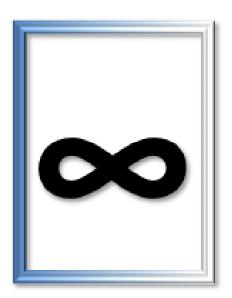
amask v1.0

AMASK: Affinity MASKs for parallel processes/threads Document Revision 1.0 June 11, 2017



Kent Milfeld milfeld@tacc.utexas.edu

High Performance Computing Texas Advanced Computing Center The University of Texas at Austin

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Abstract

Amask is a set of tools for application developers and users to discover the affinity masks of application processes (MPI ranks or OpenMP threads) so they can determine where the processes can run. Amask has the following components:

- Stand-alone executables to report default masks of OpenMP, MPI or Hybrid executions in an interactive or batch environment.
- API for instrumenting applications to report affinity masks from within a program.
- Utilities: timers, set process/thread affinity, create loads (for top viewing)

Our intention is to create a tool that provides simple-to-understand affinity information. Bug reports and feedback on usability and improvements are welcome; send to milfeld@tacc.utexas.edu with amask in the subject line.

If you use amask, cite:

github.com/TACC/amask, "amask: Affinity Mask", Texas Advanced Computing Center (TACC), Kent F. Milfeld. [1]

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1 Installation

Amask is easy to build. Execute make to build the stand-alone executables and a library for those who want to instrument their application.

Download the github repository by clicking on the "Download ZIP" file, and expanding it in a convenient location.

```
unzip amask-master.zip
```

You can also clone the git repository:

```
git clone https://github.com/TACC/amask
```

This will create a top level directory called amask, with subdirectories docs and src. Change directory to amask. Edit the Makefile to include an appropriate MPI compiler and OpenMP flag (the defaults are for Intel). Execute make.

The executables will be placed in the amask/bin directory and the library will be placed in amask/lib. Include .../amask/bin in your PATH variable.

¹https://github.com/TACC/amask

2 Process Affinity Masks for a Parallel Execution

Execute one of the stand-alone executables, amask_omp, amask_mpi, or amask_hybrid in an OpenMP, MPI or hybrid environment, respectively, to obtain the expected affinity mask for each process/thread for your program in the same environment. That is, before you execute your program, execute the appropriate stand-alone affinity program to get a listing of the masks for each OpenMP thread and/or MPI process that would occur for your program execution (provided the application doesn't adjust affinity).

In Listing 2.1 a batch job runs the amask_mpi executable before running an application, to observe the affinity mask for each rank that the application (my_mpiapp) will have. By removing the my_mpiapp execution line and requesting less time for the batch job, you can quickly discover the default affinity mask for the parallel environment. You can adjust affinity environment variables and quickly assess their impact on the affinity masks for your application.

Listing 2.1: Listing affinity masks for MPI environment

If a site allows users interactive access to batch nodes (see idev utility @TACC), then the amask commands can be run interactively. Listing 2.2 shows interactive executions to discover the affinity masks that any pure OpenMP, pure MPI or a hybrid application would have for the environment. (The environment includes the OMP variables, number of MPI tasks requested, and MPI affinity environment variables set in the mpi launcher: mpirun, or ibrun at TACC.)

Listing 2.2: Listing OpenMP/MPI and Hybrid masks

```
1 export OMP_NUM_THREADS=2; amask_omp # pure OpenMP
3 4 mpirun -np 4 amask_mpi # pure MPI
5 6 export OMP_NUM_THREADS=2; mpirun -np 4 amask_hybrid # Hybrid
```

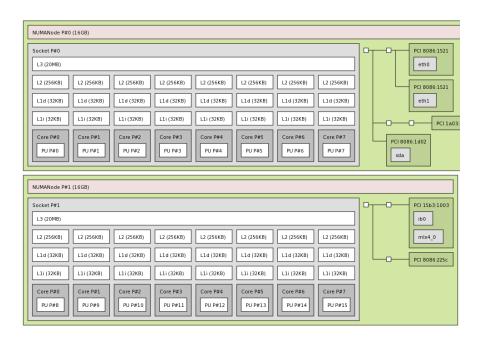


Figure 2.1: Hardware information from Istopo

How to read the output:

The output in Listing 2.3 is for an 8-rank MPI execution on a platform with 2 sockets, 8 cores/socket, no hyperthreading, and processor ids 0-7 on socket 1 and 8-15 on socket 2. One can derive these details from 1stopo. Figure 2.1 shows an 1stopo report for this system (Sandy Bridge compute node on the TACC Stampede machine.) The same information can be extracted from the contents of the /proc/cpuinfo file. The 1scpu utility provides the number of sockets, cores and SMT threads, without proc-id assignments.

Listing 2.1 shows MPI processes executing on processors (cores) 0-7. The rows of the matrix are labeled by the MPI ranks and the columns represent the machine's processor-ids (Core P# in Figure 2.1) from 0 to Nprocs-1. For this run there are 8 rows, one for each of the 8 MPI ranks (0000 through 0007). Since there are 16 processors (Nprocs), there are 16 character locations in each row, representing the mask bits for processors 0 - 15. A digit in a row position means that the mask bit is set for the the corresponding processor, and the MPI process rank listed at the beginning of the row can execute there. When multiple numbers are in a row (multiple bits set) it means that the rank can run (float) on any of those processors. (The - characters indicate the mask bits are not set.) The processor id number is determined by adding the digit in the row to the "group" number directly above in the header. (This quirky approach allows us to represent a processor's mask as a set of single characters (digits) for each bit, and allows one to determine the processor id (proc-id) number by a "look up" – adding the digit in the row to header group value above).

Listing 2.3: Default MPI Environment

```
# 8 tasks requested, MVAPICH2 MPI
2
3
   $ mpirun amask_mpi
                       # non-TACC
 4
   $ ibrun amask_mpi
                       # @TACC
5
6
        Each row of matrix is an affinity mask.
        A set mask bit = matrix digit + column group # in |...|
8
9
   rank | 0 | 10
10
   0000 0--
   0001 -1-----
11
   0002 --2-----
12
   0003 ---3-----
13
   0004 ----4-----
                         #digit + group # in header = proc-id
14
15
   0005 ----5----
   0006 -----6-----
16
   0006 -----7-----
17
```

The output in Listing 2.4 is for the same platform as 2.3, but the affinity environment has been changed with the tacc_affinity script on the ibrun launcher line (ibrun tacc_affinity amask_mpi). There are other ways to change the affinity, such as using another launcher, compiling with a different mpi library, or setting MPI environment variables. In this case tacc_affinity positions the first 4 tasks on socket 0, and allows each rank to float on any of the processors (cores) of the socket, by setting the first 8 bits of the mask for each of the 4 ranks, 0-3. Ranks 4-7 execute on the second socket and are allowed to float across all cores (proc-ids 8-15) of the socket. (tacc_affinity is not part of the amask utilities.)

Listing 2.4: TACC Tailored Environment through tacc-affinity

```
# 8 tasks requested, MVAPICH2 MPI
     # + tacc_affinity -- sets MPI affinity @TACC
3
4
   $ ibrun tacc_affinity amask_mpi
5
6
        Each row of matrix is an Affinity mask.
        A set mask bit = matrix digit + column group # in |...|
8
9
            0
                 1 10
    rank |
   0000 01234567----
10
                               # -- mask to run anyplace on socket1
   0001 01234567-----
11
                               \# rank 1 can execute on any of 0-7 cores
   0002 01234567----
                               \# rank 2 can execute on any of 0-7 cores
12
13
   0003 01234567-----
14
   0004 -----89012345
                               # -- mask to run anyplace on socket2
   0005 -----89012345
                               \# rank 5 can execute on any of cores 8-15
15
16
   0006 -----89012345
                               # Bits are for proc-ids (cores) 8,9,10-15
17
   0007 -----89012345
                               # Read as 8+0, 9+0, 0+10, 0+10, ... 10+15
```

A more challenging situation is the determination of the masks on a system with hundreds of hardware threads, such as the Intel KNL processor. At TACC, Stampede2 has 4,200 PHI 7250 processors (KNLs), each with 272 hardware threads (68 cores x 4 SMT threads/core). Before version 1.0 amask would

display rows of 272 proc-ids (processor ids=hardware threads), but that became a bit unwieldy for users to display on laptops. Now, by default, a matrix of core-ids versus process-ids is displayed. If there are n SMTs/core, there will be n rows for each process (rank/thread-id) listed (on the left). For the KNL described above there will be 4 rows displayed for each process (corresponding to proc-id 0-67, 68-135, 136-203, and 204-272) with the header displaying core-ids. This is shown in Listing 2.5.

Listing 2.5: Process masks for an 8-thread execution on 68-core KNL

```
1
 2
     export OMP_NUM_THREADS=8
3
   $ export OMP_PROC_BIND=close
4
   $ ./amask_omp
5
6
        Each row of matrix is a mask for a Hardware Thread (hwt).
 7
        CORE ID = matrix digit + column group # in |...|
 8
        A set mask bit (proc-id) = core id + add 68 to each additional row.
9
10
                                       30
11
   0000 0=
12
13
14
15
16
17
18
19
20
21
22
23
24
25
26
27
28
29
30
31
   32
33
34
35
36
37
38
39
40
41
42
```

This display is basically core occupation – something users found they would rather see than 272-character masks. For each process the digits of all hardware-thread rows combine with the group number in the header to give the core id. For a process, the sequence of rows corresponds to hardware thread ids, 0-(n-1),

where n is the number of hardware threads for each core. To obtain the proc-id for a particular row entry, add a multiple of the number of cores (68 in this case) to the core id, where the multiplier is the thread id number. E.g. In Listing 2.5 the proc-id for rank 6 is 1 (the core id) plus 2x68 (hardware thread id 2 x number of cores). When you need to work directly with the mask to determine the proc-ids, it is much easier to use the mask display by using the -ls option on command line.

2.1 amask Options

There are several useful options that can be used on the omp-, mpi-, and hybrid-affinity execution command lines.

By default, after the masks are printed to standard out, a load is created on each process for 10 seconds so that the location of the processes can be viewed with the top utility. The duration of the load can be changed with the wait option (w). Include the number of seconds (integer) as an argument to the option.

By default there is a delay between printing each map. The delay can be eliminated by specifying the fast argument for the printing option, -pf.

As described above, an argument of s with the listing -1 option forces the matrix listings to be Simultaneous Multi-Threads versus processes—that is, affinity masks are listed with proc-ids.

When SMT is detected the listing defaults to a core versus process listing (with a separate line for each hardware thread in the core). Add the header group number to the digit in each row to obtain the core-id.

The basic operations used for reporting masks by the amask executables were collected into a library, so that users could instrument there own applications to display the masks of MPI processes and OpenMP threads.

Purpose option args long option long args Separate lines for SMTs -listby -1 smt Print speed f|s -printspeed fast|slow -p Wait secs (process load) -waitsecs -w # # -h help -help

Table 2.1: amask Options for amask Executables

3 Using the amask Library

3.1 How to Print amask Reports Inside an Application

To report the masks from inside a program, include the amask API routine, amask_omp(), amask_mpi(), or amask_hybrid(), within an OpenMP parallel region, after MPI has been initialized, or within an OpenMP region of a hybrid program, respectively. These functions can be called from C, C++, and Fortran; and don't require any arguments or include files:

- amask omp()
- amask mpi()
- amask hybrid()

However, to report masks for hybrid code, it may be necessary to initialize MPI with the MPI_Init_thread() routine. View the amask codes to see how easy it is to include them in your own program. The snippets below show how they are to be used:

Listing 3.1: Invoking mask report inside code

```
Pure OpenMP code
 2
        #pragma omp parallel
 3
 4
           amask omp();
5
 6
7
 8
     // Pure MPI code
9
        MPI_Init(NULL, NULL);
10
11
           amask_mpi();
12
        MPI_Finalize();
13
14
15
     // Hybrid code
16
        MPI_Init_thread(NULL, NULL, MPI_THREAD_MULTIPLE, &provided);
17
18
       #pragma omp parallel
19
20
          amask_hybrid();
21
22
23
24
       MPI_Finalize();
```

3.2 Useful Utilities

- load cpu nsec(nsec) Puts load on process/thread for nsec seconds
- map to procid(proc-id) Sets process/thread to execute on proc-id
- gtod timer() Easy to use Get Time of Day clock
- tsc() Returns time stamp counter value

A thread or process that calls load_cpu_nsec(int nsec) will execute integer operations (a load) for nsec seconds. nsec must be zero or a positive integer. Use the cmdln_get_nsec_or_help(int * nsec, int argc, char * argv[]) function to extract an integer from the command line for nsec (e.g. mpi amask_mpi 10).

A thread or process that calls map_to_procid(int proc-id) will assign the calling thread or process to execute on proc-id, by setting the appropriate bit in the scheduling mask. For example, in a parallel region the unique thread-id returned from omp_get_thread_num() can be used in an arithmetic operation (linear, modulo, etc.) to create a unique proc-id to be executed on

The function $gtod_timer()$ returns a double precision number with the number of wall-clock seconds since the previous call. The first call sets the time to zero. The function uses the Unix gettimeofday utility, and can be called from C/C++ and Fortran. See comments in the code for more details.

The function tsc() returns an 8-byte integer (Unix) time stamp count from the rdtsc instruction. Use this to capture the difference between the counts for determining the cost of a small set of operations (instructions, code statements). This is not meant to be used as a general timer when the processor frequency may change.

3.3 Other things you should know

Build amask with the same MPI library that your application will use. On systems with multi-vendor MPI libraries, build amap for each vendor version you will use. With IMPI (Intel MPI) you can set I_MPI_DEBUG=4 to separately report the mask (and other information) for each MPI process.

Bibliography

[1] A mask is a set of executables and routines for reporting affinity information. ${\rm https://github.com/tacc/amask}$