



# Intel® Omni-Path Architecture Multiple Endpoints

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# Outline

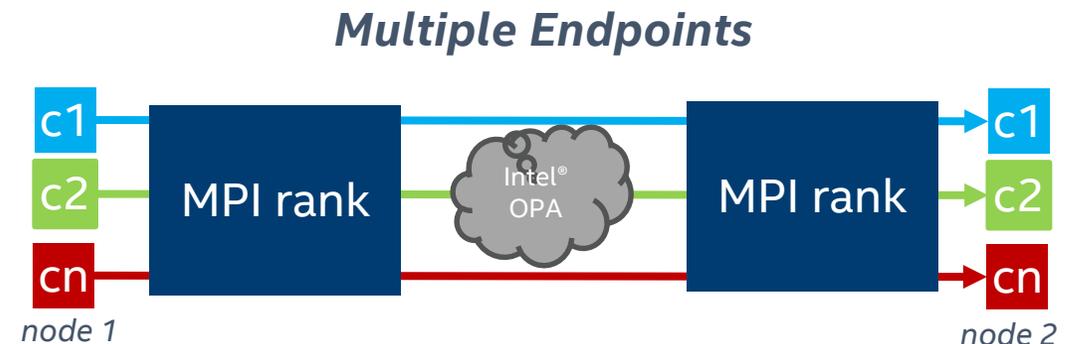
What is Multi-Endpoint (MEP)?

MEP Software Recipe

Simple Example for Allreduce

# What is Multi-Endpoint (Multi-EP)?

- Hybrid MPI/OpenMP\* codes are becoming more prevalent
- Pure MPI applications are reaching limits of scalability
- More MPI ranks = more surface to volume ratio and increased network transfer
- **Multi-EP is the ability to use more than one thread per MPI rank**
- Enables MPI communication within OpenMP regions
- Reduces fork-join<sup>1</sup> when MPI communication is required
- Increases bandwidth per MPI rank because more physical cores are used to drive the communications



1. Mattson, T. and Meadows, L., "A 'Hands-on' Introduction to OpenMP," <https://www.openmp.org/wp-content/uploads/omp-hands-on-SC08.pdf>

\*The OpenMP name and the OpenMP logo are registered trademarks of the OpenMP Architecture Review Board

# Multi-EP Software Recipe

- Requires components from:
  - Intel® Omni-Path Fabric Suite Fabric Manager (IFS) version 10.5 or newer
  - OpenFabrics Interfaces (OFI) Libfabric version 1.5 or newer
  - Intel® MPI Library 2019 (or 2019 Technical Preview)
- Example execution, using 16 nodes, 1 MPI rank per node with 4 endpoints:
  - `source /opt/intel/impi/2019.0.070/bin64/mpivars.sh release_mt -ofi_internal`
  - `export I_MPI_THREAD_SPLIT=1`
  - `export I_MPI_THREAD_RUNTIME=openmp`
  - `export PSM2_MULTI_EP=1`
  - `mpirun -np 16 -ppn 1 -hostfile 16nodes -genv I_MPI_FABRICS shm:ofi -genv OMP_NUM_THREADS=4 ./myapplication`

PSM - Performance Scaled Messaging

# Simple Example for Allreduce

```
#include <mpi.h>
#include <stdio.h>
#include <string.h>
#include <unistd.h>

int numThreads,nranks,myrank,size=1048576,niters=10240,i,ompsize
int provided; //used to confirm THREAD_MULTIPLE is supported

int main(int argc, char** argv) {

    // Initialize the MPI environment
    MPI_Init_thread(NULL, NULL, MPI_THREAD_MULTIPLE, &provided);

    // Get the number of processes
    MPI_Comm_size(MPI_COMM_WORLD, &nranks);

    // Get the rank of the process
    MPI_Comm_rank(MPI_COMM_WORLD, &myrank);
```

```
#pragma omp parallel
{
    numThreads = omp_get_num_threads();
}

int x[size],y[size];
ompsize=size/numThreads;

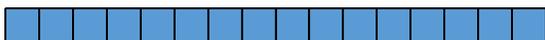
for(i=0;i<size;++i) { x[i]=i; }

MPI_Comm comm_mep[numThreads];
for(int ip=0;ip<numThreads;++ip) {
    MPI_Comm_dup(MPI_COMM_WORLD,&(comm_mep[ip]));
}
memset(y,0,sizeof(y));

#pragma omp parallel
{
    for (int iter=0;iter<niters;++iter) {
        int ip=omp_get_thread_num();
        MPI_Allreduce(x+ip*ompsize,y+ip*ompsize, ompsize, MPI_INT,
                    MPI_SUM, comm_mep[ip]);
    }
}
MPI_Finalize();
}
```

*red indicates additions necessary for Multiple Endpoints*

int x[size=1048576] array



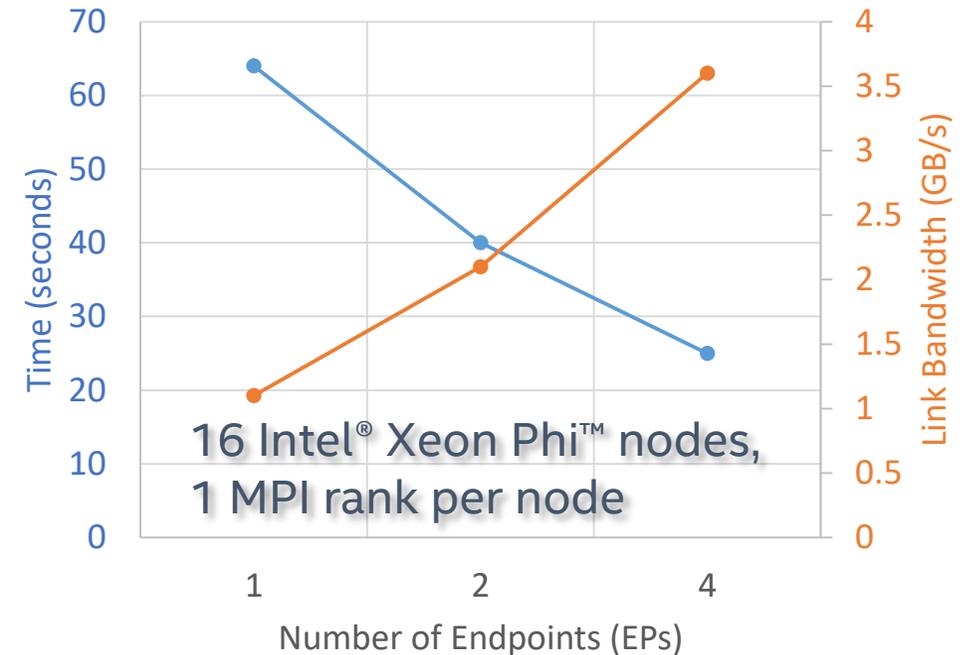
4 End points



1 2 3 4 (ompsize=262,144 each)

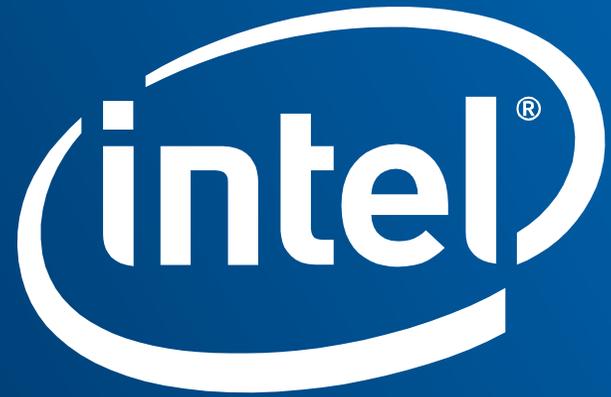
# Results & Conclusions

- **Wall clock time is reduced:**
  - 37% from 64 to 40 sec using 2 EPs
  - 60% from 64 to 25 sec using 4 EPs
- The **link bandwidth** of each node increases - higher network performance!
- Using Multi-EP with Intel® OPA only requires a **few easy steps**
- Multi-EP is being adopted by the industry to solve real world problems
  - “Multiple endpoints for improved MPI performance on a lattice QCD code” - <https://dl.acm.org/citation.cfm?id=3176375>
  - “Accelerating HPC codes on Intel(R) Omni-Path Architecture networks: From particle physics to Machine Learning” - <http://arxiv.org/abs/1711.04883>



Intel® Xeon Phi™ 7250 CPU. Intel® Turbo Boost and Hyper-Threading Technology enabled. Red Hat Enterprise Linux\* Server release 7.4 (Maipo), Kernel: 3.10.0-693.21.1.el7.x86\_64. BIOS: S72C610.86B.01.03.0018.012420182107, microcode: 0x1b6. CVE-2017-5753, CVE-2017-5715, and CVE-2017-5754 (Variants 1, 2, and 3) mitigated. Quadrant cluster mode, Flat memory mode. 16 GB MCDRAM, 96 GB DDR4 per node. Intel Fabric Suite (IFS 10.7.0.0.145). Intel® MPI Library 2019 Beta. source /opt/intel/impi/2019.0.070/bin64/mpivars.sh release\_mt -ofi\_internal; export I\_MPI\_THREAD\_SPLIT=1; export I\_MPI\_THREAD\_RUNTIME=openmp; mpirun -np 16 -ppn 1 -hostfile 16nodes -genv I\_MPI\_FABRICS shm:ofi -genv OMP\_NUM\_THREADS=4 -genv PSM2\_MULTI\_EP=1. Link bandwidth as reported by opatop tool for the highest utilized Intel® OPA Host Fabric Interface (HFI).

Performance results are based on testing as of Sept 20 2018 and may not reflect all publicly available security updates. See configuration disclosure for details. No product can be absolutely secure.



Thank you!

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