

:- HPL is a High-Performance Linpack benchmark implementation. The code solves a uniformely random system of linear equations and reports time and floating-point execution rate using a standard formula for operation count.

floating-point execution rate using a standard formula for operation count. ----- LAB —-----=> Create one machine using nat network # Turn off the firewall 1. Systemctl disable firewalld # Turn off selinux 2. vi /etc/selinux/config # This file controls the state of SELinux on the system. # SELINUX= can take one of these three values: enforcing - SELinux security policy is enforced. permissive - SELinux prints warnings instead of enforcing. disabled - No SELinux policy is loaded. SELINUX=disabled # SELINUXTYPE= can take one of three values: targeted - Targeted processes are protected, minimum - Modification of targeted policy. Only selected processes are protected. mls - Multi Level Security protection. SELINUXTYPE=targeted SELINUX='enable' to 'disabled' # do restart the system 3. Init 6 # install mathematical library Yum install blas -y Or Yum install atlas -y # Download HPL tar file https://netlib.org/benchmark/hpl/hpl-2.3.tar.qz # untar the HPL tar file tar -xvf hpl-2.3.tar.gz # Go to this path cd hpl-2.3/setup/ # see the file to all information(nothing changes)

vi Make.Linux Intel64

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for see all prebuild library files

rpm -ql atlas

```
[root@master setup]# rpm -ql atlas
/etc/ld.so.conf.d/atlas-x86_64.conf
/usr/lib64/atlas
/usr/lib64/atlas/libsatlas.so.3
/usr/lib64/atlas/libsatlas.so.3.10
/usr/lib64/atlas/libtatlas.so.3
/usr/lib64/atlas/libtatlas.so.3.10
/usr/share/doc/atlas-3.10.1
/usr/share/doc/atlas-3.10.1/README.dist
```

Download Openmpi tar file from browser

https://download.open-mpi.org/release/open-mpi/v4.1/openmpi-4.1.4.tar.gz

```
# untar openmpi file
      Tar -xvf /root/Downloads/openmpi-4.1.4.tar.gz
# Go to this file
      cd openmpi-4.1.4/
# Configure openmpi
      ./configure --prefix=/opt/openmpi-4.1.4 --enable-orterun-prefix-by-default
# if above command not run then install gcc
      yum -y install gcc gcc-c++
# Configure openmpi
      ./configure --prefix=/opt/openmpi-4.1.4 --enable-orterun-prefix-by-default
# check supporting file
       make -j 8
      The "-j" option is used with the "make" command to specify the number of
jobs (i.e., tasks or processes)
# clear all making file
      make clear all
# install make
      make install
# Export openmpi
      export PATH=/opt/openmpi-4.1.4:$PATH
# Export library
      export LD_LIBRARY_PATH=/opt/openmpi-4.1.4:$LD_LIBRARY_PATH
# copy file
      cp /opt/openmpi-4.1.4/setup/hpl-2.3 /root/Download/
# Go to path
      cd /root/Download/
# Edit this file
      vi Make.Linux_PII_CBLAS
```

```
T0Pdir
           = /root/Downloads/hpl-2.3
            = $(TOPdir)/include
INCdir
             = $(TOPdir)/bin/$(ARCH)
BINdir
LIBdir
              = $(TOPdir)/lib/$(ARCH)
HPLlib
              = $(LIBdir)/libhpl.a
TOPdir
         = /root/Downloads/hpl-2.3
MPdir
              = /opt/openmpi-4.1.4
MPinc
             = -I$(MPdir)/include
MPlib
              = $(MPdir)/lib/libmpi.so
MPdir
         = /opt/openmpi-4.1.4
MPlib
        = $(MPdir)/lib/libmpi.so
LAdir
              = /usr/lib64/atlas
LAinc
              = $(LAdir)/libsatlas.so.3 $(LAdir)/libtatlas.so.3
LAlib
LAdir = /usr/lib64/atlas
LAlib
        = $(LAdir)/libsatlas.so.3 $(LAdir)/libtatlas.so.3
cc
            = /usr/bin/qcc
CCN00PT
            = $(HPL DEFS)
CCFLAGS
            = $(HPL DEFS) -fomit-frame-pointer -03 -funroll-loops
# On some platforms, it is necessary to use the Fortran linker to find
# the Fortran internals used in the BLAS library.
LINKER = /usr/bin/gcc
LINKFLAGS = $(CCFLAGS)
CC
        = /usr/bin/gcc
LINKER
          = /usr/bin/gcc
# Go to this path
     cd /root/Downloads/hpl-2.3/bin/Linux PII CBLAS/
# check file
     Ш
[root@master Linux PII CBLAS]# ll
total 216
-rw-r--r-. 1 root root 1133 Jan 30 06:02 HPL.dat
-rwxr-xr-x. 1 root root 213576 Jan 30 06:02 xhpl
```

check all entry in HPL.dat file vi HPL.dat

```
HPLinpack benchmark input file
Innovative Computing Laboratory, University of Tennessee
HPL.out
             output file name (if any)
6
             device out (6=stdout,7=stderr,file)
             # of problems sizes (N)
29 30 34 35
             # of NBs
1 2 3 4
             NBs
0
             PMAP process mapping (0=Row-,1=Column-major)
3
             # of process grids (P x Q)
2 1 4
             Ps
2 4 1
             Qs
16.0
             threshold
3
             # of panel fact
             PFACTs (0=left, 1=Crout, 2=Right)
0 1 2
2
             # of recursive stopping criterium
2 4
             NBMINs (>= 1)
1
             # of panels in recursion
2
             NDIVs
3
             # of recursive panel fact.
0 1 2
             RFACTs (0=left, 1=Crout, 2=Right)
1
             # of broadcast
0
             BCASTs (0=1rg,1=1rM,2=2rg,3=2rM,4=Lng,5=LnM)
             # of lookahead depth
1
0
             DEPTHs (>=0)
2
             SWAP (0=bin-exch,1=long,2=mix)
64
             swapping threshold
             L1 in (0=transposed,1=no-transposed) form
0
0
             U in (0=transposed,1=no-transposed) form
1
             Equilibration (0=no,1=yes)
8~
             memory alignment in double (> 0)
```

calculate HPL Benchmarking

mpirun --allow-run-as-root -np 4 ./xhpl HPL.dat

```
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HPLinpack 2.3 -- High-Performance Linpack benchmark -- December 2, 2018
Written by A. Petitet and R. Clint Whaley, Innovative Computing Laboratory, UTK
Modified by Piotr Luszczek, Innovative Computing Laboratory, UTK
Modified by Julien Langou, University of Colorado Denver
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An explanation of the input/output parameters follows:
T/V : Wall time / encoded variant.
    : The order of the coefficient matrix A.
    : The partitioning blocking factor.
NB
    : The number of process rows.
    : The number of process columns.
Time : Time in seconds to solve the linear system.
Gflops: Rate of execution for solving the linear system.
The following parameter values will be used:
N
   : 29232
    : 232
NB
PMAP : Row-major process mapping
   : 1
    :
         1
PFACT : Right
NBMIN : 4
NDIV :
        2
RFACT : Crout
BCAST : 1ringM
DEPTH : 1
SWAP : Mix (threshold = 64)
   : transposed form
U : transposed form
EQUIL : yes
ALIGN : 8 double precision words
                      _____
- The matrix A is randomly generated for each test.
- The following scaled residual check will be computed:
    ||Ax-b||_oo / ( eps * ( || x ||_oo * || A ||_oo + || b ||_oo ) * N )
- The relative machine precision (eps) is taken to be
                                          1.110223e-16
- Computational tests pass if scaled residuals are less than
                                                       16.0
_____
-----
WR11C2R4 29232 232 1 1 74.20
                                                    2.2444e+02
```

T/V	N	NB	Р	Q		Time	Gflops
WR11C2R4	29232	232	1	1		74.20	2.2444e+02
HPL_pdgesv()	start tim	e Sat /	Aug 27	11:36:48	2022		
HPL_pdgesv()	end time	Sat A	Aug 27	11:38:02	2022		
Ax-b _oo/(eps*(A _oo* x _oo+ b _oo)*N)= 2.06977736e-03 PASSED							
Finished	1 tests	with th	he fol	lowing re	sults:		
1 tests completed and passed residual checks,							
0 tests completed and failed residual checks,							
0 tests skipped because of illegal input values.							
End of Tests.							

Result:	 	 	
===========			
=====			

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An explanation of the input/output parameters follows:

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Gflops: Rate of execution for solving the linear system.

The following parameter values will be used:

N : 29232

NB : 232

PMAP : Row-major process mapping

P : 1

Q : 1

PFACT: Right

NBMIN: 4

NDIV: 2

RFACT: Crout

BCAST: 1ringM

DEPTH: 1

SWAP : Mix (threshold = 64)

L1 : transposed form

U : transposed form

EQUIL: yes

ALIGN: 8 double precision words

- The matrix A is randomly generated for each test.
- The following scaled residual check will be computed:

- The relative machine precision (eps) is taken to be 1.110223e-16
- Computational tests pass if scaled residuals are less than 16.0

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T/V N NB P Q Time Gflops

WR11C2R4 29232 232 1 1 74.20 2.2444e+02					
HPL_pdgesv() start time Sat Aug 27 11:36:48 2022					
HPL_pdgesv() end time Sat Aug 27 11:38:02 2022					
Ax-b _oo/(eps*(A _oo* x _oo+ b _oo)*N)= 2.06977736e-03 PASSED					
======					
Finished 1 tests with the following results:					
1 tests completed and passed residual checks,					
0 tests completed and failed residual checks,					
0 tests skipped because of illegal input values.					
End of Tests.					
=======================================					