

# Hints to improve automatic load balancing with LeWI for hybrid applications

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- Loss of efficiency
- Hybrid programming models ( $MPI + X$ )
- Manual tuning of parallel codes (load-balancing, data redistribution)

# The X (in this paper)

## OpenMP

- Directives to annotate parallel code
- Fork/join model with shared memory
- Number of threads may change *between* parallel regions

## SMPSs (SMPSuperscalar)

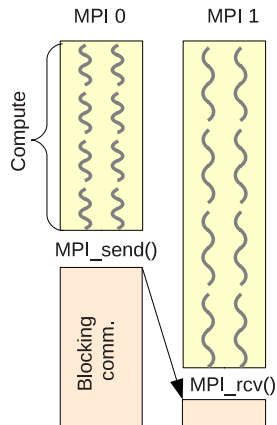
- Task as basic element
- Annotate *taskifiable* functions and their parameters (in/out/inout)
- Task graph to track dependencies
- Number of threads may change *any time*

## DLB (dynamic load balancing)

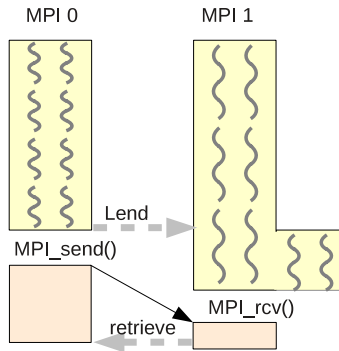
- “*Runtime interposition to [...] intercept MPI calls*”
- Balance load on the inner level (OpenMP/SMPs)
- Several load balancing algorithms

## LeWI (Lend CPU when Idle)

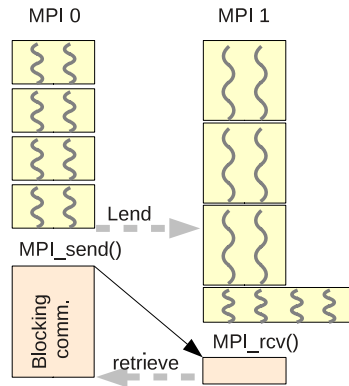
- CPUs of rank in blocking MPI call are idle
- Lend CPUs to other ranks and recover them after MPI call completes



(a) No load balancing.



(b) LeWI algorithm with SMPSSs.



(c) LeWI algorithm with OpenMP.

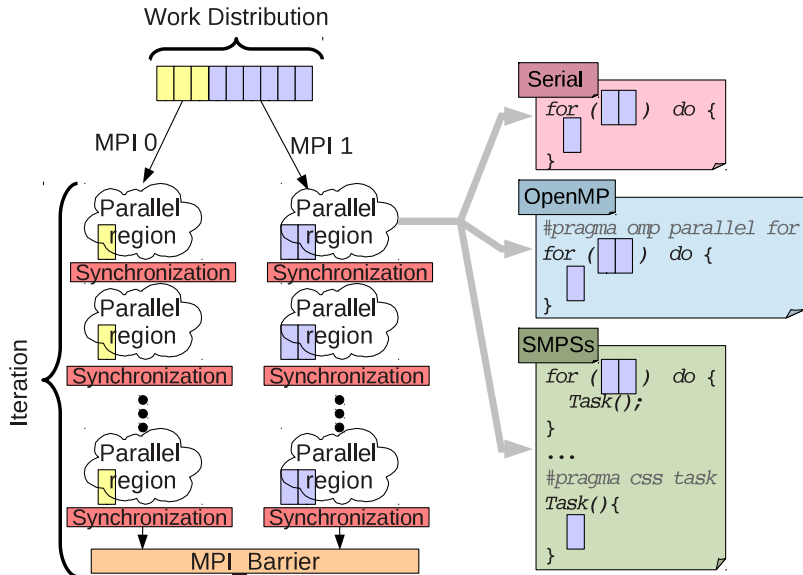
- “*Extensive performance evaluation*”
- “*Modeling parallelization characteristics that limit the automatic load balancing potential*”
- “*Improving automatic load balancing*”

- Marenstrum 2:  $2 \times$  IBM PowerPC 970MP (2 cores); 8 GiB RAM
- Linux 2.6.5-7.244-pseries64; MPICH; IBM XL C/C++ compiler w/o optimizations
- Metrics
  - $Speedup = \frac{parallel\_execution\_time}{serial\_execution\_time}$
  - $Efficiency = \frac{useful\_cpu\_time}{elapsed\_time * cpus}$  where  
 $useful\_cpu\_time = cpu\_time - (mpi\_time + openmp/smpss\_time + dlb\_time)$
  - $CPUs\_used$  to simultaneously run application code
- 3 benchmarks + 2 real applications

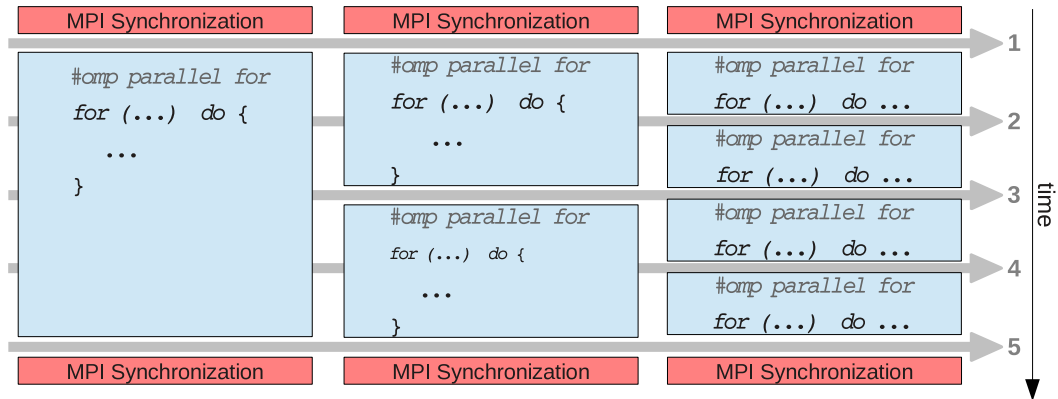
# PILS (Parallel ImbaLance Simulation)

- Synthetic benchmark
- Core: “*floating point operations without data involved*”
- Tunable parameters
  - Programming model (MPI, MPI + OpenMP, MPI + SMPs)
  - Load distribution
  - Parallelism grain ( $= \frac{1}{\#parallel\ regions}$ )
  - Iterations





# Parallelism Grain



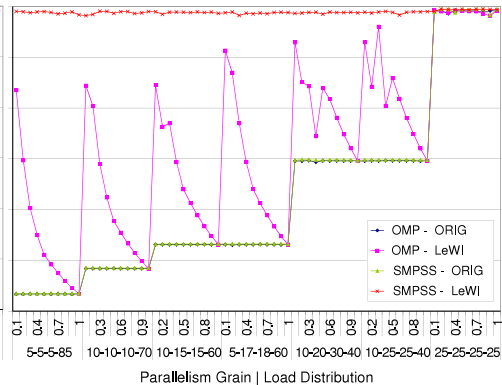
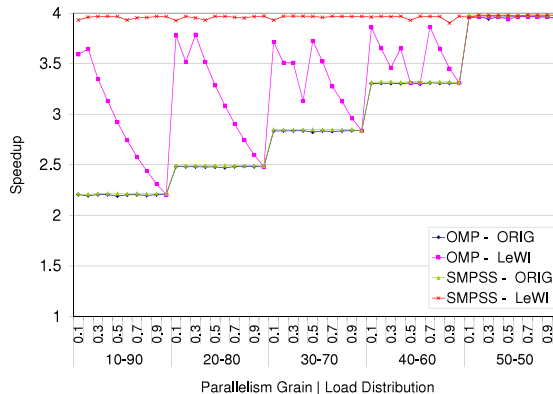
a) Parallelism Grain: 1. b) Parallelism Grain: 0.5. c) Parallelism Grain: 0.25.

- Benchmarks
  - BT-MZ: block tri-diagonal solver
  - LUB: LU matrix factorization
- Applications
  - Gromacs: molecular dynamics, MPI-only
  - Gadget: cosmological N-body/SPH (smoothed-particle hydrodynamics) simulation

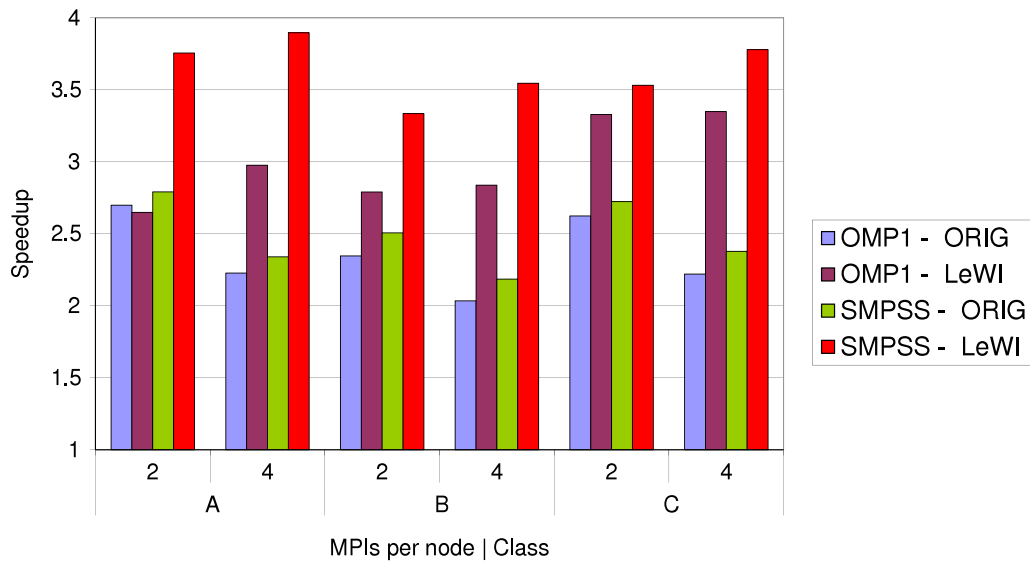
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| Application | Original version           | MPI + OpenMP | MPI + SMPs | Executed in nodes (cpus) |
|-------------|----------------------------|--------------|------------|--------------------------|
| PILS        | MPI + OpenMP<br>MPI + SMPs | X            | X          | 1 (4)                    |
| BT-MZ       | MPI + OpenMP               | X            | X          | 1, 2, 4 (4, 8, 16)       |
| LUB         | MPI + OpenMP<br>MPI + SMPs | X            | X          | 1, 2, 4 (4, 8, 16)       |
| Gromacs     | MPI                        |              | X          | 1.64 (4.256)             |
| Gadget      | MPI                        | X            |            | 200 (800)                |

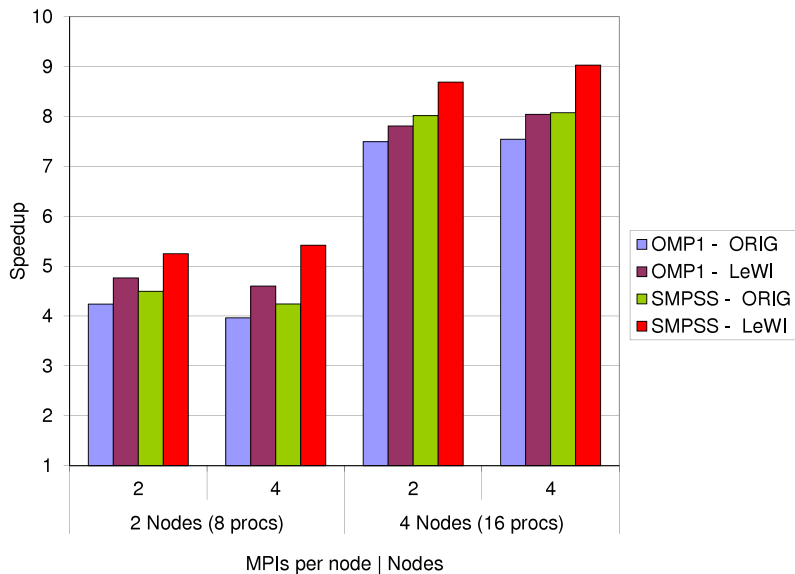
# PILS, 2 and 4 MPI processes



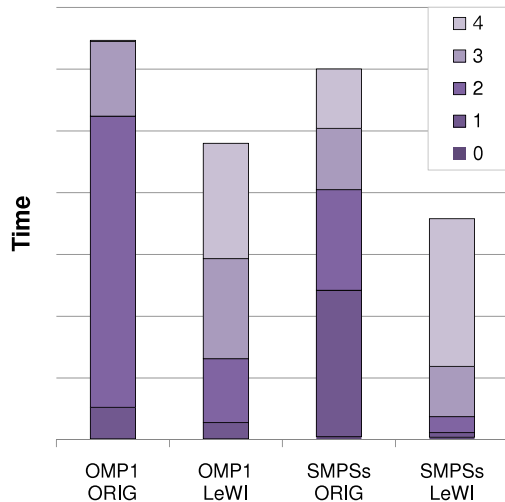
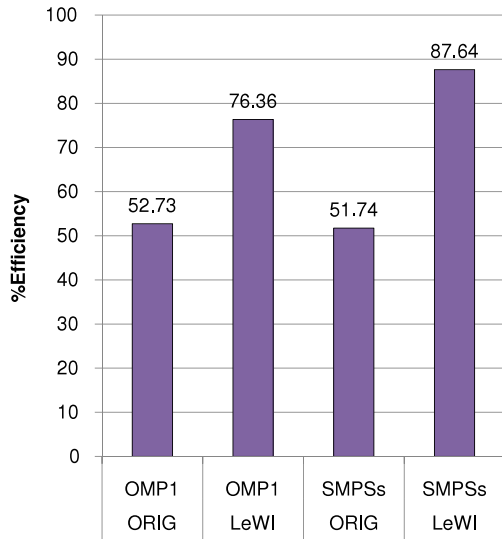
## BT-MZ; 1 node



## BT-MZ; 2,4 nodes; Class C

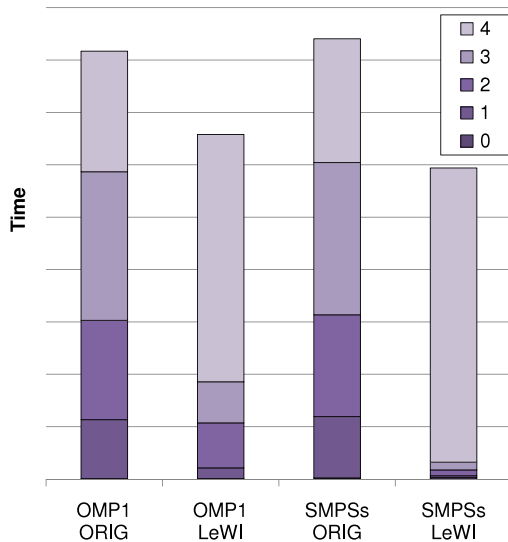
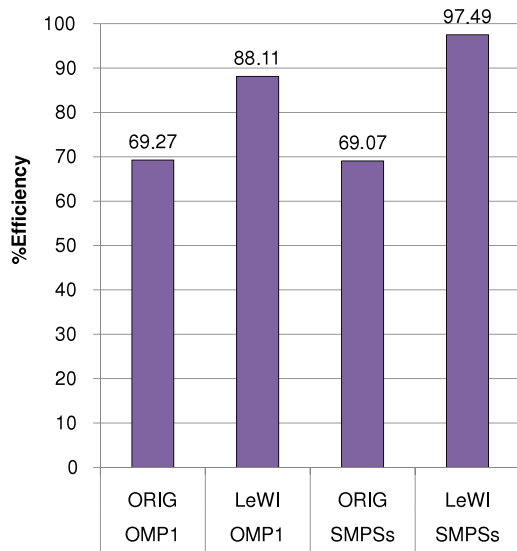


# BT-MZ; 1 node; 4 MPI processes

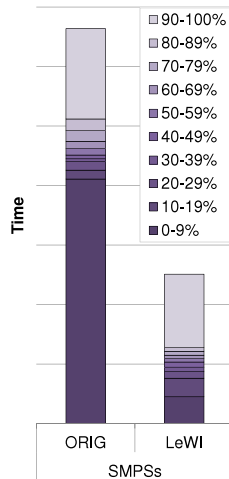
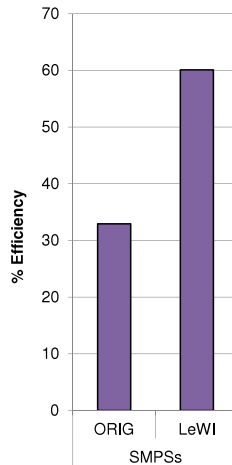
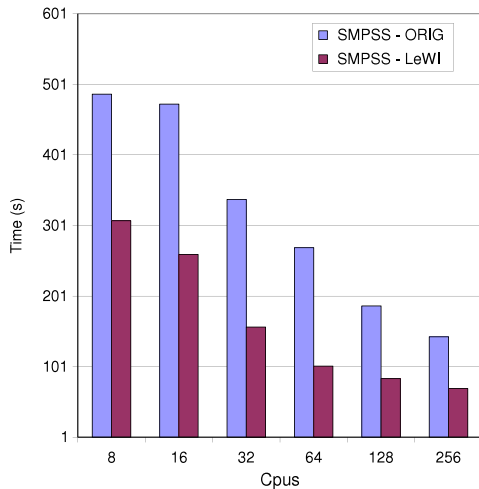




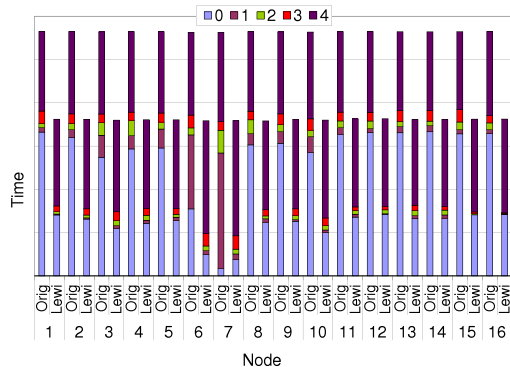
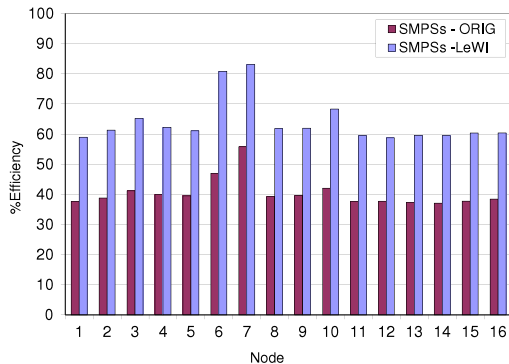
# LUB; 1 node; Block size 200



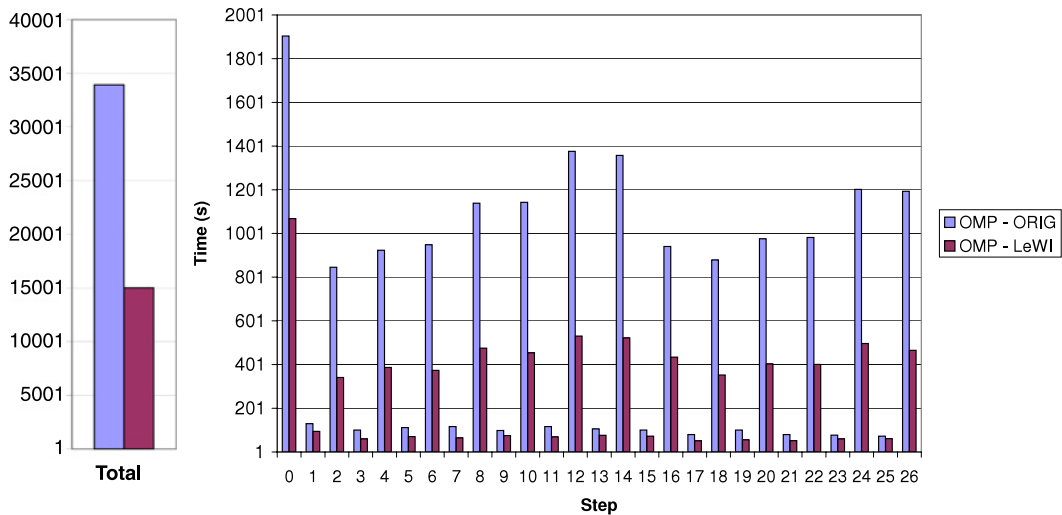
# Gromacs; 1-64 nodes + Details for 16 nodes



# Gromacs; Efficiency + CPUs used per Node

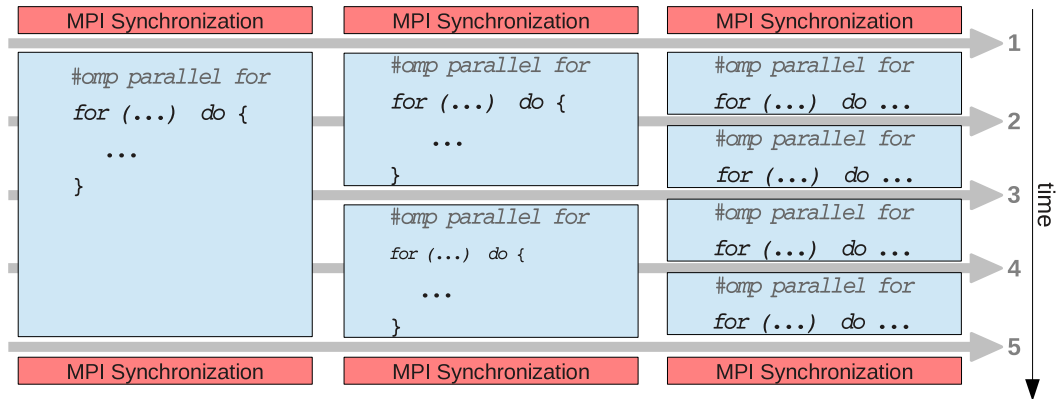


# Gadget; 200 nodes



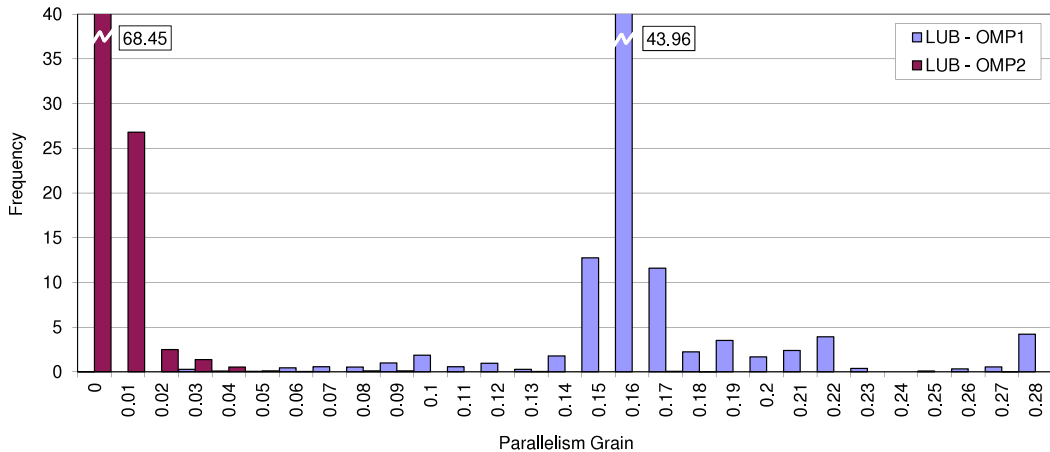
- “*Parallelism Grain in OpenMP applications*”
- “*Task duration in SMPs applications*”
- “*Distribution of MPI processes among computation nodes*”

# Parallelism Grain

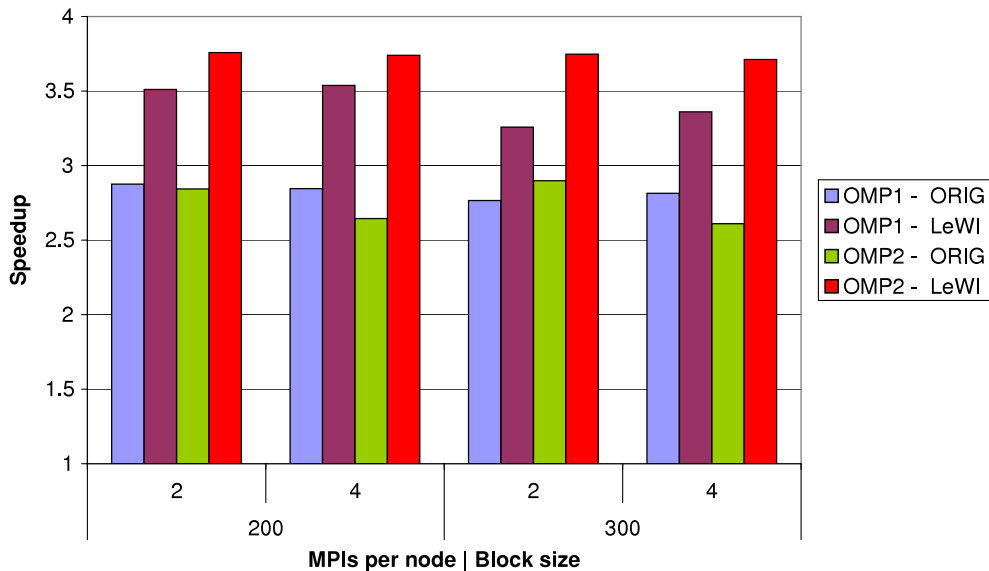


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# Modified Parallelism Grain in LUB

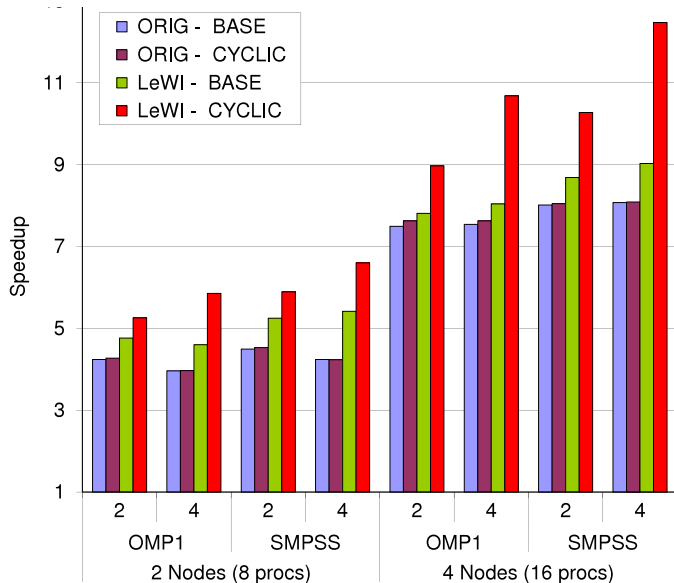


# Performance of Modified LUB

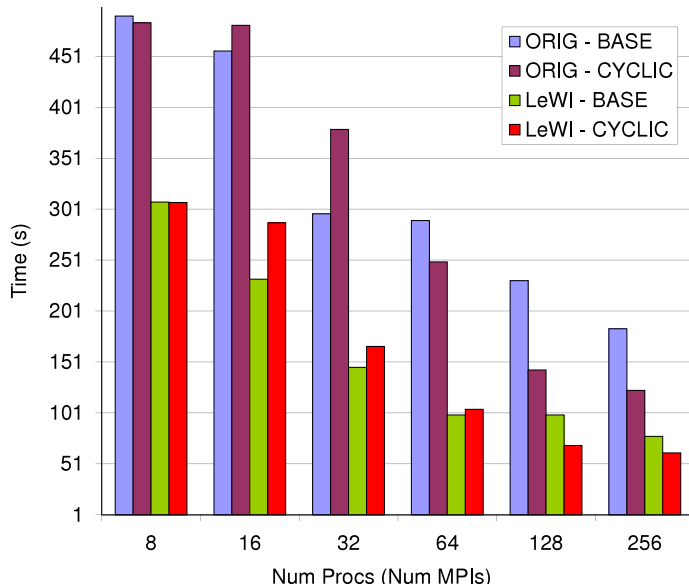




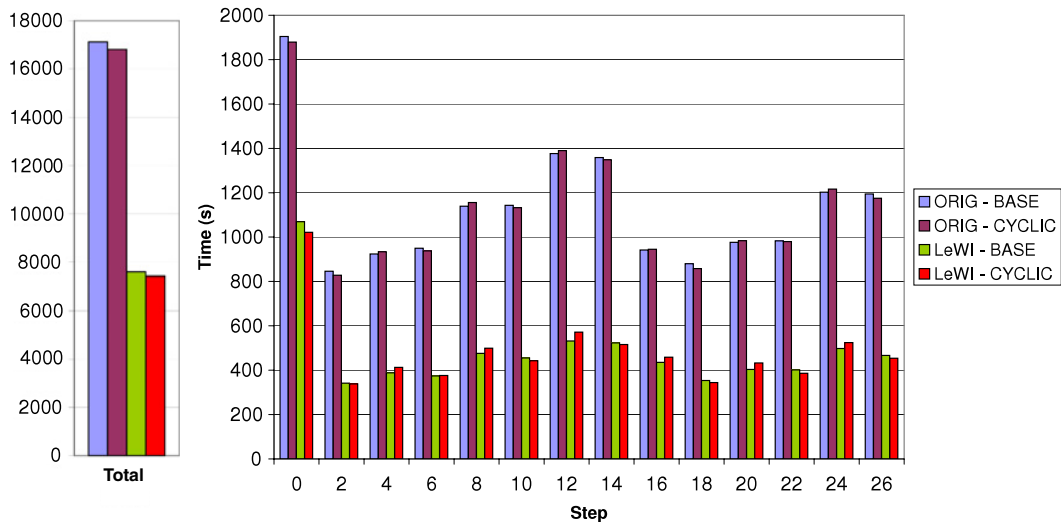
# Rank Distribution — BT-MZ



# Rank Distribution — Gromacs



# Rank Distribution — Gadget



## Summary

- DLB/LeWI can improve performance transparently
- Inter-node load imbalances not handled
- Granularity of parallelism and placement as important factors
- Optimal configuration with vs. without DLB/LeWI

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## Discussion

- Interaction with MPI
- Benchmarks (1.5 of 3 NPB-MZ, arbitrary load distribution)
- How to find “the right” granularity