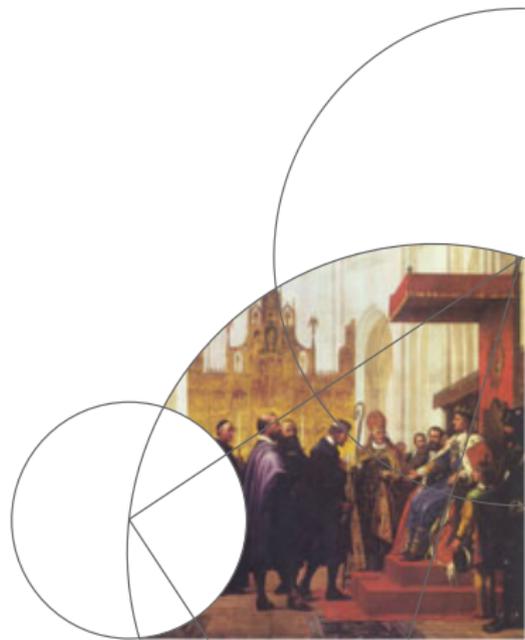




Portable Parallelization with the Bohrium Runtime System

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Design Philosophy of Bohrium

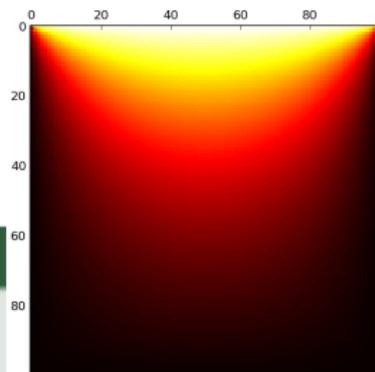
- Do science, not programming
- No bugs
 - Correctness above all
- Future proof
 - Don't optimize for today's hardware



Python / NumPy

Heat Equation – a 5-point stencil

```
import numpy
def heat2d(G, epsilon):
    delta = epsilon+1
    while delta > epsilon:
        tmp = 0.2*(G[1:-1,1:-1]+G[-2:,1:-1]+\
                  G[2:,1:-1]+G[1:-1,:2]+G[1:-1,2:])
        delta = numpy.sum(numpy.abs(tmp-center))
        center[:] = tmp
```



Python

```

# Parallelize a Python function
import multiprocessing
def parallelize(f):
    """Parallelize a Python function"""
    def wrapper(*args, **kwargs):
        pool = multiprocessing.Pool()
        return pool.map(f, args, kwargs)
    return wrapper

```

OpenMP

```

// Parallelize a C function
#include <omp.h>
void parallelize(f) {
    #pragma omp parallel
    {
        f();
    }
}

```

OpenMP+MPI

```

// Parallelize a C function with MPI
#include <mpi.h>
#include <omp.h>
int main(int argc, char** argv) {
    MPI_Init(&argc, &argv);
    MPI_Comm_rank(MPI_COMM_WORLD, &rank);
    MPI_Comm_size(MPI_COMM_WORLD, &nproc);
    #pragma omp parallel
    {
        // ...
    }
    MPI_Finalize();
}

```

+Latency Hiding

```

// Parallelize a C function with MPI and OpenMP
#include <mpi.h>
#include <omp.h>
int main(int argc, char** argv) {
    MPI_Init(&argc, &argv);
    MPI_Comm_rank(MPI_COMM_WORLD, &rank);
    MPI_Comm_size(MPI_COMM_WORLD, &nproc);
    #pragma omp parallel
    {
        // ...
    }
    MPI_Finalize();
}

```

OpenCL

```

// Parallelize a C function with OpenCL
#include <CL/cl.h>
int main(int argc, char** argv) {
    // ...
}

```





Python / NumPy

```
import numpy
def heat2d(G, epsilon):
    delta = epsilon+1
    while delta > epsilon:
        tmp = 0.2*(G[1:-1,1:-1]+G[-2,:1:-1]+\
            G[2,:1:-1]+G[1:-1,:2]+G[1:-1,2:])
        delta = numpy.sum(numpy.abs(tmp-center))
        center[:] = tmp
```

python -m bohrium heat2d.py



Bytecode

```
ADD t1, center, north
ADD t2, t1, south
FREE t1
ADD t3, t2, east
FREE t2
ADD t4, t3, west
FREE t3
MUL tmp, t4, 0.2
FREE t4
MINUS t5, tmp, center
ABS t6, t5
FREE t5
ADD_REDUCE t7, t6
FREE t6
ADD_REDUCE delta, t7
FREE t7
COPY center, tmp
FREE tmp
SYNC delta
```



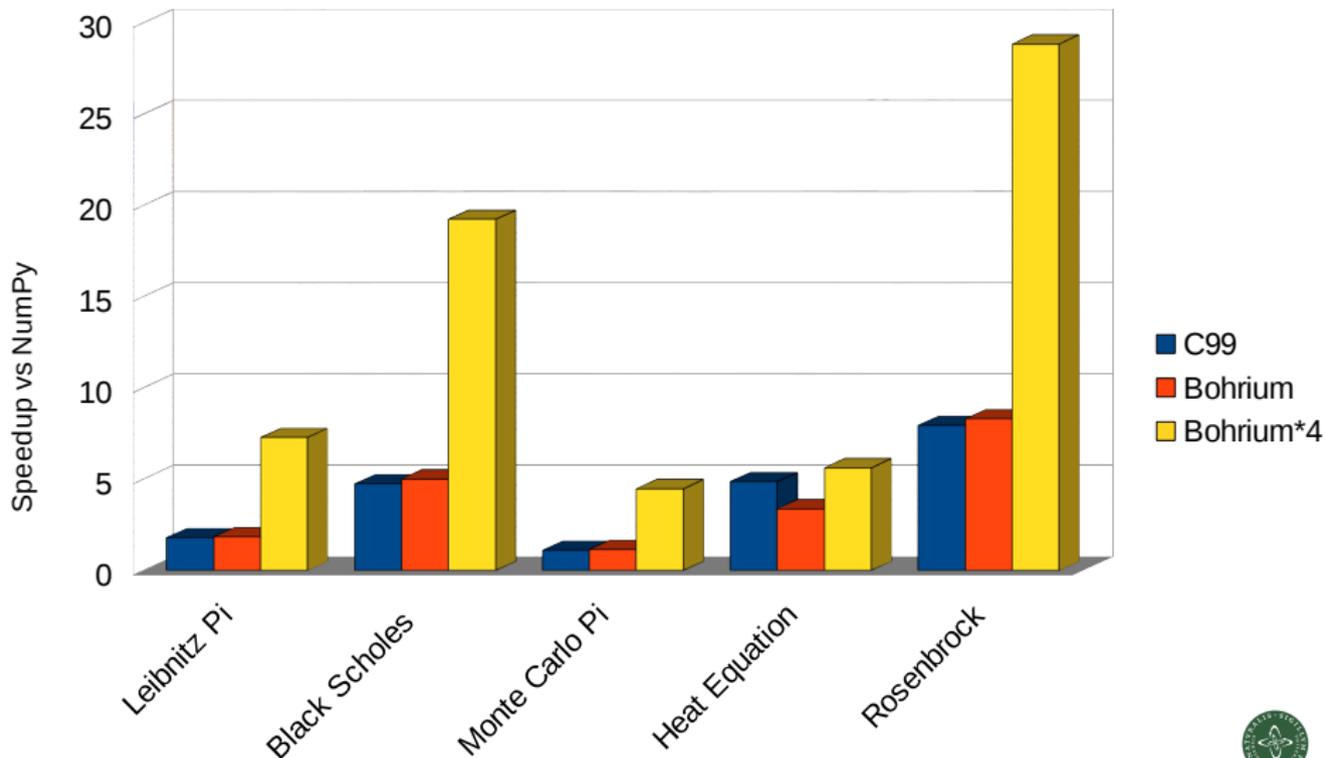
OpenMP



OpenCL



Performance – Speedup VS NumPy



Freely Available and Open Source



`www.bh107.org`

`www.github.com/bh107/bohrium`



Performance – Bohrium vs C

	C99	Bohrium		NumPy
Threads	1	1	4	1
Leibnitz Pi	1	0.99	0.25	1.79
Black Scholes	1	0.95	0.25	4.76
Monte Carlo Pi	1	0.98	0.25	1.11
Heat Equation	1	1.45	0.87	5.00
Rosenbrock	1	0.96	0.27	7.69

Table: Normalized runtime of Bohrium using C99 as baseline (**lower is better**). The timings are for full runs.



ANSI C – Sequential

Sequential

```
#include <math.h>
void heat2d(int size, double *grid, double epsilon)
{
    int gsize = size+2; //Size + borders.
    double *T = malloc(gsize*gsize*sizeof(double));
    double delta = epsilon+1;
    while(delta > epsilon)
    {
        double *a = grid;
        double *t = T;
        delta = 0;
        for(i=0; i<size; ++i)
        {
            double *up      = a+1;
            double *left   = a+gsize;
            double *right  = a+gsize+2;
            double *down   = a+1+gsize+2;
            double *center = a+gsize+1;
            double *t_center = t+gsize+1;
            for(j=0; j<size; ++j)
            {
                *t_center = (*center + *up++ + *left++ + *right++ + *down++) / 5.0;
                delta += fabs(t_center+center);
            }
            a += gsize;
            t += gsize;
        }
        memcpy(A, T, gsize*gsize*sizeof(double));
    }
}
```



ANSI C – OpenMP

Multi-core

```
#include <math.h>
void heat2d(int size, double *grid, double epsilon)
{
    int gsize = size+2; //Size + borders.
    double *T = malloc(gsize*gsize*sizeof(double));
    double delta = epsilon+1;
    while(delta > epsilon)
    {
        delta = 0;
        #pragma omp parallel for shared(grid,T) reduction(+:delta)
        for(i=0; i<size; ++i)
        {
            int a = i * gsize;
            double *up      = &grid[a+1];
            double *left   = &grid[a+gsize];
            double *right  = &grid[a+gsize+2];
            double *down   = &grid[a+1+gsize*2];
            double *center = &grid[a+gsize+1];
            double *t_center = &T[a+gsize+1];
            for(j=0; j<size; ++j)
            {
                *t_center = (*center + *up++ + *left++ + *right++ + *down++) / 5.0;
                delta += fabs(t_center-center);
            }
        }
        memcpy(grid, T, gsize*gsize*sizeof(double));
    }
}
```



ANSI C – OpenMP + MPI

Multi-core, Multiprocessor

```

#include <math.h>
#include <mpi.h>
void heat2d(int size, double *grid, double epsilon)
{
    int gsize = SIZE+2; //Size + borders.
    MPI_Comm_rank(MPI_COMM_WORLD, &myrank);
    MPI_Comm_size(MPI_COMM_WORLD, &worldsize);
    MPI_Comm comm;
    int periods[] = {0};
    MPI_Cart_create(MPI_COMM_WORLD, 1, &worldsize,
                   periods, 1, &comm);
    int l_size = SIZE / worldsize;
    if(myrank == worldsize-1)
        l_size += SIZE % worldsize;
    int l_gsize = l_size + 2; //Size + borders.

    double delta = epsilon+1;
    while(delta > epsilon)
    {
        int p_src, p_dest;
        //Send/receive - neighbor above
        MPI_Cart_shift(comm,0,1,&p_src,&p_dest);
        MPI_Sendrecv(grid+gsize, gsize, MPI_DOUBLE,
                    p_dest, 1, grid, gsize, MPI_DOUBLE,
                    p_src, 1, comm, MPI_STATUS_IGNORE);
        //Send/receive - neighbor below
        MPI_Cart_shift(comm,0,-1,&p_src,&p_dest);
        MPI_Sendrecv(grid+(l_gsize-2)*gsize,
                    gsize, MPI_DOUBLE,
                    p_dest, 1, grid+(l_gsize-1)*gsize,
                    gsize, MPI_DOUBLE,
                    p_src, 1, comm, MPI_STATUS_IGNORE);

        delta = 0;
        #pragma omp parallel for shared(grid,T) reduction(+:delta)
        for(i=0; i<size; ++i)
        {
            int a = i * gsize;
            double *up = &grid[a+1];
            double *left = &grid[a+gsize];
            double *right = &grid[a+gsize+2];
            double *down = &grid[a+1+gsize+2];
            double *center = &grid[a+gsize+1];
            double *t_center = &T[a+gsize+1];
            for(j=0; j<size; ++j)
            {
                *t_center = (*center + *up++ + *left++ + *right++ + *down++) / 5.0;
                delta += fabs(t_center+center);
            }
        }
        memcpy(grid, T, gsize*gsize*sizeof(double));
        MPI_Allreduce(delta, delta, 1, MPI_DOUBLE, MPI_SUM, MPI_COMM_WORLD);
    }
}

```



ANSI C – OpenMP + MPI + Latency Hiding

Double Buffering

```

#include <math.h>
#include <mpi.h>
void heat2d(int size, double *grid, double *epsilon)
{
    int qsize = size*2; //size * borders.
    MPI_Comm_rank(MPI_COMM_WORLD, &rank);
    MPI_Comm_size(MPI_COMM_WORLD, &worldsize);
    MPI_Comm_comm();
    int periods[] = {2};
    MPI_Cart_create(MPI_COMM_WORLD, 1, &worldsize,
        periods, 1, &comm);
    int l_size = size / worldsize;
    #pragma omp warrsize-1
    l_size = size % worldsize;
    int l_qsize = l_size + 2; //size + borders.

    double delta = *epsilon;
    while(delta > *epsilon)
    {
        int p_err, p_dest;
        MPI_Request reqq[4];
        //Initiate send/receive - neighbor above
        MPI_Cart_shift(comm, 0, 1, &p_err, &p_dest);
        MPI_Isend(grid+qsize, qsize, MPI_DOUBLE, p_dest,
            1, comm, reqq[0]);
        MPI_Irecv(grid, qsize, MPI_DOUBLE, p_err,
            1, comm, reqq[1]);
        //Initiate send/receive - neighbor below
        MPI_Cart_shift(comm, 0, -1, &p_err, &p_dest);
        MPI_Isend(grid-l_qsize-2*qsize, qsize,
            MPI_DOUBLE,
            p_dest, 1, comm, reqq[2]);
        MPI_Irecv(grid-l_qsize-1*qsize, qsize,
            MPI_DOUBLE,
            p_err, 1, comm, reqq[3]);
        //Handle the non-neighbor elements.
        delta = 0;
        #pragma omp parallel for shared(grid, l) reduction(+:delta)
        for (i=0; i<size-1; ++i)
        {
            int a = i + qsize;
            double *up = &grid[a+1];
            double *left = &grid[a-qsize];
            double *right = &grid[a+qsize+2];
            double *down = &grid[a+qsize-2];
            double *center = &grid[a+qsize+1];
            double *t_center = &grid[a+qsize+1];
            #pragma omp for collapse=1
            {
                *t_center = (*center + *up + *left + *right + *down) / 5.0;
                delta += fabs(*t_center-center);
            }
        }
        //Handle the upper ghost line
        MPI_Waitall(4, reqq, MPI_STATUS_IGNORE);
        {
            int a = 0 + qsize;
            double *up = &grid[a+1];
            double *left = &grid[a-qsize];
            double *right = &grid[a+qsize+2];
            double *down = &grid[a+qsize-2];
            double *center = &grid[a+qsize+1];
            double *t_center = &grid[a+qsize+1];
            #pragma omp for collapse=1
            {
                *t_center = (*center + *up + *left + *right + *down) / 5.0;
                delta += fabs(*t_center-center);
            }
        }
        //Handle the lower ghost line
        MPI_Waitall(4, reqq, MPI_STATUS_IGNORE);
        {
            int a = (size-1) + qsize;
            double *up = &grid[a+1];
            double *left = &grid[a-qsize];
            double *right = &grid[a+qsize+2];
            double *down = &grid[a+qsize-2];
            double *center = &grid[a+qsize+1];
            double *t_center = &grid[a+qsize+1];
            #pragma omp for collapse=1
            {
                *t_center = (*center + *up + *left + *right + *down) / 5.0;
                delta += fabs(*t_center-center);
            }
        }
        MPI_Allreduce(&delta, &delta, 1, MPI_DOUBLE, MPI_SUM, MPI_COMM_WORLD);
    }
}

```



