

# NAMD

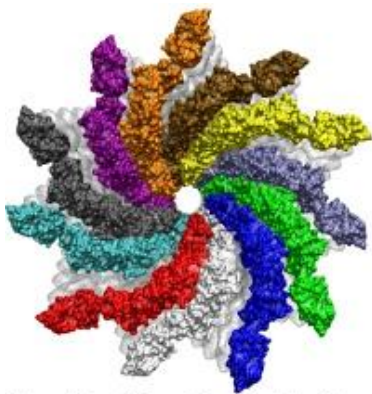
## Performance Benchmark and Profiling

February 2012

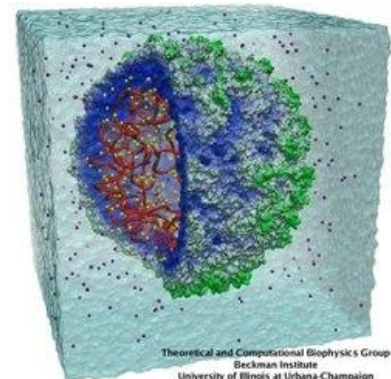
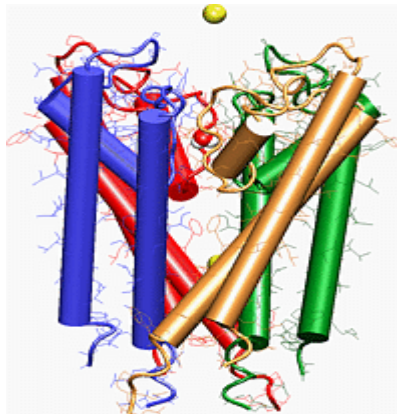


- **The following research was performed under the HPC Advisory Council activities**
  - Participating vendors: AMD, Dell, Mellanox
  - Compute resource - HPC Advisory Council Cluster Center
- **For more info please refer to**
  - [http:// www.amd.com](http://www.amd.com)
  - [http:// www.dell.com/hpc](http://www.dell.com/hpc)
  - <http://www.mellanox.com>
  - <http://www.ks.uiuc.edu/Research/namd>

- A parallel molecular dynamics code that received the 2002 Gordon Bell Award
- Designed for high-performance simulation of large biomolecular systems
  - **Scales to hundreds of processors and millions of atoms**
- Developed by the joint collaboration of the Theoretical and Computational Biophysics Group (TCB) and the Parallel Programming Laboratory (PPL) at the University of Illinois at Urbana-Champaign
- NAMD is distributed free of charge with source code



Theoretical and Computational Biophysics Group  
Beckman Institute  
University of Illinois at Urbana-Champaign



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- **The following was done to provide best practices**
  - NAMD performance benchmarking
  - Interconnect performance comparisons
  - Ways to increase NAMD productivity
  - MPI libraries comparisons
  
- **The presented results will demonstrate**
  - The scalability of the compute environment to provide nearly linear application scalability
  - The capability of NAMD to achieve scalable productivity
  - Considerations for performance optimizations

- **Dell™ PowerEdge™ R815 11-node (704-core) cluster**
- **AMD™ Opteron™ 6276 (code name “Interlagos”) 16-core @ 2.3 GHz CPUs**
- **4 CPU sockets per server node**
- **Mellanox ConnectX®-3 FDR InfiniBand Adapters**
- **Mellanox SwitchX™ 6036 36-Port InfiniBand switch**
- **Memory: 128GB memory per node DDR3 1333MHz**
- **OS: RHEL 6.2, MLNX-OFED 1.5.3 InfiniBand SW stack**
- **MPI: Open MPI 1.5.5rc2, Platform MPI 8.2**
- **Compilers: GNU Compilers 4.6**
- **Application: NAMD 2.8 (External libraries used: charm-6.2.3, fftw-2.1.3, TCL 8.3)**
- **Benchmark workload:**
  - ApoA1 bloodstream lipoprotein particle model (92,224 atoms, periodic, PME, 12A cutoff)
  - ATPase benchmark (327,506 atoms, periodic, PME)



- **HPC Advisory Council Test-bed System**
- **New 11-node 704 core cluster - featuring Dell PowerEdge™ R815 servers**
  - Replacement system for Dell PowerEdge SC1435 (192 cores) cluster system following 2 years of rigorous benchmarking and product EOL
    - System to be redirected to explore HPC in the Cloud applications
- **Workload profiling and benchmarking**
  - Characterization for HPC and compute intense environments
  - Optimization for scale, sizing and configuration and workload performance
  - Test-bed Benchmarks
    - RFPs
    - Customers/Prospects, etc
  - ISV & Industry standard application characterization
  - Best practices & usage analysis



# About Dell PowerEdge™ Platform Advantages

## Best of breed technologies and partners

Combination of AMD Opteron™ 6200 series platform and Mellanox ConnectX®-3 InfiniBand on Dell HPC

Solutions provide the ultimate platform for speed and scale

- Dell PowerEdge R815 system delivers 4 socket performance in dense 2U form factor
- Up to 64 core/32DIMMs per server – 1344 core in 42U enclosure

## Integrated stacks designed to deliver the best price/performance/watt

- 2x more memory and processing power in half of the space
- Energy optimized low flow fans, improved power supplies and dual SD modules

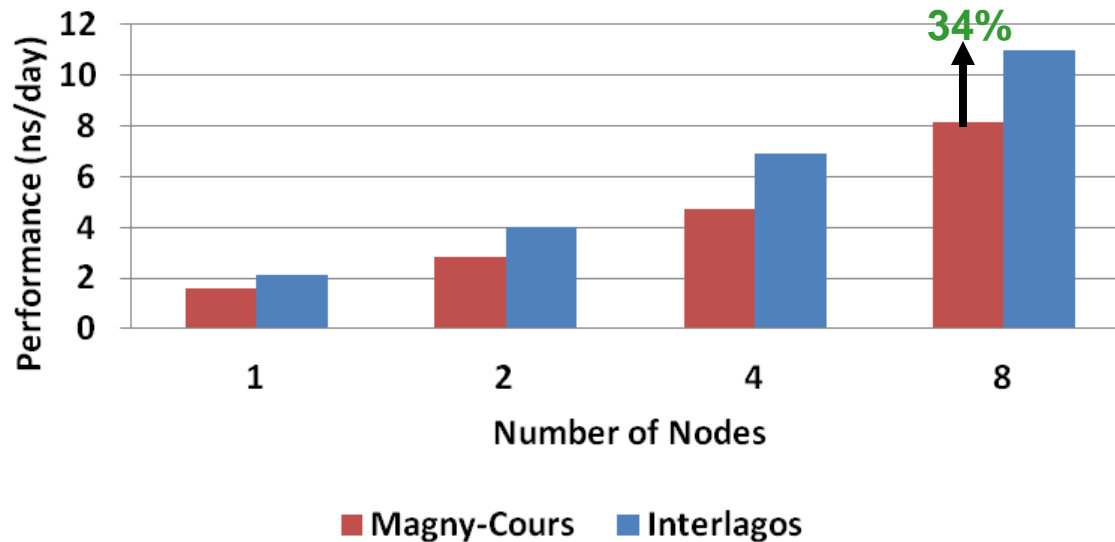
## Optimized for long-term capital and operating investment protection

- System expansion
- Component upgrades and feature releases



- **Interlagos CPUs provides better performance than Magny-Cours CPUs**
  - Up to 34% gain in performance with Open MPI versus Magny-Cours CPUs
- **Processors used:**
  - Magny-Cours: AMD Opteron™ 6174 (2200MHz)
  - Interlagos: AMD Opteron™ 6276 (2300MHz)

**NAMD Performance**  
(Apoa1, Open MPI)



*Higher is better*

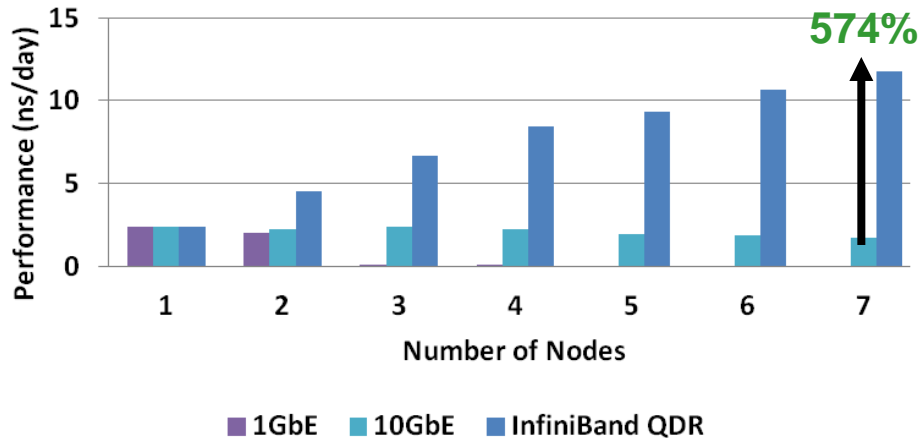
**64 Cores/Node**



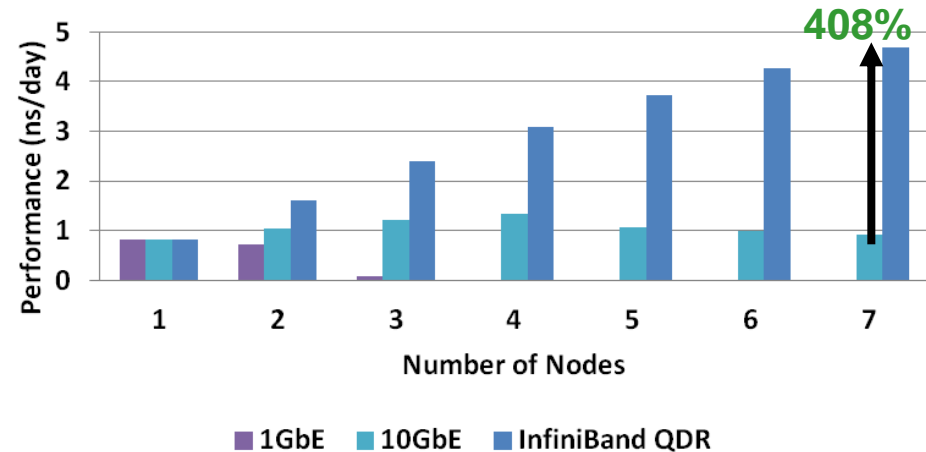
# NAMD Performance – Interconnects

- **InfiniBand enables better scalability for NAMD**
  - Showing unlimited continuous gain to 7-node
- **Ethernet does not allow good scalability**
  - The performance of 1GbE plummet after 2 nodes (128 processes)
  - Both 1GbE and 10GbE do not show gain in productivity
  - The effect of MPI communications overwhelms the Ethernet network

**NAMD Performance  
(Apoa1)**



**NAMD Performance  
(f1atpase)**

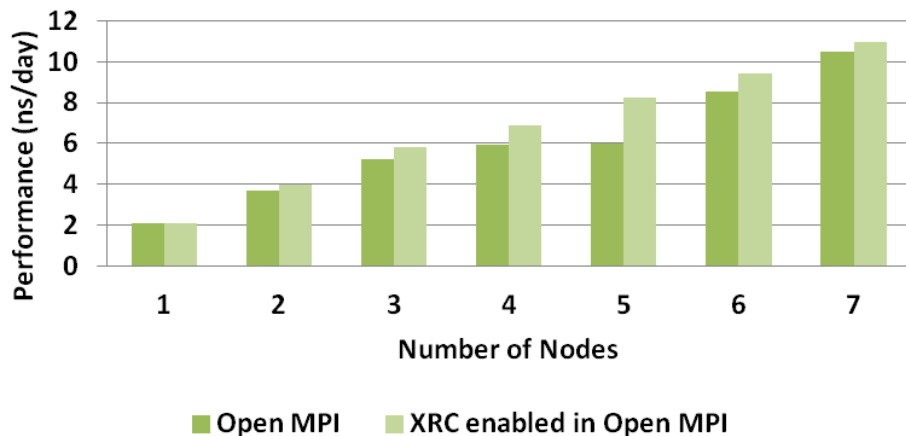


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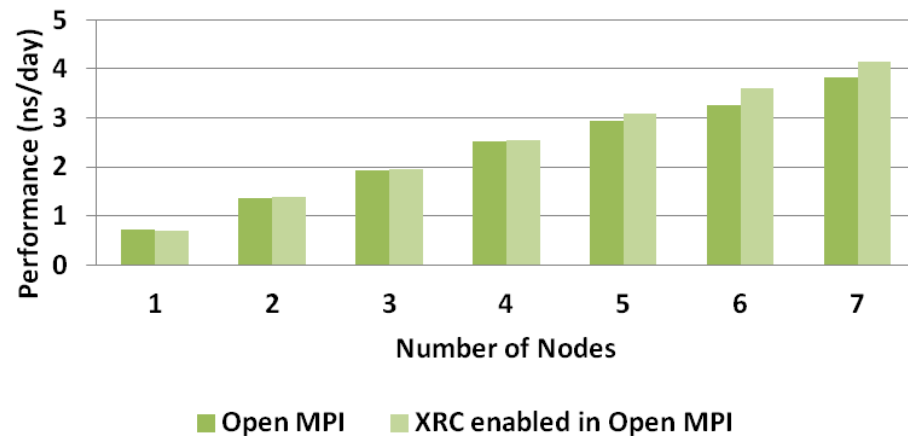
**64 Cores/Node**

- **Using XRC in Open MPI allows better performance and scalability**
  - Stands for eXtended Reliable Connection
  - Reduces memory footprint and is essential for scaling
  - Flags used: `-mca btl_openib_receive_queues X,128,256,192,128:X,2048,256,128,32:X,12288,256,128,32:X,65536,256,128,32`
- **Open MPI optimization flags used in both cases:**
  - `-bind-to-core -mca btl openib,sm,self`

**NAMD Performance  
(Apoa1)**



**NAMD Performance  
(f1atpase)**



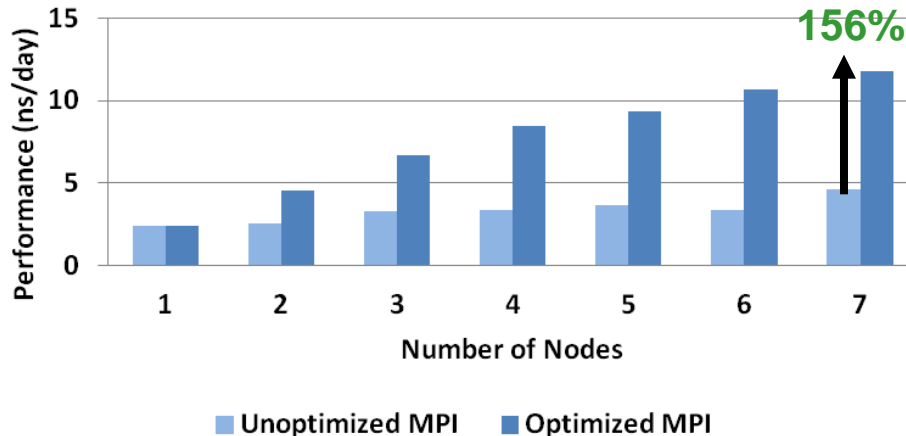
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**64 Cores/Node**

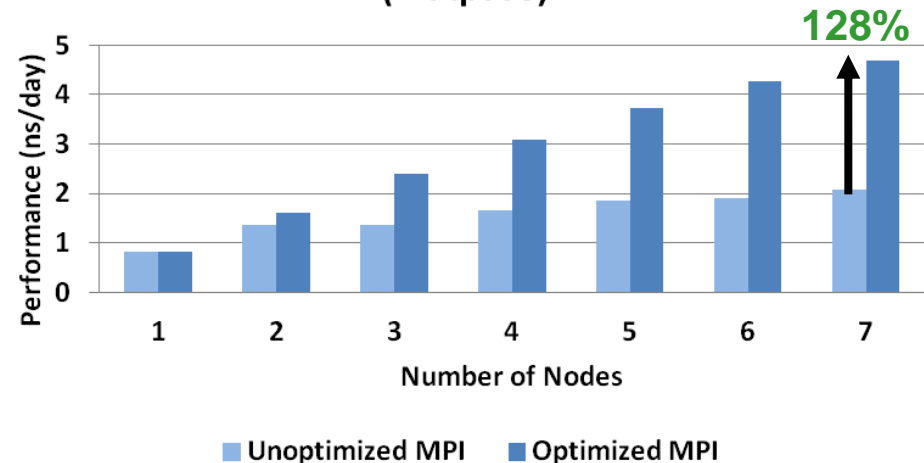
# NAMD Performance – Platform MPI Tuning

- **Using SRQ (Shared Receive Queue) allow Platform MPI to scale**
  - Reduces memory footprint and is essential for scaling
- **Explicit mapping of CPU cores to ensure MPI ranks are placed sequentially**
  - To ensure CPU core enumeration, check with “lstopo” from hwloc or “numactl --hardware”
- **Extra flags in optimized case for SRQ, RDMA params and core bindings:**
  - `-srq -e MPI_RDMA_MSGSIZE=32768,32768,4194304 -e MPI_RDMA_NSQRQRECV=2048 -e MPI_RDMA_NFRAGMENT=128 -cpu_bind=v,map_cpu:0,4,8,12,16,20,24,28,32,36,40,44,48,52,56,60,2,6,10,14,18,22,26,30,34,38,42,46,50,54,58,62,3,7,11,15,19,23,27,31,35,39,43,47,51,55,59,63,1,5,9,13,17,21,25,29,33,37,41,45,49,53,57,61`

### NAMD Performance (Apoa1)



### NAMD Performance (f1atpase)



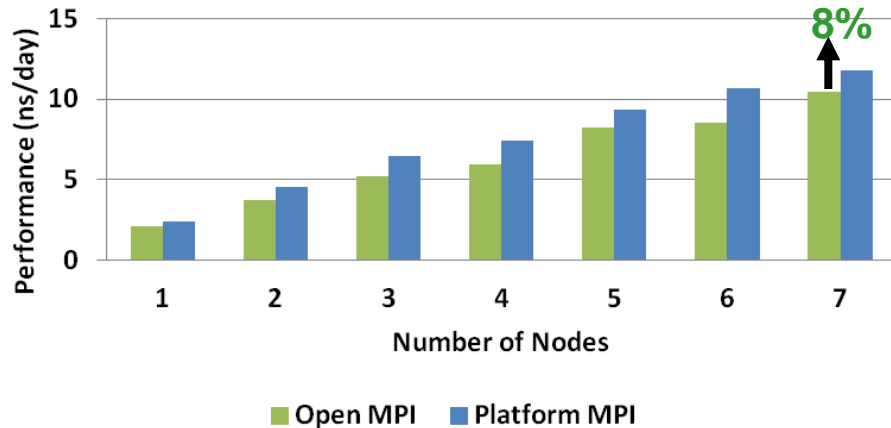
*Higher is better*

**64 Cores/Node**

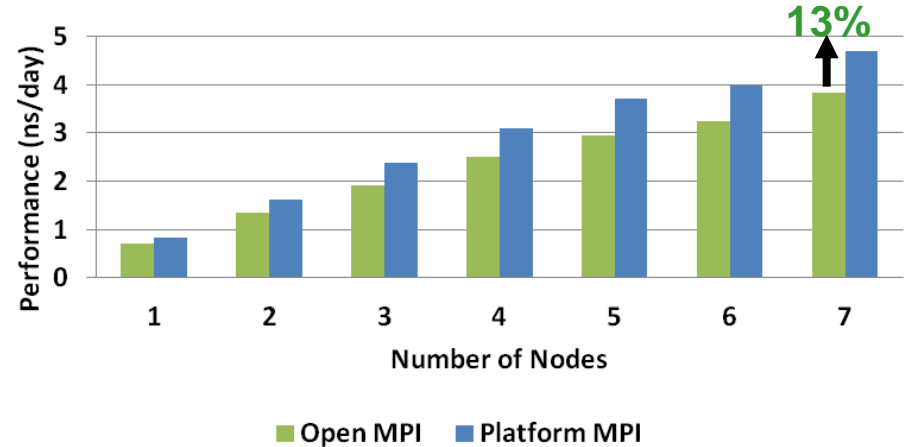
# NAMD Performance – MPI Implementations

- **The Platform MPI performs better than Open MPI**
  - Up to 8% better than Open MPI for apoa1 at 7-node
  - Up to 13% better than Open MPI for f1atpase at 7-node
- **The tuned flags are used for both Platform MPI and Open MPI**

**NAMD Performance  
(Apoa1)**



**NAMD Performance  
(f1atpase)**



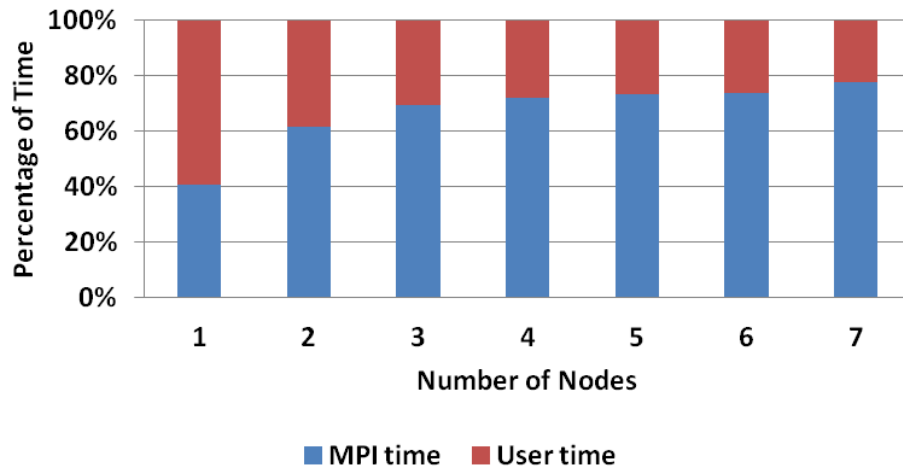
*Higher is better*

**64 Cores/Node**

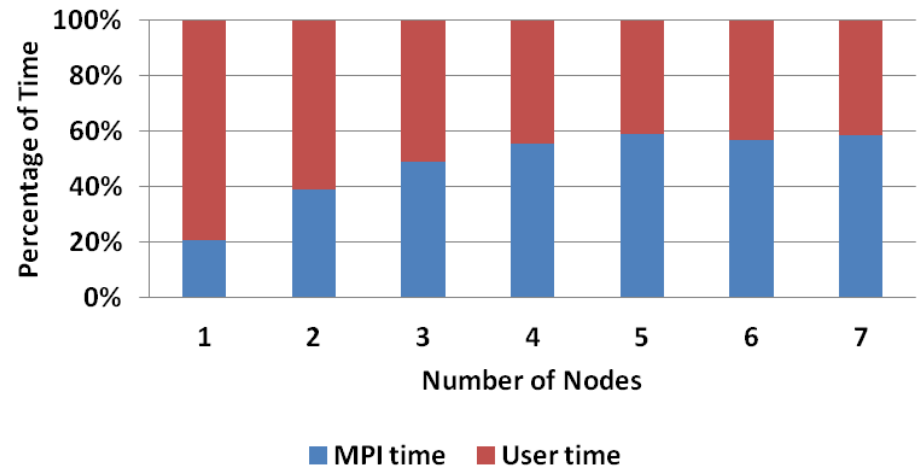
# NAMD Profiling – MPI/User Time Ratio

- **NAMD becomes highly communicative starting from 2-3 nodes**
  - Due to the high core counts per node
  - The stmv contains more CPU and MPI communication times than apoa1
- **MPI communication time dominates the overall time**
  - Shows low latency interconnect such as InfiniBand is required for good scalability

**NAMD Profiling  
(Apoa1)**  
MPI/User Time Ratio



**NAMD Profiling  
(stmv)**  
MPI/User Time Ratio



*Higher is better*

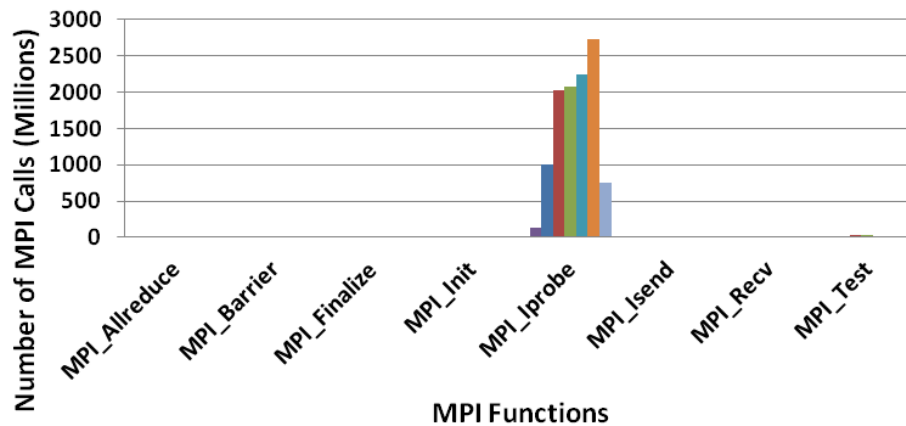
**64 Cores/Node**



# NAMD Profiling – Number of MPI Calls

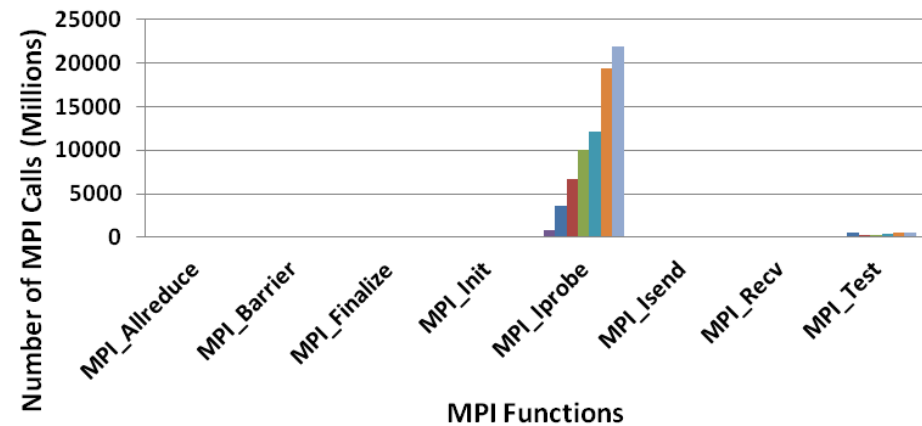
- **The most used MPI function is MPI\_Iprobe**
  - MPI\_Iprobe is used for testing non-blocking messages
  - Accounted for 97% of all MPI calls

**NAMD Profiling  
(Apoa1)  
Number of MPI Calls**



■ 1 Node ■ 2 Nodes ■ 3 Nodes ■ 4 Nodes  
■ 5 Nodes ■ 6 Nodes ■ 7 Nodes

**NAMD Profiling  
(stmv)  
Number of MPI Calls**

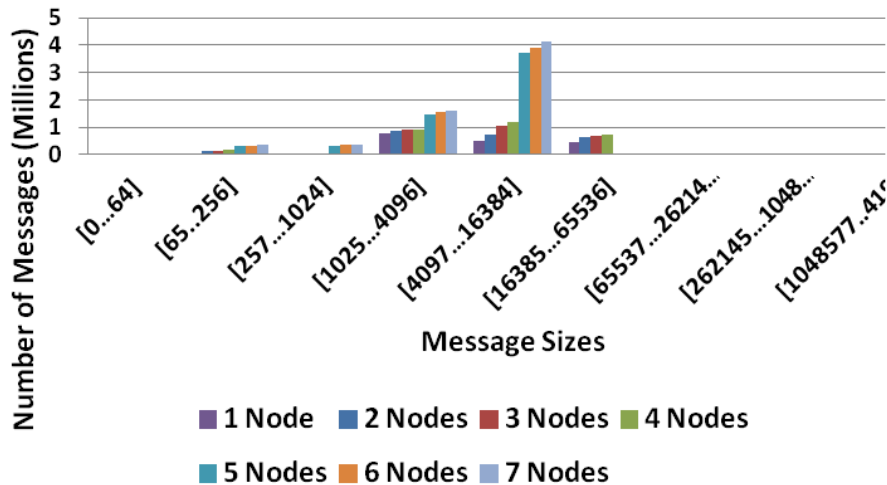


■ 1 Node ■ 2 Nodes ■ 3 Nodes ■ 4 Nodes  
■ 5 Nodes ■ 6 Nodes ■ 7 Nodes

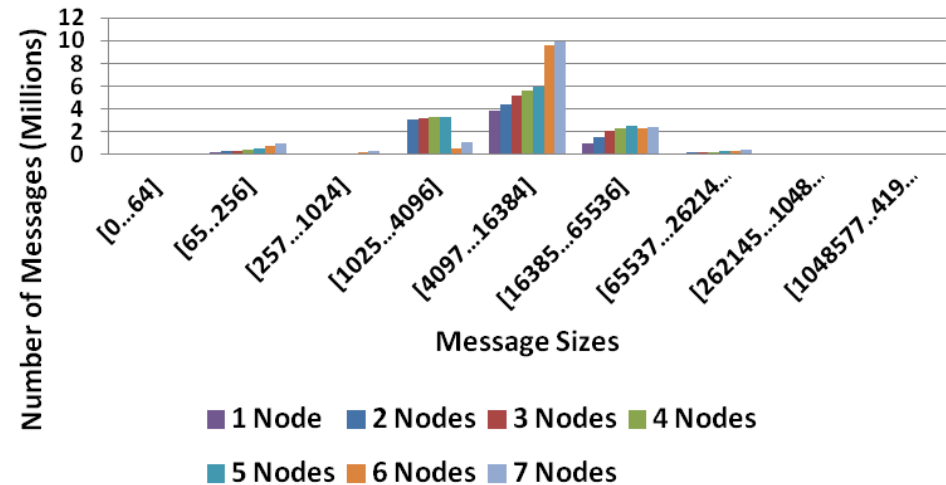
# NAMD Profiling – MPI Message Sizes

- Majority of the MPI message sizes are
  - in the range from 4KB to 16KB
- The increase in messages accelerates starting around 5-6 nodes

**NAMD Profiling**  
(Apoa1)  
MPI Message Sizes



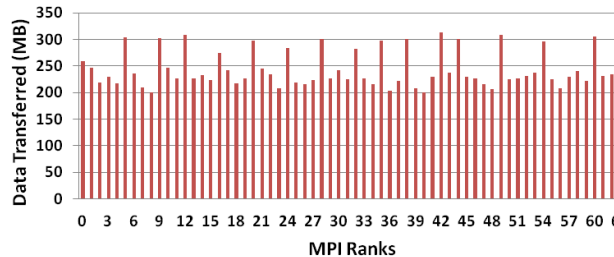
**NAMD Profiling**  
(stmv)  
MPI Message Sizes



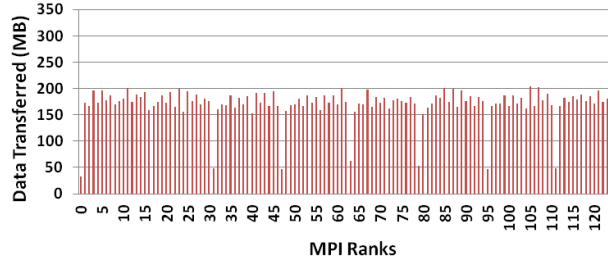
# NAMD Profiling – Data Transfer Per Process

- **Data transferred to each MPI rank is showing some variance**
  - But overall data transfer is roughly the same on a per-node basis
- **As the cluster scales, less data is driven to each rank and each node**
  - 300-600MB per rank in 1-node job versus 150-200MB per rank in a 4-node job

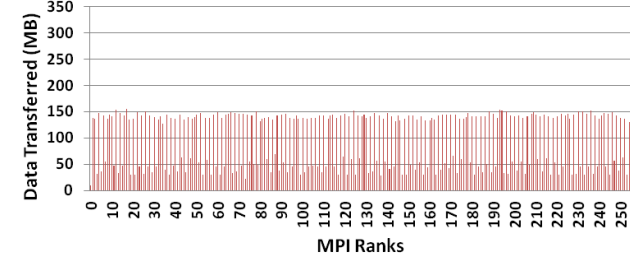
**NAMD Profiling  
(Apoa1, 1-node)  
Data Transferred by Ranks**



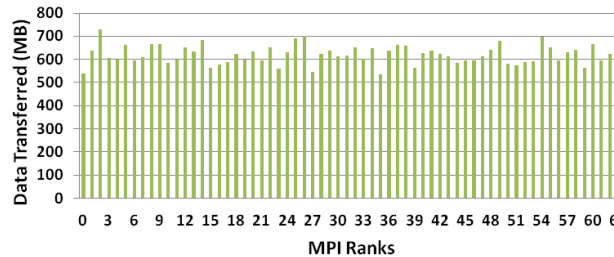
**NAMD Profiling  
(Apoa1, 2-node)  
Data Transferred by Ranks**



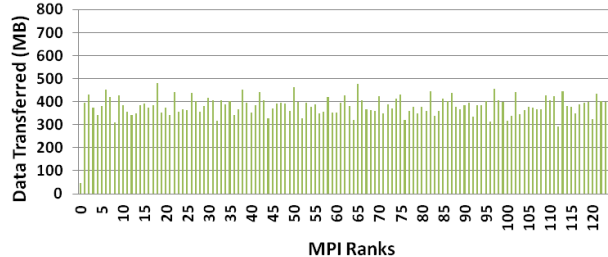
**NAMD Profiling  
(Apoa1, 4-node)  
Data Transferred by Ranks**



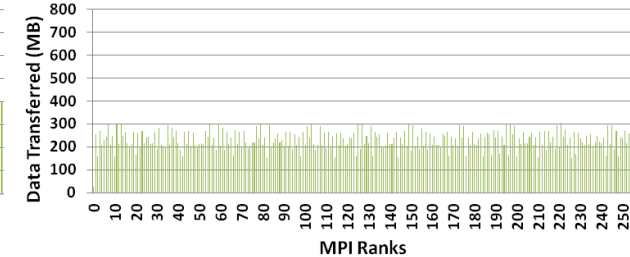
**NAMD Profiling  
(f1atpase, 1-node)  
Data Transferred by Ranks**



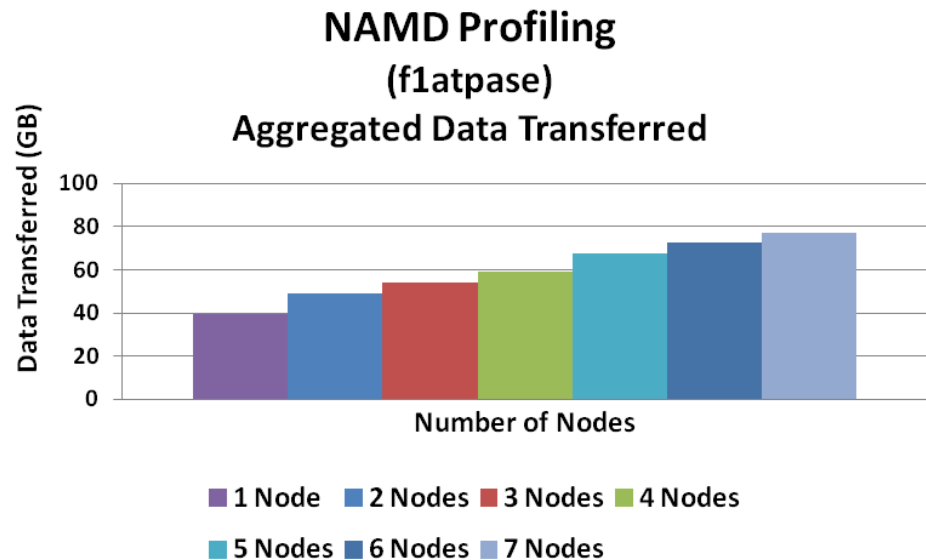
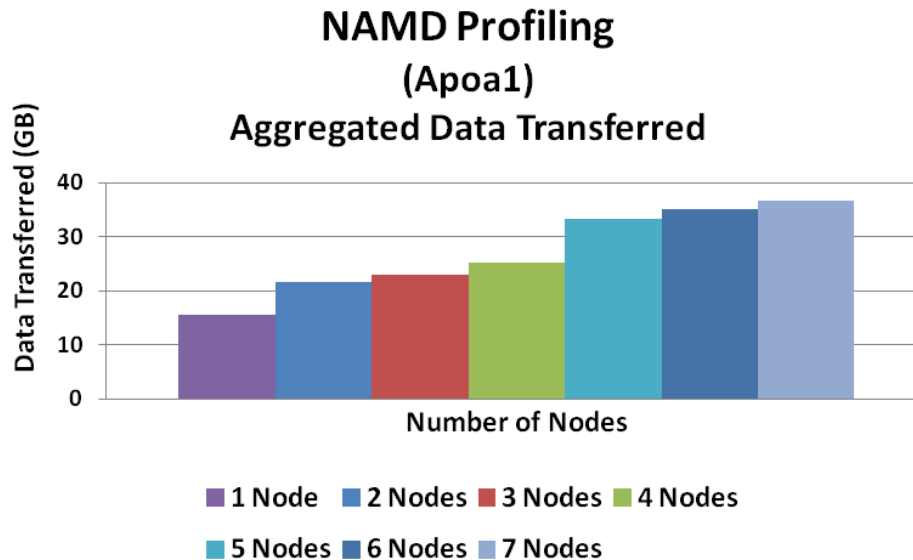
**NAMD Profiling  
(f1atpase, 2-node)  
Data Transferred by Ranks**



**NAMD Profiling  
(f1atpase, 4-node)  
Data Transferred by Ranks**



- **Aggregated data transfer refers to:**
  - Total amount of data being transferred in the network between all MPI ranks collectively
- **The total data transfer increases as the cluster scales**
- **Demonstrates the importance of scalable network interconnect**
  - InfiniBand can deliver bandwidth needed to push data in 40GB+ across the network



- **Interlagos provides higher performance than Magny-Cours CPUs**
  - Up to 34% performance gain with Open MPI
    - AMD Opteron™ 6276 (code name “Interlagos”) 16-core @ 2.3 GHz CPUs
    - AMD Opteron™ 6174 (code name “Magny-Cours”) 12-core @ 2.2GHz CPUs
- **Mellanox ConnectX®-3 proves significantly higher scalability for NAMD**
  - 4x to 5x higher performance versus 10GbE
- **Open MPI and Platform MPI benefit from tuned parameters**
  - Having XRC and SRQ enabled the MPIs to scale at large core counts
- **The tuned Platform MPI performs better than the tuned Open MPI**
  - By 8-13% on 2 different datasets



# Thank You

## HPC Advisory Council



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