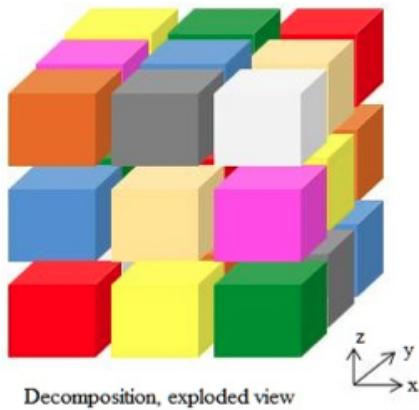

NPB Application Kernels

A 1D partitioning distributed approach

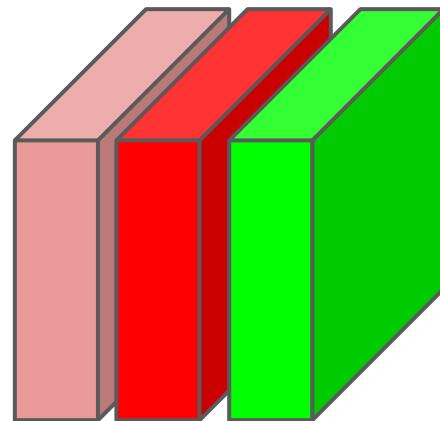
Brook Kassa, Prerna Narayanpure, Verónica Vergara L.

1D-MPI-NPB

Original distributed NPB
uses multidimensional
data partitioning



The original shared
memory NPB uses a 1D
data partition



Goal: Implement a distributed version of NPB
that uses a 1D data partitioning scheme

NPB Application Kernels

The NPB application kernels simulate processes on real CFD applications. The three solvers are:

- **Block Tridiagonal (BT)**: solves multiple, independent systems of non-diagonally dominant, block tridiagonal equations with 5 by 5 block size
- **Lower Upper diagonal (LU)**: employs a symmetric successive over-relaxation (SSOR) numerical scheme to solve a regular, sparse, 5 by 5 lower and upper triangular systems.
- **Scalar Pentadiagonal (SP)**: solves three sets of uncoupled scalar pentadiagonal systems of equations in x, y and z direction.

NPB Application Kernels

- *Number of lines of code*

App. Kernel	Main Routine	Total
<i>BT</i>	328	9162
<i>SP</i>	245	4902
<i>LU</i>	199	5957

- *Problem Size*
 - 1) Class 'S' - 12 x 12 x 12
 - 2) Class 'A' - 64 x 64 x 64
 - 3) Class 'C' - 162 x 162 x 162

Translation Steps: for loop

```
#pragma omp for
for(k=0; k< Nz; k++)
    for(j=0; j< Ny; j++)
        for(i=0; i< Nx; i++)
            rhs[k][j][i] =
...
...
```



```
size = # proc
rank = my proc ID
kgap = Nz/size
kstart = rank * kgap
kend = kstart + kgap

for(k=kstart; k<kend; k++)
    for(j=0; j< Ny; j++)
        for(i=0; i< Nx; i++)
            rhs[k][j][i] = ...
```

Translation Steps

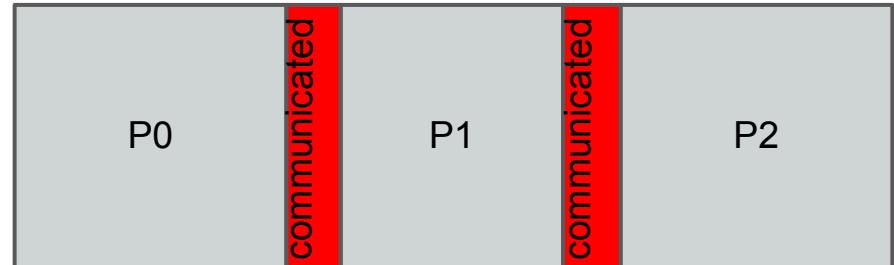
```
for (k=0; k < Nz; k++)  
  for (i=0; i < Nx; i++)  
    for (j=0; j < Ny; j++)  
      rsd[k][i][j] = ...
```

```
...
```

```
MPI_Allgather (rsd[kstart][0][0], rsd[0][0][0])
```



```
if(my_id == 0)  
  MPI_Send(&rsd[kend - 2][0][0], ..., rank+1, ...)  
else  
  MPI_Recv(&rsd[kstart - 2][0][0], ..., rank-1, ...)  
  
... = rsd[k-2][i][j] + ...
```



Translation Steps : Other transformation

- Reduction operation

```
#pragma omp atomic  
sum = sum + x
```



```
MPI_Allreduce(gsum,  
sum)
```

- Master construct

```
#pragma omp master  
time_start(trhs)
```



```
if(rank == root)  
time_start(trhs)  
MPI_Barrier...
```

- Flush construct

```
#pragma omp flush  
(isynch)
```



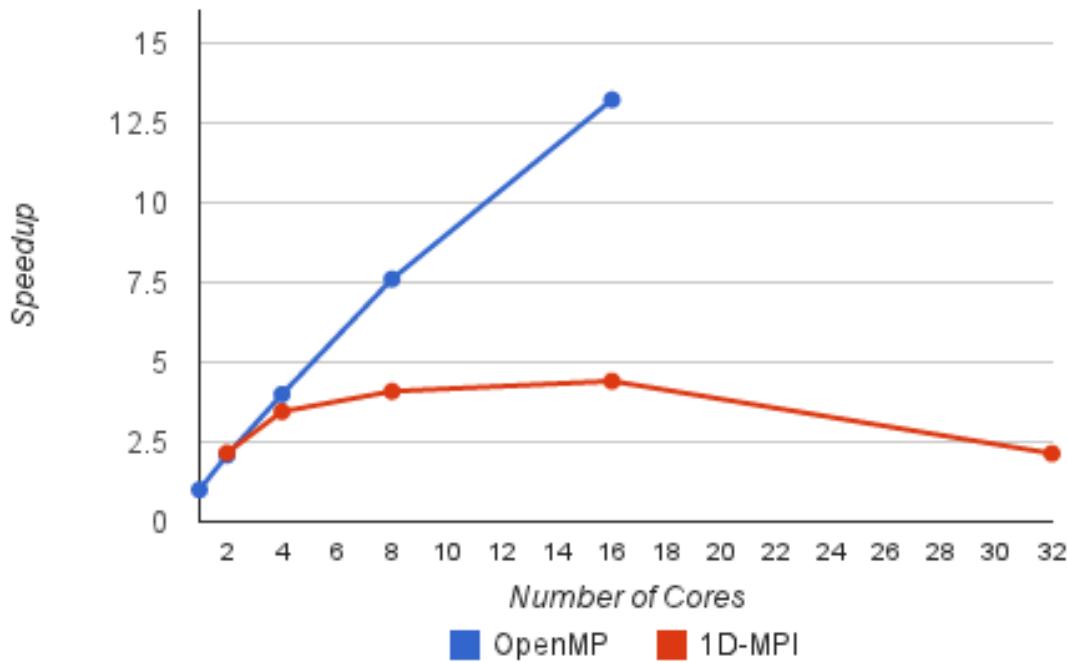
```
MPI_Bcast(isynch, root)
```

Results : BT Benchmark

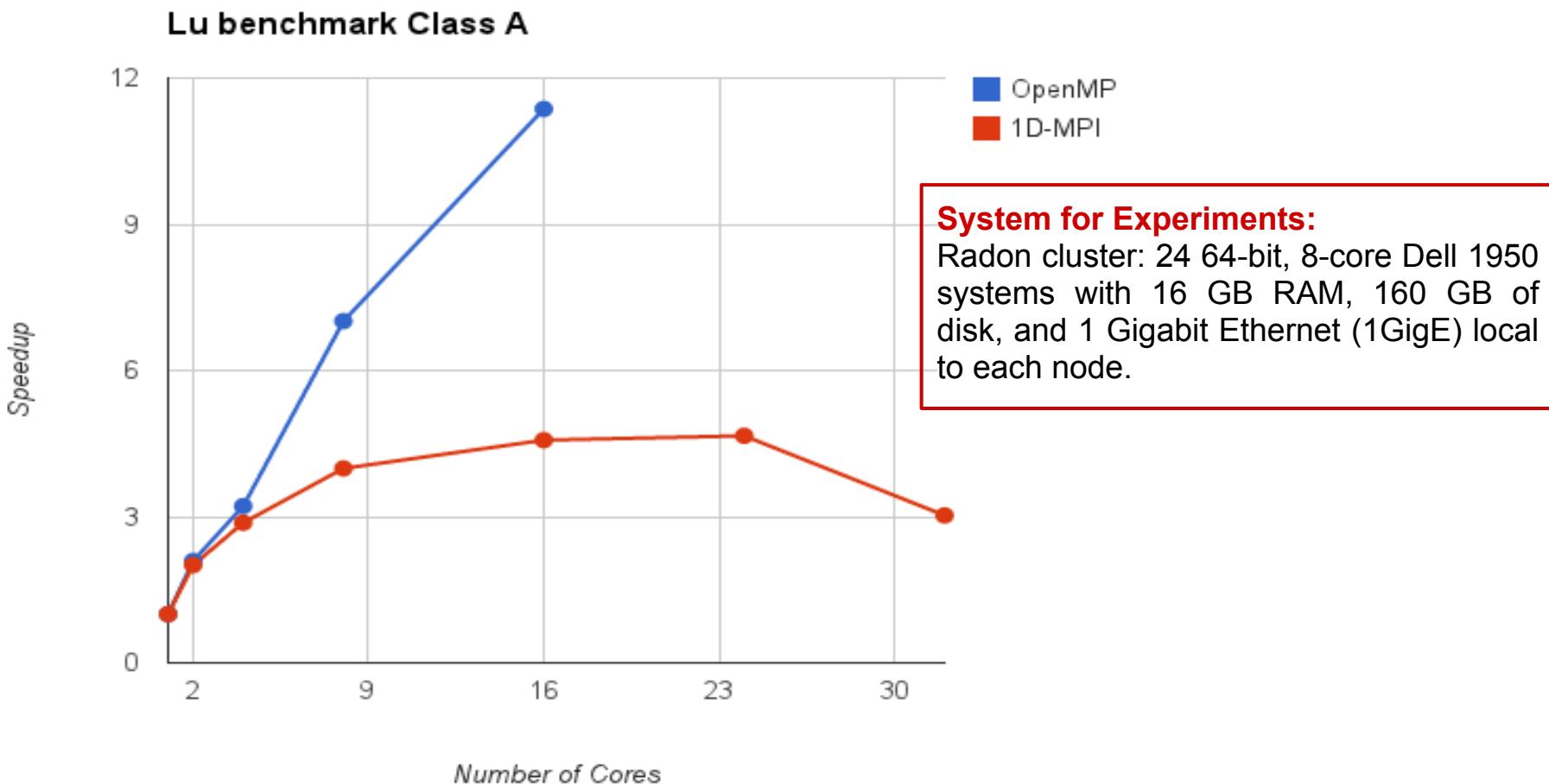
System for Experiments:

Carter cluster with two 8-core Intel Xeon-E5 nodes with 32 GB of memory and 56 Gbps FDR Infiniband interconnects.

BT benchmark Class A



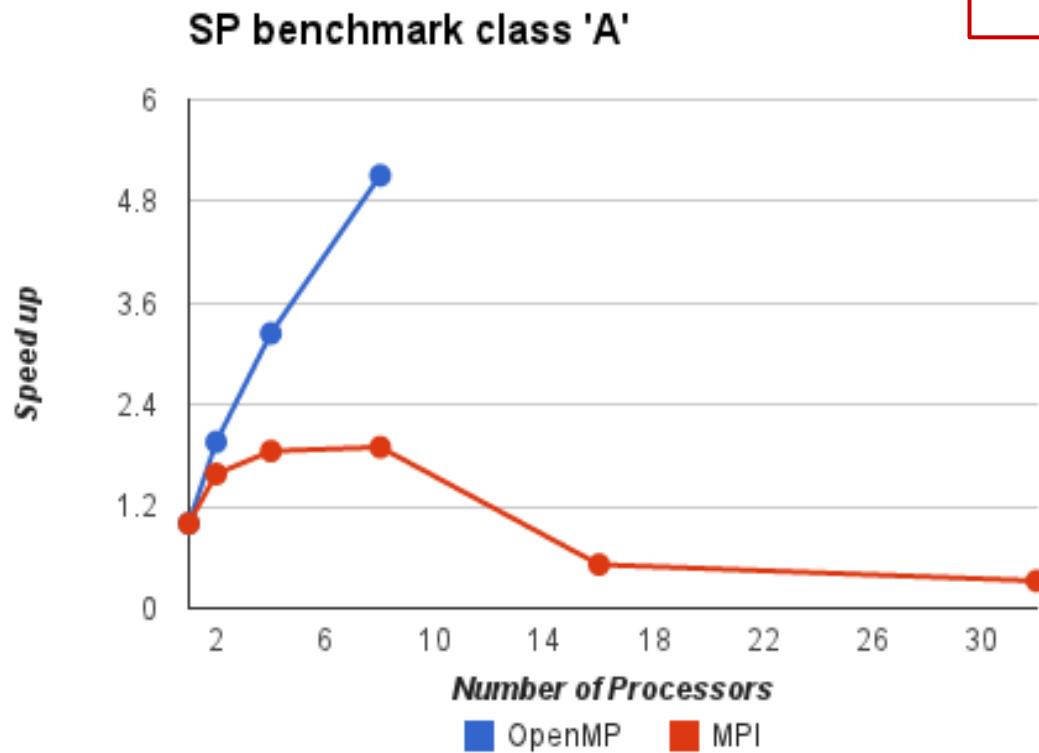
Results : LU Benchmark



Results : SP Benchmark

System for Experiment

ACSL cluster with 10 nodes each having 2 quad core AMD opteron, 16GB RAM and GigE network.

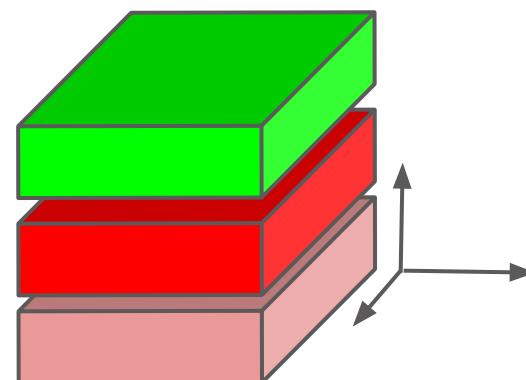
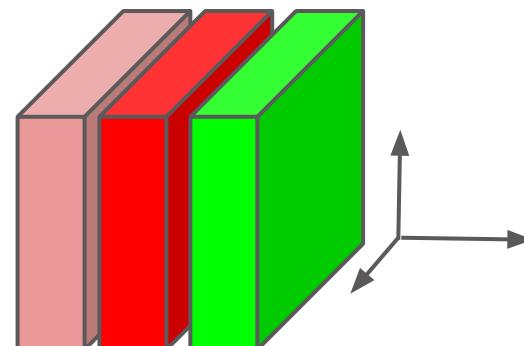


Challenges

- Overhead for MPI communication specially with collective communication.
- OpenMP parallelization across different dimensions

```
#pragma omp for
for(k=0; k< Nz; k++)
    for(j=0; j< Ny; j++)
        for(i=0; i< Nx; i++)
            rhs[k][j][i] = ...
```

for(j=**jstart**; j<**jend**; j++)
 for(k=0; k< Nz; k++)
 for(i=0; i< Nx; i++)
 rhs[k][j][i] = ...



Because of some true dependencies some matrix had to be transposed and partitioned on second dimension of the array.

Future Work

- Increase problem size to class C
- Increase computation and communication overlap by interleaving non-blocking point-to-point communications
- Implement algorithmic changes that will allow us to exploit further parallelism

Questions?

Results : BT Benchmark (with MPI)

