

Exploiting Multicore Processors and GPUs with OpenMP and OpenCL

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BSC

Clusters, Clouds, and Grids for Scientific Computing
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StarSs, OpenMP and OpenMPT

Programming models for accelerators

Handling accelerators and heterogeneity

Examples and results

Conclusions

StarSs

CellSs

SMPSSs

GPUSs

GridSs

ClusterSs

ClearSpeedSs

StarSs

- A “node” level programming model
- C/Fortran + directives
- Nicely integrates in hybrid MPI/StarSs
- Natural support for heterogeneity

• Portability

- “Same” source code runs on “any” machine
 - Optimized task implementations will result in better performance.
- “Single source” for maintained version of a application

Programmability

- Incremental parallelization/restructure
- Abstract/separate algorithmic issues from resources
- Disciplined programming

• Performance

- Asynchronous (data-flow) execution and loc awareness
- Intelligent Runtime: specific for each type of target platform.
 - Automatically extracts and exploits parallelism
 - Matches computations to resources

MAS
94

NANOS
~1996

GridSs
~2002

NANOS++
~2008

CompSs
~2007

CellSs
~2006

GPUsSs
~2009

SMP
~2009

SMPSS V1
~2007

```
add3 (float A[BS], float B[BS],  
      float C[BS]);  
  
cale_add (float sum, float A[BS],  
           float B[BS]);  
  
accum (float A[BS], float *sum);
```

```
=0; i<N; i+=BS) // C=A+B  
d3 (&A[i], &B[i], &C[i]);  
  
=0; i<N; i+=BS) // sum(C[i])  
um (&C[i], &sum);  
  
=0; i<N; i+=BS) // B=sum*A  
le_add (sum, &E[i], &B[i]);  
  
=0; i<N; i+=BS) // A=C+D  
d3 (&C[i], &D[i], &A[i]);  
  
=0; i<N; i+=BS) // E=G+F  
d3 (&G[i], &F[i], &E[i]);
```

```

ma css task input(A, B) output(C)
    radd3 (float A[BS], float B[BS],
           float C[BS]);
ma css task input(sum, A) inout(B)
    scale_add (float sum, float A[BS],
                float B[BS]);
ma css task input(A) inout(sum)
    accum (float A[BS], float *sum);

```

```

=0; i<N; i+=BS) // C=A+B
    radd3 ( &A[i], &B[i], &C[i]);

=0; i<N; i+=BS) // sum(C[i])
    sum (&C[i], &sum);

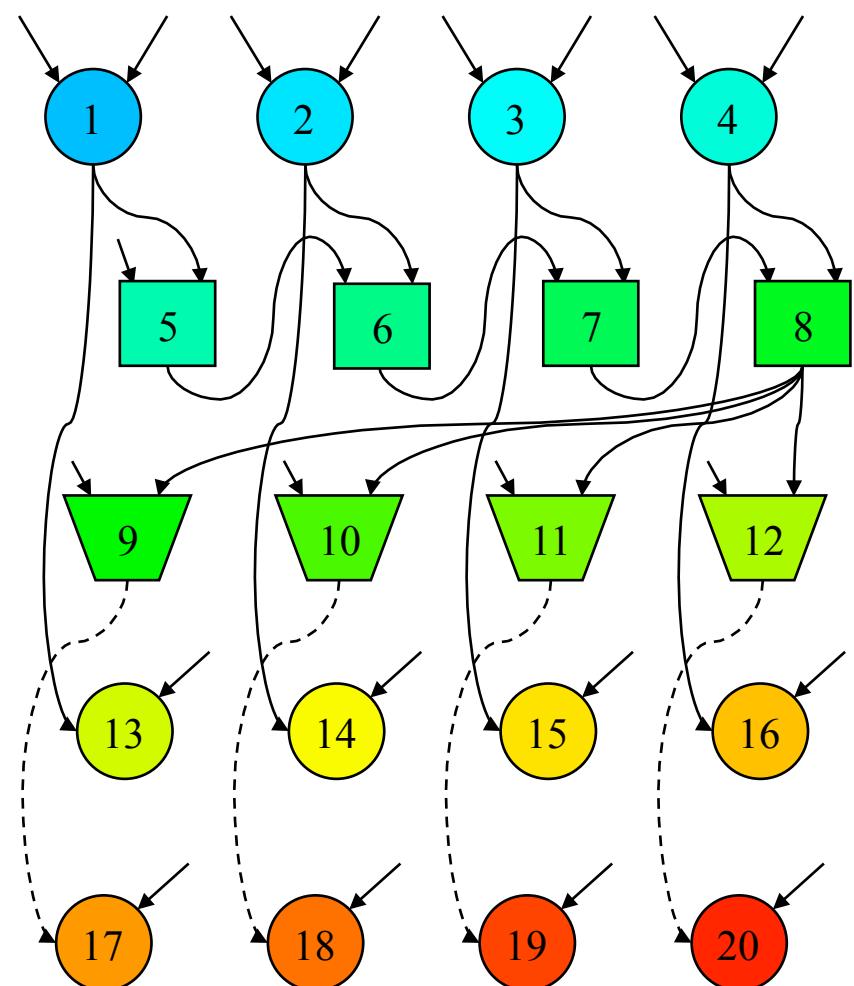
=0; i<N; i+=BS) // B=sum*A
    scale_add (sum, &E[i], &B[i]);

=0; i<N; i+=BS) // A=C+D
    radd3 (&C[i], &D[i], &A[i]);

=0; i<N; i+=BS) // E=G+F
    radd3 (&G[i], &F[i], &E[i]);

```

Compute dependences @ task instantiation



```

ma css task input(A, B) output(C)
add3 (float A[BS], float B[BS],
      float C[BS]);
ma css task input(sum, A) inout(B)
cale_add (float sum, float A[BS],
           float B[BS]);
ma css task input(A) inout(sum)
accum (float A[BS], float *sum);

```

```

=0; i<N; i+=BS) // C=A+B
d3 (&A[i], &B[i], &C[i]);

=0; i<N; i+=BS) // sum(C[i])
um (&C[i], &sum);

=0; i<N; i+=BS) // B=sum*A
le_add (sum, &E[i], &B[i]);

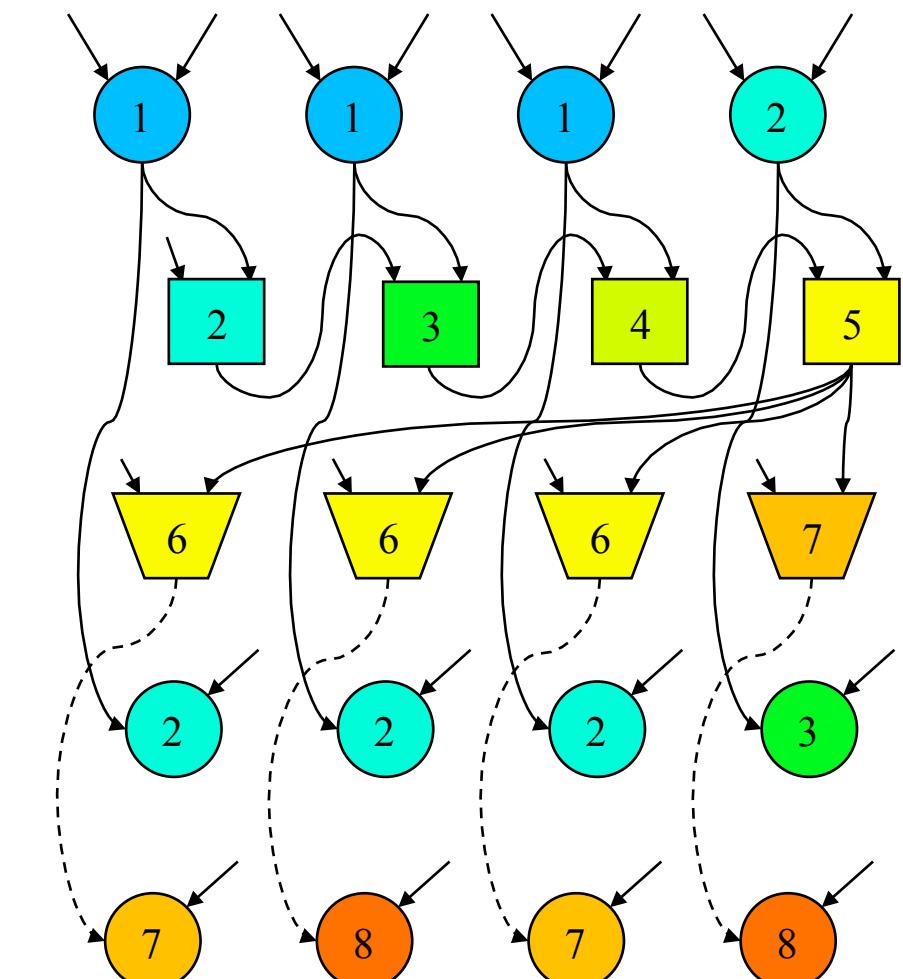
=0; i<N; i+=BS) // A=C+D
d3 (&C[i], &D[i], &A[i]);

=0; i<N; i+=BS) // E=G+F
d3 (&G[i], &F[i], &E[i]);

```

Decouple
how we write
from
how it is executed

Execute



```

na css task input(A, B) output(C)
radd3 (float A[BS], float B[BS],
      float C[BS]);
na css task input(sum, A) inout(B)
cale_add (float sum, float A[BS],
          float B[BS]);
na css task input(A) inout(sum) reduction(sum)
accum (float A[BS], float *sum);

```

```

=0; i<N; i+=BS)           // C=A+B
d3 (&A[i], &B[i], &C[i]);

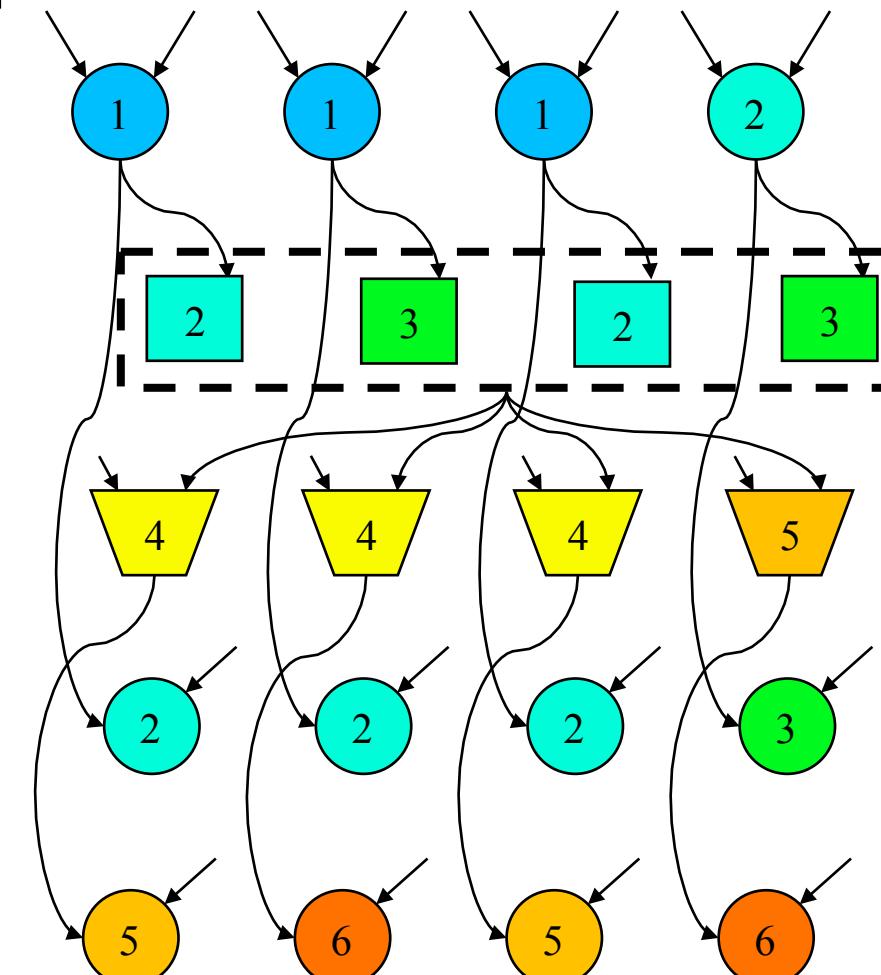
=0; i<N; i+=BS)           // sum(C[i])
um (&C[i], &sum);

=0; i<N; i+=BS)           // B=sum*A
le_add (sum, &E[i], &B[i]);

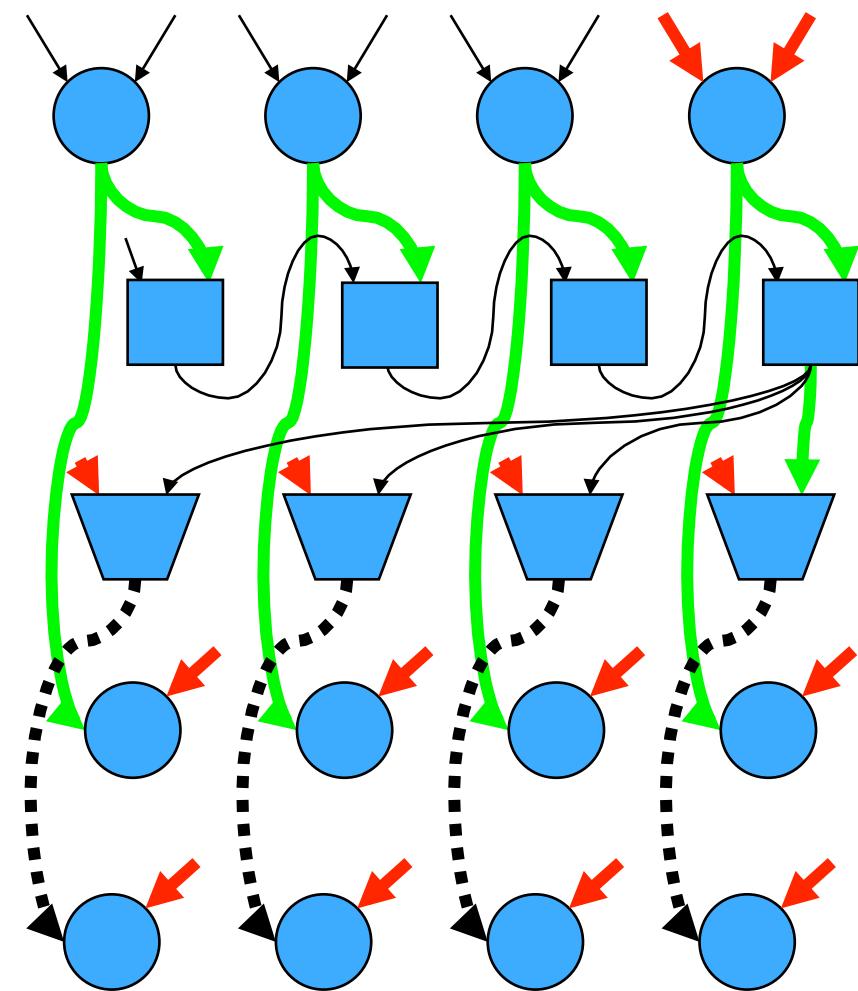
=0; i<N; i+=BS)           // A=C+D
d3 (&C[i], &D[i], &A[i]);

=0; i<N; i+=BS)           // E=G+F
d3 (&G[i], &F[i], &E[i]);

```



- Flat global address space seen by programmer
- Flexibility to dynamically traverse dataflow graph “optimizing”
 - Concurrency. Critical path
 - Memory access: data transfers performed by run time
- Opportunities for
 - Prefetch
 - Reuse
 - Eliminate antidependences (rename)
 - Replication management
 - Coherency/consistency handled by the runtime



```
pragma css task [input ( parameters ) ] \  
[output ( parameters ) ] \  
[inout ( parameters )] \  
[target device( [cell, smp, cuda] ) ] \  
[implements ( task_name ) ] \  
[reduction ( parameters ) ] \  
[ highpriority ]
```

```
pragma css wait on ( data_address )
```

```
pragma css barrier
```

```
pragma css mutex lock ( variable )
```

```
pragma css mutex unlock( variable )
```

parameters: parameter [, parameter]*
parameter: *variable_name* {[dimension]}

MP

t parallelism.

oin

provide some more flexibility

ality information. Global Addressing

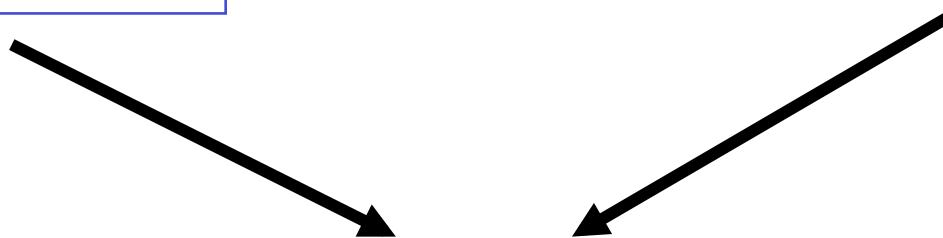
g

Sta

Implicit parallel
“atomic” ta

Explicit data access informa
Local addres

Single work gener

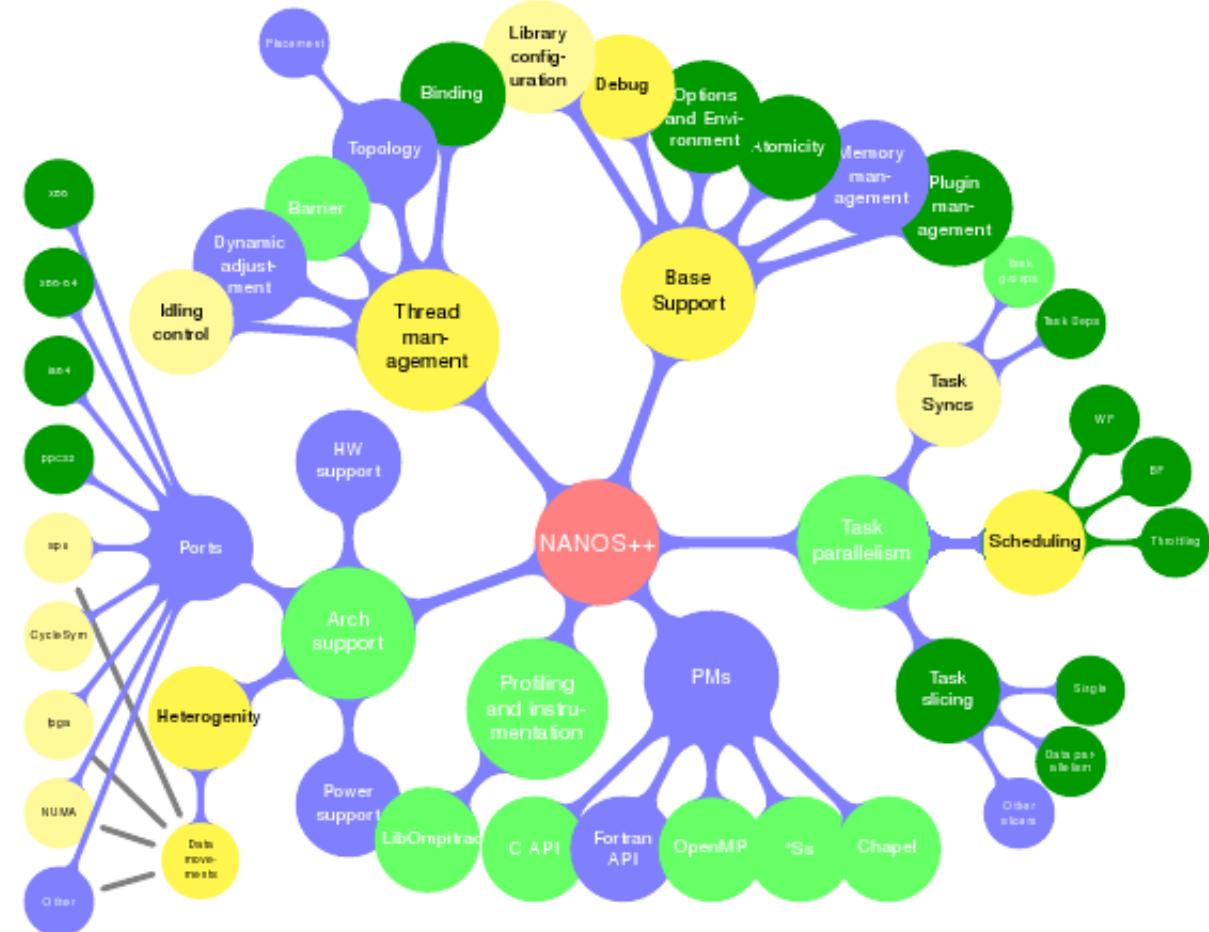


OpenMPT

, et al, “A Proposal to Extend the OpenMP Tasking Model for Heterogeneous Architectures” IWOMP 2009 & IJPP

et al, “Extending the OpenMP Tasking Model to Allow Dependent Tasks” IWOMP 2008, LNCS & IJPP

- Compiler (mercurium) and runtime (NANOS++)
- Support/integrate: OpenMP, StarSs, Chapel, ...



Status:

- Just started
- In progress
- Advanced
- Completed

Implicit kernel execution submission

```
/*Enqueue a kernel run call */
status = clEnqueueNDRangeKernel(
    commandQueue,
    kernel,
    2,
    NULL,
    globalThreads,
    localThreads,
    0,
    NULL,
    &events[0]);

if(!sampleCommon->checkVal(
    status,
    CL_SUCCESS,
    "clEnqueueNDRangeKernel failed."))
    return SDK_FAILURE;

/* wait for the kernel call to finish execution */
status = clWaitForEvents(1, &events[0]);
if(!sampleCommon->checkval(
    status,
    CL_SUCCESS,
    "clWaitForEvents failed."))
    return SDK_FAILURE;
```

Implicit kernel synchronization

execute the kernel over the entire range of our 1d input data set
using the maximum number of work group items for this device

```
global = count;
err = clEnqueueNDRangeKernel(commands, kernel, 1, NULL, &global, &local, 0, NULL,
);
(err)

printf("Error: Failed to execute kernel!\n");
return EXIT_FAILURE;
```

Host program deleting:

- Comments
- Some return code checks
- Checks of block sizes supported by device

```
eofday(&timev1, NULL);

ROR(clGetPlatformIDs(16, platforms, &num_platforms));
ROR(clGetPlatformInfo(platforms[0], CL_PLATFORM_PROFILE
e_buffer,&param_value_size_ret));
ROR(clGetDeviceIDs(platforms[0], device_type, 1, &device
ROR(clGetDeviceInfo(device_list,CL_DEVICE_PREFERRED_VEC
dth,NULL));
ROR(clGetDeviceInfo(device_list,CL_DEVICE_MAX_WORK_ITEM
ROR(clGetDeviceInfo(device_list,CL_DEVICE_GLOBAL_MEM_SI
ROR(clGetDeviceInfo(device_list,CL_DEVICE_LOCAL_MEM_SIZ
ROR(clGetDeviceInfo(device_list,CL_DEVICE_MAX_MEM_ALLOC
ROR(clGetDeviceInfo(device_list, CL_DEVICE_EXTENSIONS,
ROR(clGetDeviceInfo(device_list,CL_DEVICE_MAX_COMPUTE_U

: (ntasks == -1) ? max_compute_units : ntasks;
size = (cl_ulong) (max_alloc_size / 8);
groups = array_size / (vector_width * local_work_group_si
= clCreateContext(0, (cl_uint) 1, &device_list, NULL, N
e = clCreateCommandQueue(context,device_list, CL_QUEUE_
array_size;
array_size = (array_size < 16) ? 16 : array_size;
malign((void **) &rawbuf, 128, sizeof(cl_double) * 7 *
= &rawbuf[0 * sizeof(cl_double) * malloc_array_size]
= &rawbuf[1 * sizeof(cl_double) * malloc_array_size]
= &rawbuf[2 * sizeof(cl_double) * malloc_array_size]
= &rawbuf[3 * sizeof(cl_double) * malloc_array_size]
```

```
cpflag_fptr = (unsigned int *) cpflag; S0_fptr = (float *) S0; K_fptr = (float *)
r_fptr = (float *) r; sigma_fptr = (float *) sigma; T_fptr = (float *) T;
answer_fptr = (float *) answer;
int memsize = array_size * (double_flag ? sizeof(double) : sizeof(float));
memobjs[0] = clCreateBuffer(context, CL_MEM_USE_HOST_PTR, memsize, cpflag, &rc);
memobjs[1] = clCreateBuffer(context, CL_MEM_USE_HOST_PTR, memsize, S0, &rc);
memobjs[2] = clCreateBuffer(context, CL_MEM_USE_HOST_PTR, memsize, K, &rc);

memobjs[3] = clCreateBuffer(context, CL_MEM_USE_HOST_PTR, memsize, r, &rc);
memobjs[4] = clCreateBuffer(context, CL_MEM_USE_HOST_PTR, memsize, sigma, &rc);
memobjs[5] = clCreateBuffer(context, CL_MEM_USE_HOST_PTR, memsize, T, &rc);
memobjs[6] = clCreateBuffer(context, CL_MEM_USE_HOST_PTR, memsize, answer, &rc);

bs_source = load_program (kernel_source_file, &rc);

program = clCreateProgramWithSource (context, 1, (const char**) (&bs_source), NULL

if (cod
-DRANGE_L
if (cod
char nu
sprintf
strcat(
strcat(
rc = clEnqueueNDRangeKernel(cmd_queue, kernel, 1, NULL,
global_work_size[0] = (size_t) (n_workgroups*local_work_group_size);
local_work_size[0] = (size_t) local_work_group_size;
sprintf
strcat(
strcat(
if(rc != CL_SUCCESS) {
fprintf(stderr, "Executing the kernel failed...rc=%d\n",rc);
clWaitForEvents((cl_uint) 1, event);
rc = cl
return 0;
kernel }

rc = clSetKernelArg(kernel, 0, sizeof(cl_mem), (void *) &memobjs[0]);
rc = clSetKernelArg(kernel, 1, sizeof(cl_mem), (void *) &memobjs[1]);
rc = clSetKernelArg(kernel, 2, sizeof(cl_mem), (void *) &memobjs[2]);
rc = clSetKernelArg(kernel, 3, sizeof(cl_mem), (void *) &memobjs[3]);
rc = clSetKernelArg(kernel, 4, sizeof(cl_mem), (void *) &memobjs[4]);
rc = clSetKernelArg(kernel, 5, sizeof(cl_mem), (void *) &memobjs[5]);
rc = clSetKernelArg(kernel, 6, sizeof(cl_mem), (void *) &memobjs[6]);

gettimeofday(&timev4, NULL);
```

```
bsop_ref (unsigned int4 cpflag, float4
          float4 r, float4 sigma, float4 T
t4 d1, d2, Nd1, Nd2, expval, k1, n1, k2
          accum2, candidate_answer1, candidate
          flag1, flag2;
```

```
log(S0/K) + (r + HALF * sigma*sigma)*T
= (sigma * sqrt(T));
al = exp(ZERO - r * T);
d1 = sigma * sqrt(T);
1 = (d1 < ZERO);
2 = (d2 < ZERO);
fabs(d1);
fabs(d2);
ONE / (ONE + NCDF * d1);
ONE / (ONE + NCDF * d2);
m1 = A4 + A5 * k1;
m2 = A4 + A5 * k2;
m1 = k1 * accum1 + A3;
m2 = k2 * accum2 + A3;
m1 = k1 * accum1 + A2;
m2 = k2 * accum2 + A2;
m1 = k1 * accum1 + A1;
m2 = k2 * accum2 + A1;
m1 = k1 * accum1;
m2 = k2 * accum2;
exp(ZERO - HALF * d1 * d1);
exp(ZERO - HALF * d2 * d2);
= INV_ROOT2PI;
= INV_ROOT2PI;
candidate_answer1 = ONE - n1 * accum1;
candidate_answer2 = ONE - n2 * accum2;
= SELECT(candidate_answer1, (ONE - candidate
= SELECT(candidate_answer2, (ONE - candidate
= S0 * Nd1 - K * expval * Nd2;
```

```
__kernel __attribute__((reqd_work_group_size(LWGSIZE, 1, 1)))
void bsop_kernel (__global unsigned int4 *dm_cpflag,
                  __global float4 *dm_S0,
                  __global float4 *dm_K,
                  __kernel __attribute__((reqd_work_group_size(1, 1, 1)))
void bsop_kernel (__global const FIXED *cpflag_dm,
                  __global const FLOAT *S0_dm,
                  __global const FLOAT *K_dm,
                  __global const FLOAT *r_dm,
                  __global const FLOAT *sigma_dm,
                  __global const FLOAT *T_dm,
                  __global FLOAT *answer_dm,
                  int task_id,
                  __local FIXED *lm_cpflag,
                  __local FLOAT *lm_S0,
                  __local FLOAT *lm_K,
                  __local FLOAT *lm_r,
                  __local FLOAT *lm_sigma,
                  __local FLOAT *lm_T,
                  __local FLOAT *answer,
                  int n, size_t stride) {
    int i, j;
    event_t event;
    stride >>= STRIDESHIFT;

    for (j = n*task_id; j < n*(task_id+1); j+=stride) {
        event = async_work_group_copy(lm_cpflag, (cpflag_dm +
        stride, (event_t) 0);
        event = async_work_group_copy(lm_S0, (S0_dm+j), stride,
                                      ...
                                      wait_group_events(1, &event);
        for(i = 0; i < stride; i++)
            answer[i] = bsop_ref(lm_cpflag[i], lm_S0[i], lm_K[i],
r[i], lm_sigma[i], lm_T[i]);
```

```
op_test (int double_flag, cl_ulong array_size, cl_ulong local_work_group_size)
```

```
, bsize;
```

```
neofday(&timev5, NULL);
```

```
i=0; i<array_size; i+=local_work_group_size) {  
    ze= ((i+local_work_group_size)>array_size) ? array_size - i : local_work_group_size  
    bsop_ref_float (&bsize, &cpflag_fptr[i], &S0_fptr[i], &K_fptr[i], &r_fptr[i],  
                    &sigma_fptr[i], &T_fptr[i],  
                    &answer_fptr[i]);
```

```
a css barrier
```

```
neofday(&timev6, NULL);
```

```
float bsop_reference_float (unsigned int cpflag, float S0, float K, float r, float sigma, float T) {  
    float d1, d2, c, p, Nd1, Nd2, expval, answer;  
    d1 = logf(S0/K) + (r + 0.5*sigma*sigma)*T;  
    d1 /= (sigma * sqrt(T));  
    expval = exp(-r * T);  
    d2 = d1 - sigma * sqrt(T);  
    Nd1 = Nf(d1); Nd2 = Nf(d2);  
    c = S0 * Nd1 - K * expval * Nd2;  
    p = K * expval * (1.0 - Nd2) - S0 * (1.0 - Nd1);  
    answer = cpflag ? c : p;  
    return answer;  
}
```

```
s task input(size, cpflag_fptr[size], S0_fptr[size], K_fptr[size], r_fptr[size], \  
            sigma_fptr[size], T_fptr[size]) \  
    output (answer_fptr[size])
```

```
ref_float (cl_ulong size, unsigned int * cpflag_fptr, float * S0_fptr, float * K_fptr,  
          float * r_fptr, float * sigma_fptr, float * T_fptr, float * answer_fptr)
```

```
0; i < size; i++) {
```

```
_fptr[i] = bsop_reference_float (cpflag_fptr[i], S0_fptr[i],
```

```

c bsop_test (int double_flag, cl_ulong array_size, cl_ulong local_work_group_size)

signed long long esp;
c i;

ctimeofday(&timev5, NULL);

c ii;
for (ii = 0; ii < array_size; ii+=local_work_group_size) {
cna omp task private (i)
    for (i=ii; (i<ii+local_work_group_size) && (i<array_size); i+=OCLN) {
        answer_fptr[i] = bsop_reference_float(cpflag_fptr[i], S0_fptr[i], K_fptr[i],
                                              r_fptr[i], sigma_fptr[i], T_fptr[i]);
    }
}
cna omp taskwait

ctimeofday(&timev6, NULL);

return 0;

```

```

float bsop_reference_float (unsigned int cpflag, float S0, float K, float r, float sigma, float T) {
    float d1, d2, c, p, Nd1, Nd2, expval, answer;
    d1 = logf(S0/K) + (r + 0.5*sigma*sigma)*T;
    d1 /= (sigma * sqrt(T));
    expval = exp(-r * T);
    d2 = d1 - sigma * sqrt(T);
    Nd1 = Nf(d1); Nd2 = Nf(d2);
    c = S0 * Nd1 - K * expval * Nd2;
    p = K * expval * (1.0 - Nd2) - S0 * (1.0 - Nd1);
    answer = cpflag ? c : p;
    return answer;
}

```

- Same main program, just another **implementation of the task**
- Using **OpenCL clean SIMD code** but not OpenCL kernel declarations
- No need for manual overlap between computation and transfers

```

float4 bsop_ref (unsigned int4 cpflag, float4 S0, float4 K,
                 float4 r, float4 sigma, float4 T) {
float4 d1, d2, Nd1, Nd2, expval, k1, n1, k2, n2, accum1,
        accum2, candidate_answer1, candidate_answer2, call,
int4 flag1, flag2;

d1 = log(S0/K) + (r + HALF * sigma*sigma)*T;
d1 /= (sigma * sqrt(T));
expval = exp(ZERO - r * T);
d2 = d1 - sigma * sqrt(T);
flag1 = (d1 < ZERO);
flag2 = (d2 < ZERO);
d1 = fabs(d1);
d2 = fabs(d2);
k1 = ONE / (ONE + NCDF * d1);
k2 = ONE / (ONE + NCDF * d2);
accum1 = A4 + A5 * k1;
accum2 = A4 + A5 * k2;

```

```

gma css task input(size, cpflag_fptr[size], S0_fptr[size], K_fptr[size], r_fptr[size], \
                   sigma_fptr[size], T_fptr[size]) \
                   output (answer_fptr[size])

bsop_ref_float (cl_ulong size, unsigned int * cpflag_fptr, float * S0_fptr, float * K_fptr,
                  float * r_fptr, float * sigma_fptr, float * T_fptr, float * answer_fptr)

typedef __attribute__((vector_size(16))) float float4;
typedef __attribute__((vector_size(16))) int int4;
float4 bsop_ref (int4, float4, float4, float4, float4, float4);

for (int i = 0; i < size; i+=4) {
    *((float4 * &answer_fptr[i])) = bsop_ref (*((int4 *) &cpflag_fptr[i]), *((float4 *) &S0_fptr[i]),
                                                 *((float4 *) &K_fptr[i]), *((float4 *) &r_fptr[i]),
                                                 *((float4 *) &sigma_fptr[i]), *((float4 *) &T_fptr[i]));
}

```

main program code

```
int main(void) {
// Allocate and initialize the matrices
    Matrix M = AllocateMatrix(WIDTH, WIDTH, 1);
    Matrix N = AllocateMatrix(WIDTH, WIDTH, 1);
    Matrix P = AllocateMatrix(WIDTH, WIDTH, 0);

// M * N on the device
    MatrixMulOnDevice(M, N, P);

// Free matrices
    FreeMatrix(M);
    FreeMatrix(N);
    FreeMatrix(P);
return 0;
}
```

multiplication (host side)

```
MatrixMulOnDevice(float* M, float* N, float* P, int Width)

size = Width * Width * sizeof(float);
float* Md, Nd, Pd;

Allocate and Load M, N to device memory
daMalloc(&Md, size);
daMemcpy(Md, M, size, cudaMemcpyHostToDevice);

daMalloc(&Nd, size);
daMemcpy(Nd, N, size, cudaMemcpyHostToDevice);

Allocate P on the device
daMalloc(&Pd, size);

Setup the execution configuration
m3 dimGrid(1, 1);
m3 dimBlock(Width, Width);

Launch the device computation threads!
MatrixMulKernel<<<dimGrid, dimBlock>>>(Md, Nd, Pd, Width);
daMemcpy(P, Pd, size, cudaMemcpyDeviceToHost);

Free device matrices
daFree(Md); cudaFree(Nd); cudaFree(Pd);
```

Matrix multiplication (device side)

```
__global__ void MatrixMulKernel(Matrix M, Matrix N, Matrix P)
{
    // 2D Thread ID
    int tx = threadIdx.x;
    int ty = threadIdx.y;

    // Pvalue is used to store the element of the matrix
    // that is computed by the thread
    float Pvalue = 0;
    for (int k = 0; k < M.width; ++k)
    {
        float Melement = M.elements[ty * M.pitch + k];
        float Nelement = Nd.elements[k * N.pitch + tx];
        Pvalue += Melement * Nelement;
    }
    // Write the matrix to device memory
}
```

program code

```
main( void ){  
    ...  
    for ( i = 0; i < N; i++ ) {  
        for ( j = 0; j < N; j++ ) {  
            for ( k = 0; k < N; k++ ) {  
                matmul_tile ( A[i][k], B[k][j], C[i][j] );  
            }  
        }  
    }  
}
```

Main program:

- No explicit data transfers or allocation
- No explicit execution configuration
- The same StarSs main program can be used

Task (device)

```
__global__ void matmul_cuda ( float * A, float * B, float * C, int wA, int  
wB ) {  
    int bx = blockIdx.x;    int by = blockIdx.y;  
    int tx = threadIdx.x;  int ty = threadIdx.y;  
  
    int aBegin = wA * BLOCK_SIZE * by;  
    int aEnd = aBegin + wA - 1;    int aStep = BLOCK_SIZE;  
    int bBegin = BLOCK_SIZE * bx;      int bStep = BLOCK_SIZE * wB;  
    float Csub = 0;  
  
    for( int a = aBegin, b = bBegin; a <= aEnd; a += aStep, b += bStep ) {  
        __shared__ float As[ BLOCK_SIZE ][ BLOCK_SIZE ];  
        __shared__ float Bs[ BLOCK_SIZE ][ BLOCK_SIZE ];  
        As[ ty ][ tx ] = A[ a+wA * ty + tx ];  
        Bs[ ty ][ tx ] = B[ b+wB * ty + tx ];  
        __syncthreads();  
        for( int k = 0; k < BLOCK_SIZE; k++ ) {  
            Csub += As[ ty ][ k ] * Bs[ k ][ tx ];  
        }  
        __syncthreads();  
    }  
}
```

(host side)

```
ma css task input(A[BS][BS], B[BS][BS]) inout( C[BS][BS] )  
ma css target device (CUDA)  
matmul_tile (float *A, float *B, float *C ){  
matmul_cuda <<<dimGrid dimBlock>>>(A, B, C, BS, BS);
```

main program code

```
void matmul ( int m, int l, int n, int mDIM , in
  int nDIM ,
  float ** A, float ** B, float ** C)
{
  for (i = 0;i < mDIM ; i++) {
    for (j = 0; j < nDIM ; j++) {
      for (k = 0; k < lDIM ; k++) {
        matmul_block (A[i* lDIM +k],B[k* nDIM +
          C[i* nDIM +j ] );
      }
    }
  }
  # pragma omp taskwait
}
```

x multiplication: multiple kernel implementations

```
: int NB = 512;
# pragma omp target device (smp , cell ) copy_deps
# pragma omp task inout ([ NB*NB] C) input ([ NB*NB] A, [NB*NB] B)
matmul_block ( float * A, float * B, float * C)

openCL kernel

# pragma omp target device ( cuda ) copy_deps implement (matmul_block)
matmul_block_gpu ( float * A, float * B, float * C)

CUDA kernel
```

Dependences: not all arguments in directionality clauses

```
    pragma omp task inout(C[BS][BS])  
    d matmul( float *A, float *B, float *C) {  
        // original sequential code  
    }
```

Heterogeneous devices

```
    pragma omp target device(cuda) implements(matmul)  
    copy_in(A[BS][BS], B[BS][BS], C[BS][BS]); copy_out(C[BS][BS])  
    d matmul_cuda( float *A, float *B, float *C) {  
        // optimized kernel for cuda  
    }
```

Different implementations

```
    pragma omp target device(cell) implements(matmul)  
    copy_in(A[BS][BS], B[BS][BS], C[BS][BS]); copy_out(C[BS][BS])  
    d matmul_spe( float *A, float *B, float *C);
```

Separation dependences/transfers

```
    at *A[NB][NB], *B[NB][NB], *C[NB][NB];
```

```
main( void ){  
    (int i = 0; i < NB; i++)  
    for (int j = 0; j < NB; j++)  
        for (int k = 0; k < NB; k++)  
            matmul (A[i][k], B[k][j], C[i][j]);
```

```

factorization A = LU, overwriting A with the triangular factors */
u_getrf( float *A );

angular system solve B := B * inv(A), with A upper triangular */
u_trsm_right( float *A, float *B );

angular system solve B := inv(A) * B, with A unit lower triangular */
u_trsm_left( float *A, float *B );

matrix multiplication C = C - A * B */
u_gemm( float *A, float *B, float *C );

parse_LU() {
    int k = 0; k < NB; k++ ) {
        #pragma omp task inout( A[k][k][0:BS-1][0:BS-1] )
        getrf( A[k][k] );
        ( int i = k+1; i < NB; i++ )
        if ( A[i][k] != NULL )
            #pragma omp task input( A[k][k][0:BS-1][0:BS-1] ) \
                inout( A[i][k][0:BS-1][0:BS-1] )
            lu_trsm_right( A[k][k], A[i][k] );

        ( int j = k+1; j < NB; j++ ) {
            if ( A[k][j] != NULL )
                #pragma omp task input( A[k][k][0:BS-1][0:BS-1] ) \
                    inout( A[k][j][0:BS-1][0:BS-1] )
                lu_trsm_left( A[k][k], A[k][j] );
            for ( int i = k+1; i < NB; i++ )
                if ( A[i][k] != NULL ) {
                    if ( A[i][j] == NULL)
                        A[i][j] = allocate_clean_block();
                    #pragma omp task input( A[i][k][0:BS-1][0:BS-1], \
                        A[k][j][0:BS-1][0:BS-1] ) \
                        inout( A[i][j][0:BS-1][0:BS-1] )
                    lu_gemm( A[i][k], A[k][j], A[i][j] );
                }
        }
    }
}

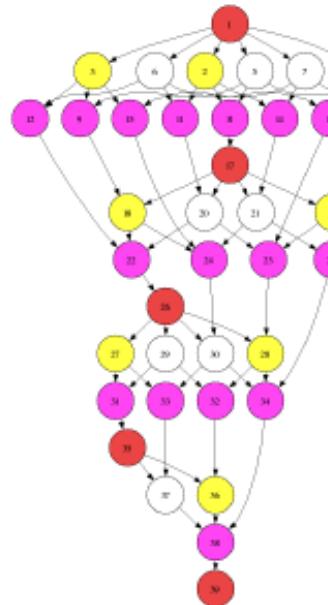
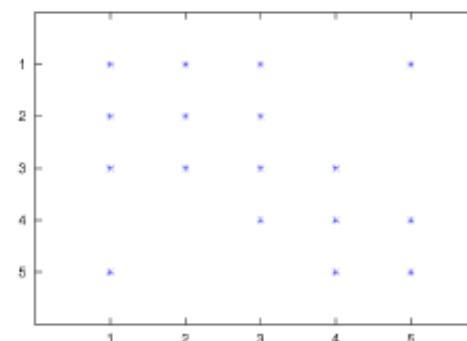
```

Inline directives:

saves manual outlining !!!

Tasks have no name

→ not multiple implementations



```

Cholesky factorization A = LL^T, overwriting the lower triangle of A with L */
gma omp task inout( A[0:BS-1][0:BS-1] )
| chol_potrf( float *A );

Triangular system solve B = B * inv(A)^T, with A lower triangular */
gma omp target device( cuda ) copy_in ( A[0:BS-1][0:BS-1], B[0:BS-1][0:BS-1] ) \
| copy_out( B[0:BS-1][0:BS-1] )
gma omp task input( A[0:BS-1][0:BS-1] ) inout( B[0:BS-1][0:BS-1] )
| chol_trsm_right( float *A, float *B );

Matrix multiplication C = C - A * B^T */
gma omp target device( cuda ) copy_in ( A[0:BS-1][0:BS-1], B[0:BS-1][0:BS-1], \
| C[0:BS-1][0:BS-1] ) \
| copy_out( C[0:BS-1][0:BS-1] )
gma omp task input( A[0:BS-1][0:BS-1], B[0:BS-1][0:BS-1] ) \
| inout( C[0:BS-1][0:BS-1] )
| chol_gemm( float *A, float *B, float *C );

Symmetric rank-BS update C = C - A * A^T */
gma omp target device( cuda ) copy_in ( A[0:BS-1][0:BS-1], C[0:BS-1][0:BS-1] ) \
| copy_out( C[0:BS-1][0:BS-1] )
gma omp task input( A[0:BS-1][0:BS-1] ) inout( C[0:BS-1][0:BS-1] )
| chol_syrk( float *A, float *C );

Cholesky() {
    i, j, k;

    ( k = 0; k < NB; k++ ) {
        chol_potrf( A[k][k] );
        for ( i = k+1; i < NB; i++ )
            chol_trsm_right( A[k][k], A[i][k] );

        for ( i = k+1; i < NB; i++ ) {
            for ( j = k+1; j < i; j++ )
                chol_gemm( A[i][k], A[j][k], A[i][j] );
            chol_syrk( A[i][k], A[i][i] );
        }
    }
}

```

Annotated function declaration
ALL instances become tasks

```
a[N];  
  
#pragma omp task output(a[0:N/2-1])  
a(a); //Task A: generate the bottom half of the array  
  
#pragma omp task output(a[N/2:N-1])  
b(a); //Task B: generate the upper half of the array  
  
#pragma omp task input(a[0:N-1])  
c(a); //Task C: use the array a
```

6 Simple example of array sections

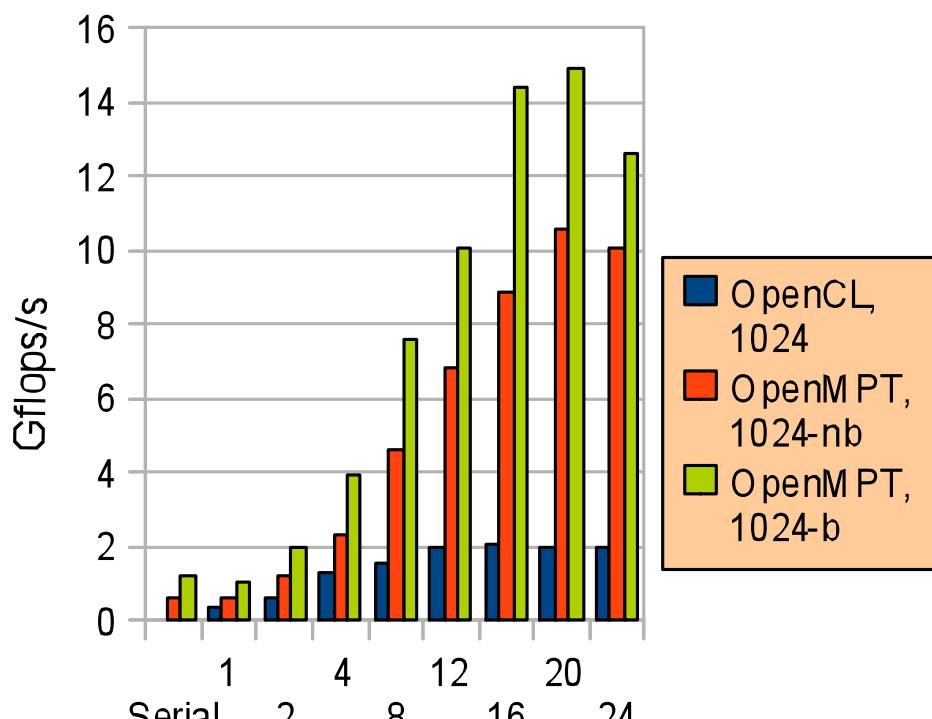
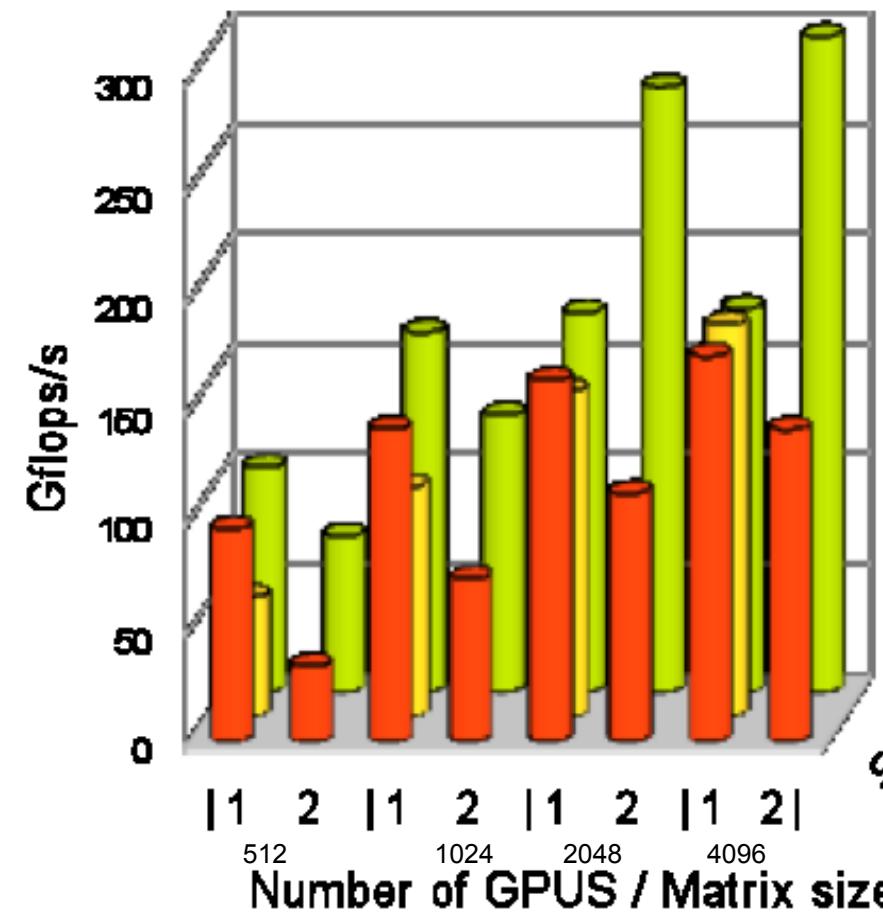
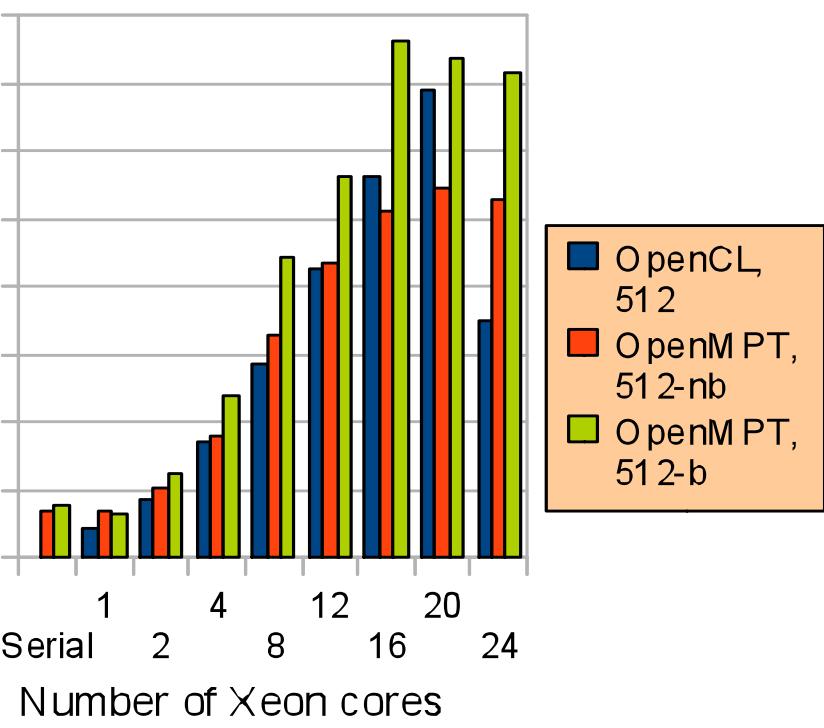
```
d f (int *a, int **b, int c[N][N])  
  
#pragma omp task input(a) //refers to the pointer a  
fa(a);  
#pragma omp task input([N] a) \  
  output([N] a[N/2:]) \  
  inout([N][N] b[:,1:10]) // refers to the N elements pointed by a  
  // refers to the elements N/2 to N-1  
  // refers to the submatrix [0..N-1][1..10]  
fb(a,b); // pointed by b  
  
#pragma omp task input(c) \  
  inout([N][N] c) //refers to the matrix of NxN pointed by c  
fc(c);
```

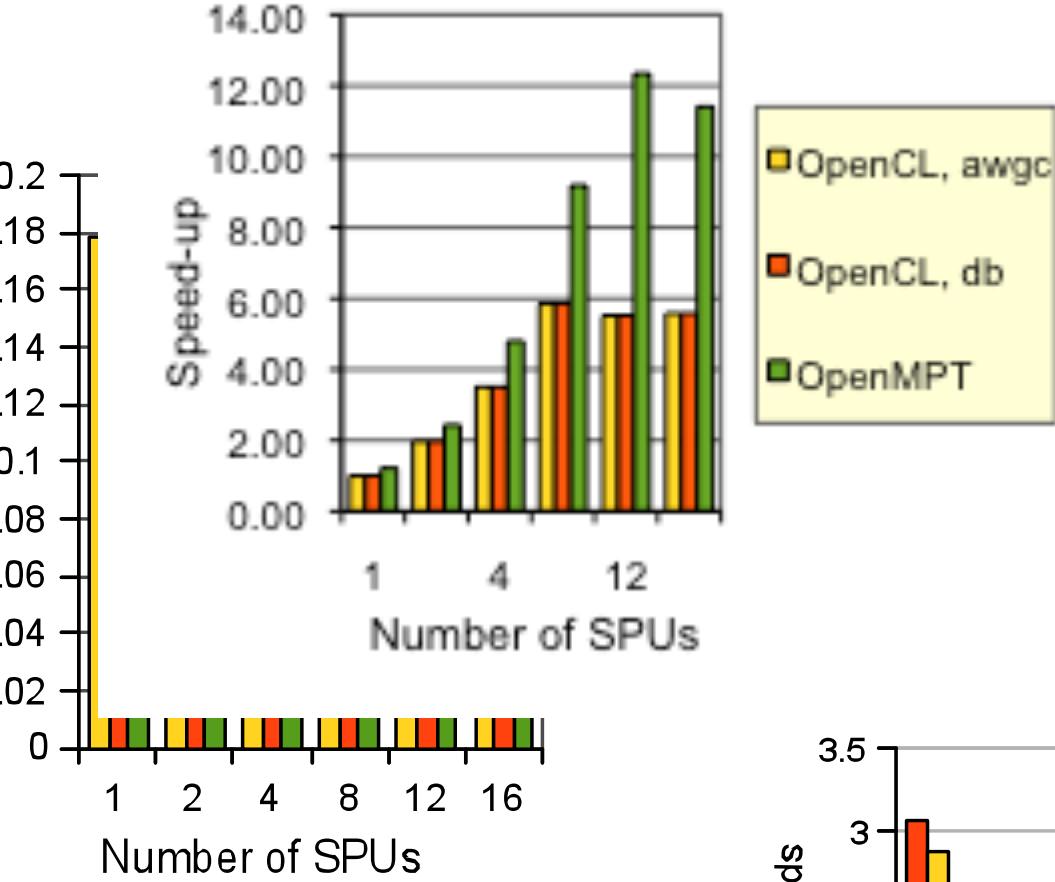
• Models

Model	Cell /B.E.	x86_64	GPUs
opencl	IBM OpenCL SDK	AMD/ATI OpenCL SDK	NVIDIA OpenCL SDK
StarSs	CellSs	Nanos++	Nanos++ (CUDA kernels)
		SMPSS	
OpenMPT	-	Nanos++	Nanos++ (CUDA kernels)

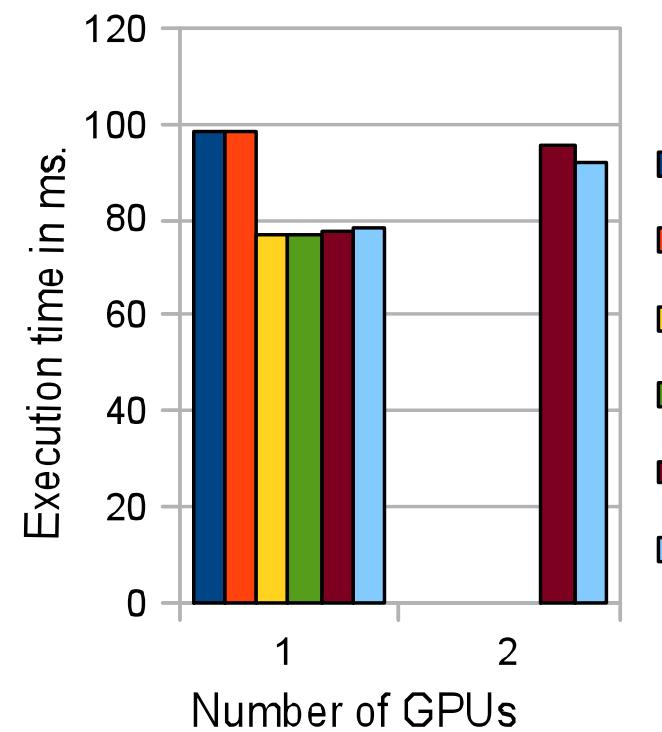
• Benchmarks

- Matrix Multiply
- BlackScholes – Computes pricing of European-style options
- Perlin Noise – Computes an image filled with noise to improve realistic view of moving graphics
- Julia set – Computes a set of images of the Julia Set fractal

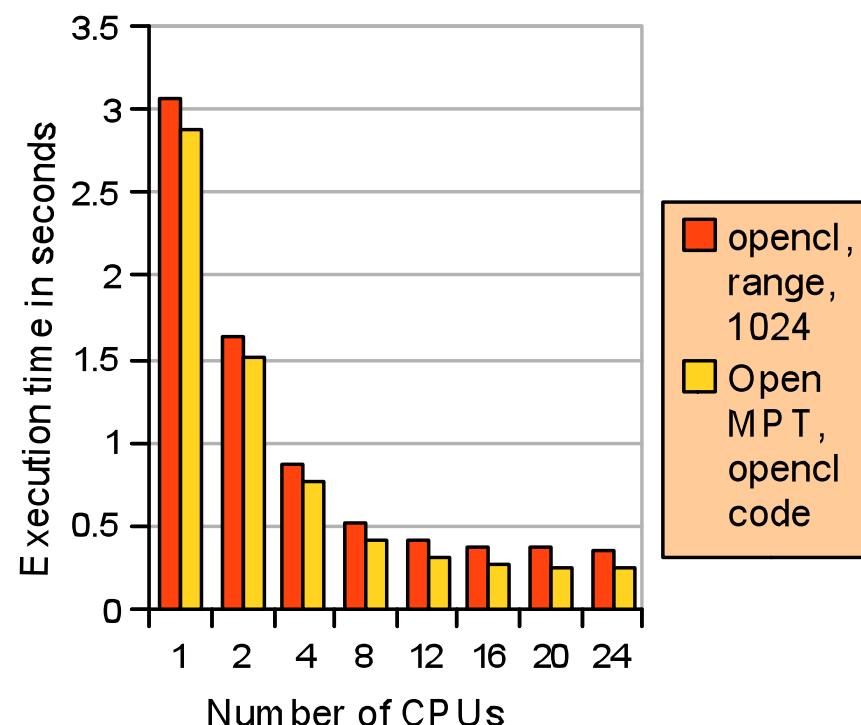




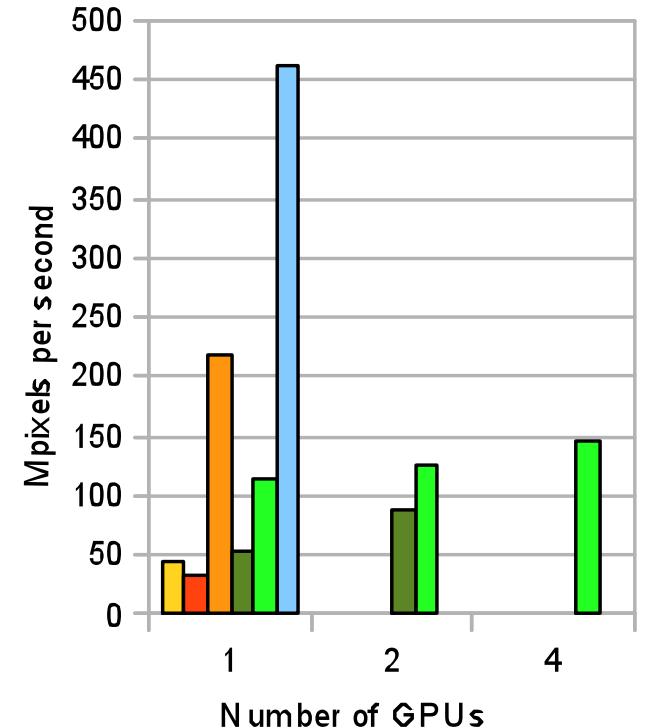
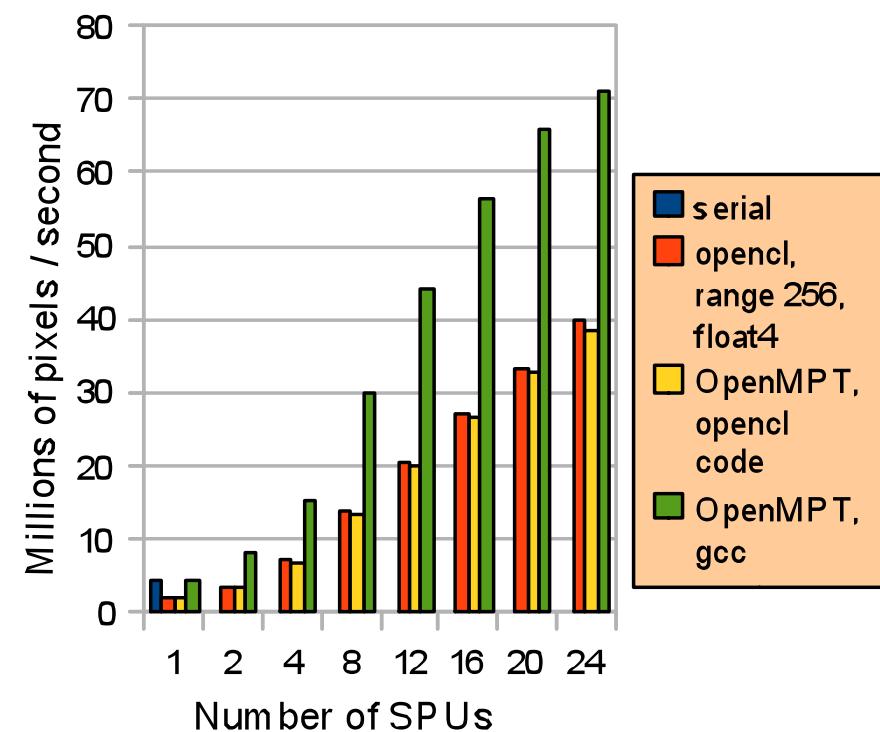
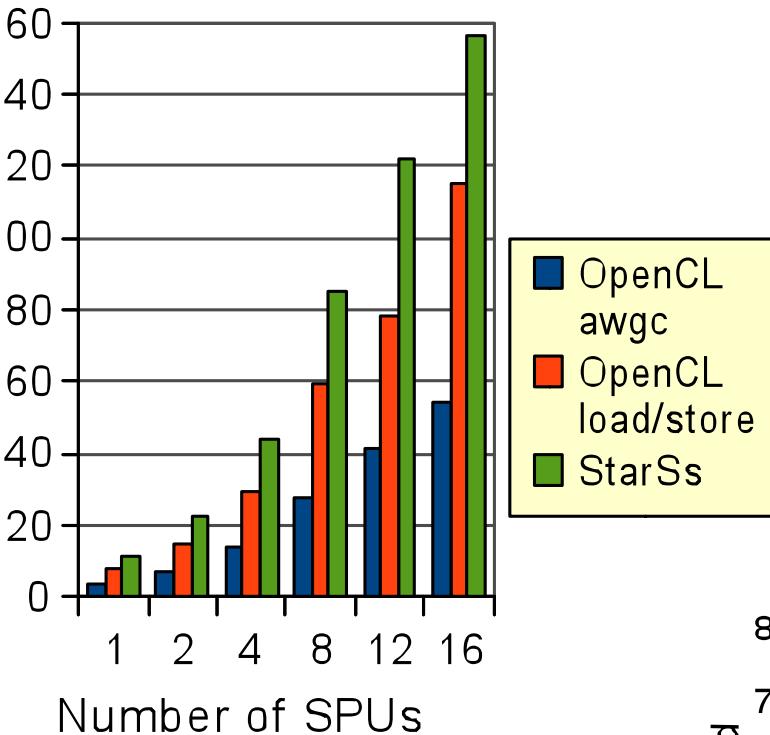
Cell BE

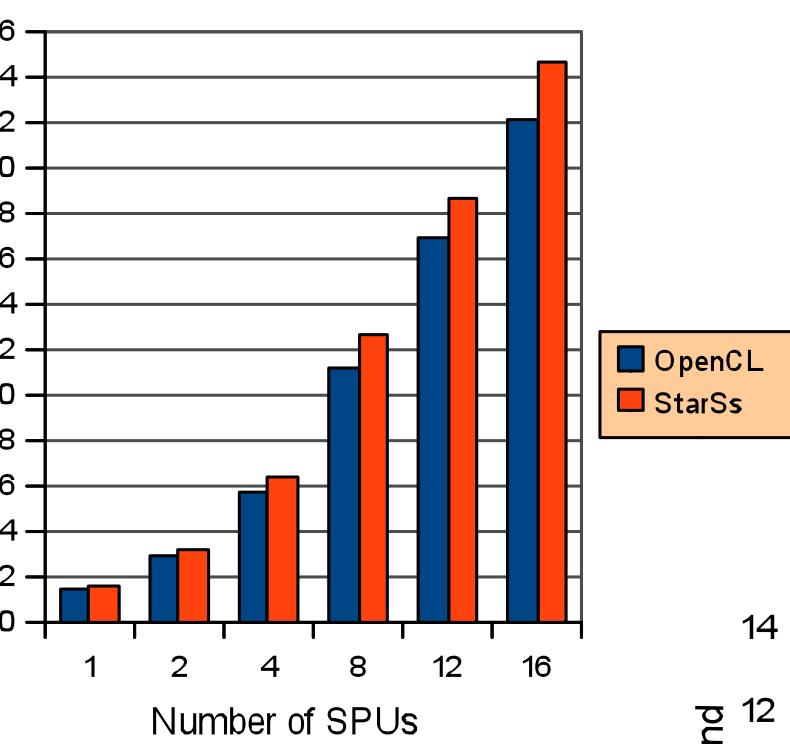


**Nvidia GPU
GTX 285**

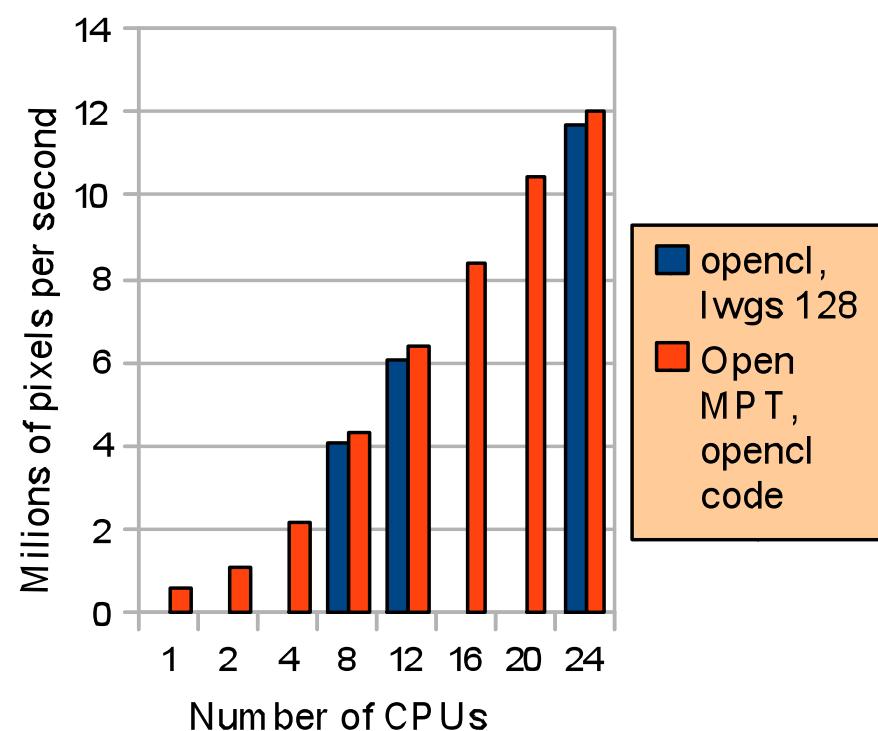
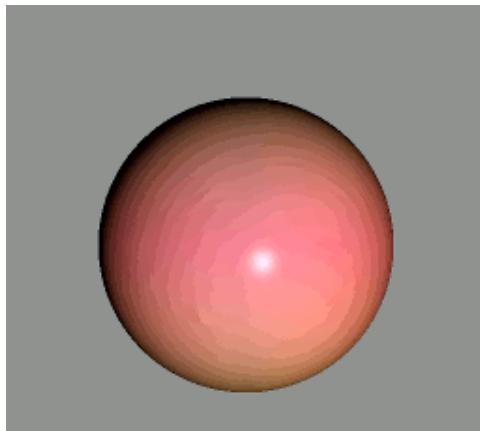


Intel Xeon

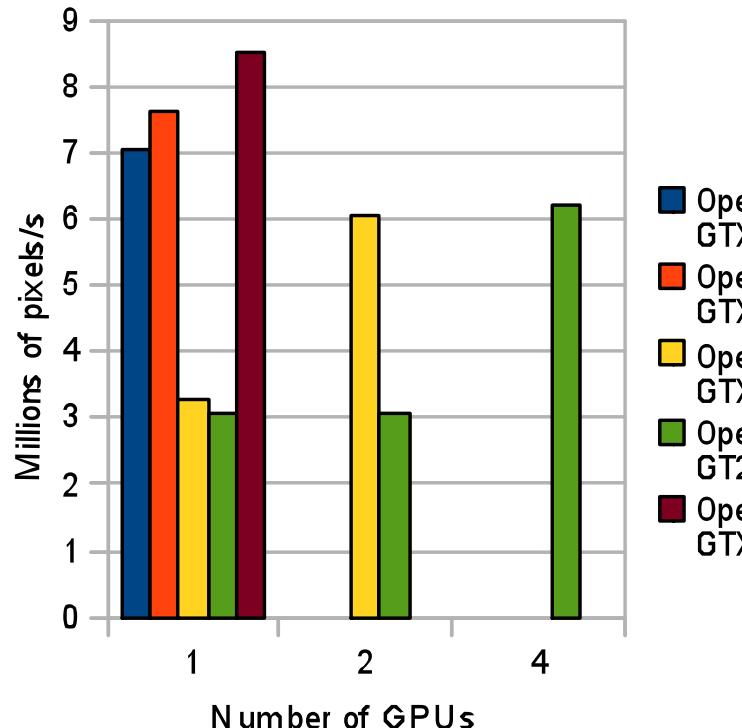




Cell BE



Intel Xeon



Nvidia GPU
GTX 285

- OpenMPT integrates ideas from StarSs and OpenMP
 - Support for task dependences, enabling data-flow like execution and exploitation of local memory
 - Support for heterogeneity, increasing the portability of applications by means of specialized kernels for each architecture
 - Based on source code to source compilation and intelligent runtimes
- Distributed as open source:
 - StarSs releases:
 - CellSs: www.bsc.es/cellsuperscalar
 - SMPSS: www.bsc.es/smpsuperscalar
 - OpenMPT: http://www.bsc.es/plantillaG.php?cat_id=328