="eth1" # GbE interface to use in the nodes export IPOIB\_INTERFACE="ib0" # IPoIB interface to use in the nodes export CPUS\_PER\_NODE='grep "^physical id" /proc/cpuinfo | sort -u | wc -l' # Available CPUs per node export CORES\_PER\_CPU='grep "^core id" /proc/cpuinfo | sort -u | wc -l' # Available cores per CPU export CORES\_PER\_NODE=\$(( \$CPUS\_PER\_NODE \* \$CORES\_PER\_CPU )) # Available cores per node export MEMORY\_PER\_NODE=\$(('grep MemTotal /proc/meminfo | awk '{print \$2}''/1024)) # Available memory per node export MAPPERS\_PER\_NODE=\$(( \$CORES\_PER\_NODE / 2 - 1 )) # Maximum number of map tasks per node export REDUCERS\_PER\_NODE=\$(( \$CORES\_PER\_NODE / 2 - 1 )) # Maximum number of reduce tasks per node export HEAPSIZE='op\_int "\$MEMORY\_PER\_NODE \* 0.75 / ( \$MAPPERS\_PER\_NODE + \$REDUCERS\_PER\_ NODE )"' # Heap volume size per map/reduce task (MB) export MASTER\_HEAPSIZE='op\_int "\$MEMORY\_PER\_NODE / 4"' # Heap volume size per master daemon (MB) export BLOCKSIZE=\$((128\*1024\*1024)) # HDFS block size (Bytes) export REPLICATION\_FACTOR=3 # Number of block replications export SSD="false" # Enable SSD optimization (only applies to RDMA-Hadoop) #Hadoop specific parameters export IO\_SORT\_FACTOR=100 # Number of streams to merge at once while sorting files export IO\_SORT\_MB=\$(( \$HEAPSIZE / 3 )) # Total amount of buffer memory to use while sorting files (MB) export SHUFFLE\_PARALLELCOPIES=20 # Default number of parallel transfers run by reduce during the copy(shuffle) phase #MAHOUT export MAHOUT\_HEAPSIZE=\$MASTER\_HEAPSIZE # Heap volume size for Mahout master process ( MB)

**experiment-conf.sh** This file sets the problem size of the benchmarks and the number of times each one is executed. Moreover, it also contains the METHOD\_COMMAND variable, which contains the action to run in batch mode during the command benchmark. Additionally, METHOD\_PREPARE\_COMMAND is called to set up the input datasets needed for METHOD\_COMMAND. This enables to perform accurate performance and resource utilization monitoring, without taking into account the data generation or the copy to HDFS. This file also allows to configure specific timeouts for each benchmark.

#!/bin/sh

```
export NUM_EXECUTIONS=1 # Number of times each benchmark is executed
#Wordcount, sort & terasort
export DATASIZE=$(( 1024 * 1024 * 1024)) # Size of each input dataset (Bytes)
#TestDFSI0
export DFSI0_N_FILES=100 # Number of files to generate
export DFSIO_FILE_SIZE=100 # Size of each file (MB)
#PageRank
export PAGERANK_ITERATIONS=10 # Number of iterations to obtain the results
#Bayes
export BAYES_INPUT=$PWD/wikixml # Input data set (optional)
## Command
# export METHOD_COMMAND= # Command to run in batch mode
# export METHOD_PREPARE_COMMAND= # Command to run to set up input datasets
## TIMEOUT
# export TESTDFSI0_TIMEOUT="0"
# export WORDCOUNT_TIMEOUT="0"
# export SORT_TIMEOUT="0"
# export TERASORT_TIMEOUT="0"
# export PAGERANK_TIMEOUT="0"
# export BAYES_TIMEOUT="0"
# export COMMAND_TIMEOUT="0"
```

solutions.lst This file contains the solutions to be used in the experiment.

#Hadoop-1.2.1-GbE #Hadoop-1.2.1-IPoIB #Hadoop-1.2.1-UDA #Hadoop-2.5.1-GbE #Hadoop-2.5.1-IPoIB #Hadoop-2.5.1-UDA Hadoop-2.7.1-GbE #Hadoop-2.7.1-IPoIB #Hadoop-2.7.1-UDA #RDMA-Hadoop-0.9.9-GbE #RDMA-Hadoop-0.9.9-IPoIB #RDMA-Hadoop-2-0.9.7-GbE #RDMA-Hadoop-2-0.9.7-IPoIB #DataMPI-0.6.0-HDFS-GbE #DataMPI-0.6.0-HDFS-IPoIB #Spark-1.4.1-YARN-GbE #Spark-1.4.1-YARN-IPoIB

benchmarks.lst This file contains the benchmarks to be used in the experiment.

```
#testdfsio # Tests the read and write throughput of HDFS
wordcount # Counts the number of times each word appears in the input data set
#sort # Sorts the input data set
#terasort # Sorts 100B-sized < key, value > tuples
#pagerank # Ranks websites by counting the number and quality of the links to each one
#bayes # Performs a classification algorithm, based on Bayes' Theorem
#command # Executes user-defined actions (interactive or batch)
```

**cluster\_sizes.lst** This file contains the cluster sizes with which the user wants to run the experiments. Additionally, the cluster size can be set to the maximum number of nodes available.

#3 #5 #9 #13 MAX

#### 3.2 Execution

The following command starts the experiments:

bash MREv/bin/run.sh

#### 3.3 Results

The results from the execution will be found in the **\$OUT\_DIR** directory, having the structure shown in Figure 1.

#### 3.3.1 Log & configuration

MREv creates separate log and configuration directories for each framework and stores them at {cluster\_size}/{framework}. For example, the configuration directory of Hadoop-2.7.1-IPoIB with 5 nodes is report\_MREv\_08\_10\_12-00-00/5/Hadoop-2.7.1-IPoIB/etc/hadoop and its log directory is report\_MREv\_08\_10\_12-00-00/5/Hadoop-2.7.1-IPoIB/log. Both directories can be used to check the configuration generated by MREv and the execution of the workloads. Moreover, this feature enables to run simultaneous evaluations of the same framework using different configurations.

#### 3.3.2 Performance

The performance results in terms of time are available in the graphs subdirectory. For example, for the Wordcount benchmark, they can be found in the report\_MREv\_08\_10\_12-00-00/graphs/wordcount.eps file. For each cluster size, the graph depicts the average, maximum and minimum execution times taken by each framework to perform the workload.

#### 3.3.3 Resource Utilization

The resource utilization results from the execution of a benchmark can be found at {cluster\_size}/{framework}/{benchmark}\_{num\_execution}/stat\_records. For example, the values of the first execution of Wordcount using Hadoop-2.7.1-IPoIB on 5 nodes are at report\_MREv\_08\_10\_12-00-00/5/Hadoop-2.7.1-IPoIB/wordcount\_1/stat\_records. This directory contains one subdirectory for the values of each cluster node, plus another one for the average

MREv_OUT
report_MREv_08_10_12-00-00
Log Execution log
_ summary Experiments configuration and main results
5Output directory for cluster size 5
Hadoop-2.7.1-GbEOutput directory for Hadoop-2.7.1-GbE
Hadoop-2.7.1-IPoIB Output directory for Hadoop-2.7.1-IPoIB
etc
hadoop Hadoop configuration directory
Logs
wordcount_1Output directory for the 1 <sup>st</sup> execution of Wordcount
elapsed_timeElapsed seconds
output
stat_records
log Stat graphs generation log
avg Average statistics directory
CPU utilization results graph
dsk_sda_rw_stat.epsDisk read/write results graph
dsk_sda_util_stat.epsDisk utilization results graph
mem_stat.eps
net_ib0_stat.eps IPoIB network results graph
_ cpu_stat.datCPU results data file
node-0Node 0 (master) statistics directory
wordcount_2 Output directory for the $2^{nd}$ execution of Wordcount
wordcount_2 Output directory for the 2 execution of wordcount
Hadoop-2.7.1-UDAOutput directory for Hadoop-2.7.1-UDA
9Output directory for cluster size 9
graphs
LogGraph generation log
wordcount.eps
wordcount.dat Time results for the Wordcount benchmark (data file)

Figure 1: MREv output directory structure

values among the slave nodes. The resource utilization graphs include CPU utilization (cpu\_stat.eps), CPU load (cpu\_load\_stat.eps), memory usage (mem\_stat.eps), disk read/write (dsk\_sda\_rw\_stat.eps), disk utilization (dsk\_sda\_util\_stat.eps) and network send/recv (net\_eth1\_stat.eps, net\_ib0\_stat.eps). Disks (sda) and network interfaces (eth1, ib0) are automatically detected by MREv. For some resources, like CPU utilization, there are different visualization modes that allow to see the results individually (with lines, cpu\_stat.eps) or as a whole (with stacked values, cpu\_stat\_stacked.eps).

### A About Open Grid Scheduler/Grid Engine

As most supercomputers use a batch-queuing system for distributed resource management, MREv is aware of the environment variables and connection constraints that are typically present in these cases. The correct behaviour of MREv under this kind of systems has been tested with the Open Grid Scheduler/GE (OGS/GE).

MREv will detect the PE\_HOSTFILE environment and use it to read the compute nodes. In this case, no HOSTFILE variable will be needed, although it can also be set. Moreover, the ssh connections used to launch DataMPI and Hadoop daemons do not work properly under OGS/GE, so MREv performs a light them to enable their execution.

### **B** About Environment Modules

MREv is aware of the use of Modules for dynamically modifying the user's environment. If available, MREv will use it for loading the Java and MPI environment variables.

### C System Requirements

The following packages need to be installed:

- Gnuplot 4.4
- Java JRE 1.7
- $\bullet\,$  MPI  $^1$

 $<sup>^1\</sup>mathrm{DataMPI}$  has been tested using Mvapich2