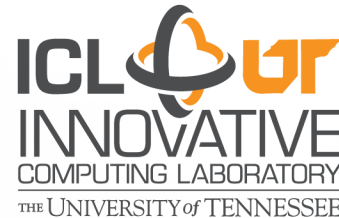


Creating a new operation with DPLASMA: a step by step guide

Anthony Danalis

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PaRSEC & DPLASMA

- PaRSEC is the runtime engine
 - It can be used to implement affine DLA operations
 - Or any operation that can be expressed as a static PTG
 - Or any operation that has a fixed DAG upon startup
 - Or anything else!
- DPLASMA is PLASMA implemented on top of PaRSEC
 - It's just sweet.

PaRSEC team and (re)sources

- ***George Bosilca***
- ***Aurelien Bouteiller***
- ***Anthony Danalis***
- ***Jack Dongarra***
- Mathieu Faverge
- ***Thomas Herault***
- ***Pierre Lemarinier***
- **Stephanie Moreaud**

ICL project site:

icl.cs.utk.edu/projectsdev/parsec/

Mercurial repository:

bitbucket.org/bosilca/dplasma

Getting started

- Download icl.cs.utk.edu/projectsfiles/dague/dplasma-1.0.0.tar.bz2
 - Top level view:
 - CMakeLists.txt
 - CTestConfig.cmake
 - Doxyfile
 - License.txt
 - cmake_modules
 - contrib
 - data_dist
 - **dplasma**
 - include
 - src
 - tests
 - tools
- 
- CMakeLists.txt
 - cores
 - include
 - lib
 - testing

Example serial code

```
for (i=0; i<N; i++)  
    for (j=0; j<N; j++)  
        for (k=0; k<N; k++)  
            c[i][j] += a[i][k] * b[k][j];
```

Code after tiling (via pluto)

```
for ( t2=0; t2 <= floord(N-1,32); t2++ )  
  for ( t3=0; t3<=floord(N-1,32); t3++ )  
    for ( t4=0; t4<=floord(N-1,32); t4++ )  
      for ( t5=32*t2; t5<=min(N-1,32*t2+31); t5++ )  
        for ( t6=32*t3; t6<=min(N-1,32*t3+31); t6++ )  
          for ( t7=32*t4; t7<=min(N-1,32*t4+31); t7++ )  
            c[t5][t6] += a[t5][t7] * b[t7][t6];
```

Code after tiling (via pluto)

for (t2=0; t2 <= floord(N-1,32); t2++)
 for (t3=0; t3<=floord(N-1,32); t3++)
 for (t4=0; t4<=floord(N-1,32); t4++)
 for (t5=32*t2; t5<=min(N-1,32*t2+31); t5++)
 for (t6=32*t3; t6<=min(N-1,32*t3+31); t6++)
 for (t7=32*t4; t7<=min(N-1,32*t4+31); t7++)
 c[t5][t6] += a[t5][t7] * b[t7][t6];

Tile

Tiled based kernel

```
for ( t2=0; t2 <= floord(N-1,32); t2++ )  
    for ( t3=0; t3<=floord(N-1,32); t3++ )  
        for ( t4=0; t4<=floord(N-1,32); t4++ )  
            zmm(a, b, c, t2, t3, t4, N);
```


Updating files

- Code of zmm() goes in dplasma/cores/core_zmm.c
- Prototype in dplasma/cores/dplasma_zcores.h
- dplasma/cores/CMakeLists.txt needs updating:

```
set(SOURCES
    core_ztrdv.c
    core_zmm.c
    core_zhetrf2_nopiv.c
    core_zhebut.c
    ${GPU_KERNEL_SOURCES}
)
```

Calling function: zgemm

```
int zgemm(PLASMA_desc a, PLASMA_desc b,  
          PLASMA_desc c, int N_sz, int ub )  
{  
    for (t2=0;t2<=ub;t2++)  
        for (t3=0;t3<=ub;t3++)  
            for (t4=0;t4<=ub;t4++)  
                Insert_Task(zmm);  
}
```

Calling function: zgemm

```
int zgemm(PLASMA_desc a, PLASMA_desc b,  
          PLASMA_desc c, int N_sz, int ub )  
{  
    for (t2=0;t2<=ub;t2++)  
        for (t3=0;t3<=ub;t3++)  
            for (t4=0;t4<=ub;t4++)  
                Insert_Task(zmm);  
}
```

This code resides in
a temporary file, say
zgemm.c

Calling function: zgemm

```
int zgemm(PLASMA_desc a, PLASMA_desc b,  
          PLASMA_desc c, int N_sz, int ub )
```

```
{  
    for (t2=0;t2<=ub;t2++)  
        for (t3=0;t3<=ub;t3++)  
            for (t4=0;t4<=ub;t4++)  
                Insert_Task(zmm);  
}
```

This code resides in
a temporary file, say
zgemm.c

Insert_Task() Quark Interface

```
QUARK_Insert_Task(junk, CORE_zmm_quark, junk,  
                  junk, a[t2][t4], INPUT,  
                  junk, b[t4][t3], INPUT,  
                  junk, c[t2][t3], INOUT,  
                  junk, &(t2),      VALUE,  
                  junk, &(t3),      VALUE,  
                  junk, &(t4),      VALUE,  
                  junk, &(N_sz),    VALUE,  
                  0);
```

Insert_Task() Generic Interface

```
Insert_Task( zmm,  
            a[t2][t4], INPUT,  
            b[t4][t3], INPUT,  
            c[t2][t3], INOUT,  
            t2,      VALUE,  
            t3,      VALUE,  
            t4,      VALUE,  
            N_sz,    VALUE);
```

New Quark to Old Quark

For Dyn. Sc. Quark codes, use Mathieu's "plasma2parsec.sh" & "core.h" (on zgemm.c)

Example from core.h:

```
#define QUARK_CORE_zsssm(quark, task_flags, m1, n1, m2, n2, k, ib, nb, A1, lda1, A2, lda2, L1, ld11, L2, ld12, IPIV) {\n    QUARK_Insert_Task((quark), CORE_zsssm_quark, (task_flags),\n        sizeof(int),                &(m1),    VALUE,\n        sizeof(int),                &(n1),    VALUE,\n        sizeof(int),                &(m2),    VALUE,\n        sizeof(int),                &(n2),    VALUE,\n        sizeof(int),                &(k),     VALUE,\n        sizeof(int),                &(ib),    VALUE,\n        sizeof(PLASMA_Complex64_t)*nb*nb, (A1),    INOUT,\n        sizeof(int),                &(lda1),   VALUE,\n        sizeof(PLASMA_Complex64_t)*nb*nb, (A2),    INOUT | LOCALITY,\n        sizeof(int),                &(lda2),   VALUE,\n        sizeof(PLASMA_Complex64_t)*ib*nb, (L1),    INPUT,\n        sizeof(int),                &(ld11),   VALUE,\n        sizeof(PLASMA_Complex64_t)*ib*nb, (L2),    INPUT,\n        sizeof(int),                &(ld12),   VALUE,\n        sizeof(int)*nb,              (IPIV),    INPUT,\n    0);}
```



Customization

- You can name the matrix variables

```
#pragma zsssm A1 A2 L1 L2 IPIV
```

- You can provide additional conditions

```
#pragma PARSEC_INVARIANT B.mt==minMT
```

- Invariant expression may contain:

```
||, &&, ==, !=, <=, <, (, ), +, -, vars, ints
```


JDF Generation

- `tools/q2j/q2j -anti zgemm.c > zgemm.jdf`
- q2j parameters:
 - `-shmem`
 - `-phony_tasks`
 - `-line_numbers`
 - `-anti`
 - `-v`
 - `file_1.c [file_2.c ... file_N.c]`

JDF Generation

- `tools/q2j/q2j -anti zgemv.c > zgemv.jdf`
- q2j parameters:
 - **-shm** (controls read/write pseudotasks)
 - **-phony_tasks** (controls phony IN/OUT tasks)
 - **-line_numbers** (adds #line in body)
 - **-anti** (finalizes antidependencies – can be very slow)
 - **-v** (more verbose warnings, slightly more)
 - **file_1.c [file_2.c ... file_N.c]** (input file(s))

.q2jrc

- generate_line_numbers = 0
- produce_shmem_jdf = 0
- finalize_antideps = 0
- verbose_warnings = 0
- add_phony_tasks = 0
- direct_output = 0

Q2J limitations

- Codes must be affine*
- Loops must be somewhat canonical
 - (i++, not i+=7)
 - Induction variables only in loop headers
 - No `for(i=0...){ j++; A[i][j] =... }`
- General format must look like PLASMA/Quark codes
- Codes must be self-contained*

Not self-contained codes

```
PLASMA_zgelqs_Tile_Async(){  
    plasma_parallel_call_3(plasma_pztile_zero,  
        PLASMA_desc, plasma_desc_submatrix(descB, descA.m, 0, descA.n-descA.m, descB.n),  
        PLASMA_sequence*, sequence,  
        PLASMA_request*, request);  
  
    plasma_parallel_call_9(plasma_pztrsm,  
        PLASMA_enum, PlasmaLeft,  
        PLASMA_enum, PlasmaLower,  
        PLASMA_enum, PlasmaNoTrans,  
        PLASMA_enum, PlasmaNonUnit,  
        PLASMA_Complex64_t, 1.0,  
        PLASMA_desc, plasma_desc_submatrix(descA, 0, 0, descA.m, descA.m),  
        PLASMA_desc, plasma_desc_submatrix(descB, 0, 0, descA.m, descB.n),  
        PLASMA_sequence*, sequence,  
        PLASMA_request*, request);  
}
```

plasma_desc_submatrix()

PLASMA_desc plasma_desc_submatrix(PLASMA_desc descA, int i, int j, int m, int n)

```
{  
    PLASMA_desc descB;  
    int mb, nb;  
  
    descB = descA;  
    mb = descA.mb;  
    nb = descA.nb;  
    // Submatrix parameters  
    descB.i = i;  
    descB.j = j;  
    descB.m = m;  
    descB.n = n;  
    // Submatrix derived parameters  
    descB.mt = (i+m-1)/mb - i/mb + 1;  
    descB.nt = (j+n-1)/nb - j/nb + 1;  
    return descB;  
}
```

plasma_desc_submatrix()

plasma_desc_submatrix(descB, descA.m, 0, descA.n-descA.m, descB.n)

PLASMA_desc plasma_desc_submatrix(PLASMA_desc descA, int i, int j, int m, int n)

```
{
    PLASMA_desc descB;
    int mb, nb;

    descB = descA;
    mb = descA.mb;
    nb = descA.nb;
    // Submatrix parameters
    descB.i = i;
    descB.j = j;
    descB.m = m;
    descB.n = n;
    // Submatrix derived parameters
    descB.mt = (i+m-1)/mb - i/mb + 1;
    descB.nt = (j+n-1)/nb - j/nb + 1;
    return descB;
}
```

plasma_desc_submatrix()

plasma_desc_submatrix(descB, descA.m, 0, descA.n-descA.m, descB.n)

PLASMA_desc plasma_desc_submatrix(PLASMA_desc descA, int i, int j, int m, int n)

```
{  
    descB = descA;  
    mb = descA.mb;  
    nb = descA.nb;  
    descB.i = descA.m;  
    descB.j = 0;  
    descB.m = descA.n-descA.m;  
    descB.n = descB.n;  
    descB.mt = (descA.m+descA.n-descA.m-1)/descA.mb - descA.m/descA.mb + 1;  
    descB.nt = (0+descB.n-1)/descA.nb - 0/descA.nb + 1;  
}
```


plasma_desc_submatrix()

plasma_desc_submatrix(descB, descA.m, 0, descA.n-descA.m, descB.n)

```
PLASMA_desc plasma_desc_submatrix(PLASMA_desc descA, int i, int j, int m, int n)
{
    descB = descA;
    mb = descA.mb;
    nb = descA.nb;
    descB.i = descA.m;
    descB.j = 0;
    descB.m = descA.n-descA.m;
    descB.n = descB.n;
    descB.mt = (descA.n-1)/descA.mb - descA.m/descA.mb + 1;
    descB.nt = (descB.n-1)/descA.nb + 1;
}
```

Proposed extension

```
plasma_desc_submatrix_affine(PLASMA_desc descA,  
                               int i,  
                               int j,  
                               int m,  
                               int n,  
                               int mt  
                               int nt  
                               );
```

Backend compilation daguepp

- JDF compiler: tools/dague-compiler/daguepp
- daguepp -o zgemm < zgemm.jdf
- daguepp will generate zgemm.c and zgemm.h
- .jdf, .c, and .h files should go in dplasma/lib
- We will also place the wrapper in dplasma/lib

Generated zgemm.h file

[illegible]

zgemm_wrapper.c NB interface

```
#include "zgemm.h"
dague_object_t *dplasma_zgemm_New( tiled_matrix_desc_t * desca,
                                    tiled_matrix_desc_t * descb,
                                    tiled_matrix_desc_t * descc,
                                    int N, int ub )
{
    dague_zgemm_object_t *o = dague_zgemm_new((dague_ddesc_t*)desca,
                                                (dague_ddesc_t*)descb,
                                                (dague_ddesc_t*)descc,
                                                N, ub);

    dplasma_add2arena_tile( o->arenas[DAGUE_zgemm_DEFAULT_ARENA],
                           desca->mb*desca->nb*sizeof(double),
                           DAGUE_ARENA_ALIGNMENT_SSE,
                           MPI_DOUBLE_COMPLEX, desca->mb );
    return (dague_object_t*)o;
}
```

zgemm_wrapper.c NB interface

```
void dplasma_zgemm_Destruct(dague_object_t *object)
{
    dague_zgemm_object_t *o = (dague_zgemm_object_t *)object;
    dplasma_datatype_undefine_type(
        &(o->arenas[DAGUE_zgemm_DEFAULT_ARENA]->opaque_dtt) );
    DAGUE_INTERNAL_OBJECT_DESTRUCT(o);
}
```

zgemm_wrapper.c B interface

```
int dplasma_zgemm(dague_context_t *dague,  
                  tiled_matrix_desc_t * desca,  
                  tiled_matrix_desc_t * descb,  
                  tiled_matrix_desc_t * descc,  
                  int N, int ub {  
  
    dague_handle_t *dague_zgemm = NULL;  
    dague_zgemm = dplasma_zgemm_New(ddesca, descb, descc, N, ub);  
  
    if ( dague_zgemm != NULL ) {  
        dague_enqueue( dague, (dague_handle_t*)dague_zgemm);  
        dplasma_progress(dague);  
        dplasma_zgemm_Destruct( dague_zgemm );  
    }
```

Updating files

dplasma/lib/CMakeLists.txt

```
set(JDF  
  zprint.jdf  
  zgemm.jdf  
  zlansy.jdf  
)
```

```
set(SOURCES  
  zprint_wrapper.c  
  zgemm_wrapper.c  
  zunmqr_wrapper.c  
)
```


Updating files

Add wrapper prototypes in dplasma/include/dplasma_z.h

```
dague_object_t *dplasma_zgemm_New( tiled_matrix_desc_t * desca,  
                                     tiled_matrix_desc_t * descb,  
                                     tiled_matrix_desc_t * descc,  
                                     int N, int ub )  
  
void dplasma_zgemm_Destruct(dague_object_t *object)  
  
int dplasma_zgemm(dague_context_t *dague,  
                  tiled_matrix_desc_t * desca,  
                  tiled_matrix_desc_t * descb,  
                  tiled_matrix_desc_t * descc,  
                  int N, int ub );
```

Writing testing_zgemm.c

```
#include "common.h"
#include "data_dist/matrix/two_dim_rectangle_cyclic.h"
int main(int argc, char ** argv)
{
    dague_context_t* dague;
    int iparam[IPARAM_SIZEOF];
    int ub;

    iparam_default_gemm(iparam);
    iparam_default_ibnbmb(iparam, 0, 32, 32);

    /* Initialize DAGuE */
    dague = setup_dague(argc, argv, iparam);
    PASTE_CODE_IPARAM_LOCALS(iparam);
    ub = floord(N-1,32);
```

testing_zgemm.c cont.

```
PASTE_CODE_ALLOCATE_MATRIX(ddescA, 1,  
    two_dim_block_cyclic, (&ddescA, matrix_ComplexDouble, matrix_Tile,  
        nodes, cores, rank, MB, NB, M, N, 0, 0,  
        M, N, SMB, SNB, P));
```

```
dplasma_gemm(dague,  
    (tiled_matrix_desc_t*)&ddescA,  
    (tiled_matrix_desc_t*)&ddescB,  
    (tiled_matrix_desc_t*)&ddescC,  
    N,  
    ub );
```

```
dague_data_free(ddescA.mat);  
dague_ddesc_destroy((dague_ddesc_t*)&ddescA);  
cleanup_dague(dague, iparam);
```

Updating files

- Main (testing) program goes into dplasma/testing
- dplasma/testing/CMakeLists.txt needs updating:
set(TESTS
 testing_zgemm.c
- tools/precision_generator/subs.py needs updating:
('sgemm','dgemm','cgemm','zgemm')

Obligatory voodoo

- Precision generator looks for following header:

```
/*  
 * @precisions normal z -> s d c  
 */
```

Compile & run

- `cd dplasma/testing; make -j testing_dgemm`
- `mpirun -np 4 ./testing_dgemm -N 128 -t 32 -c 1 -p 4 -q 1`
- For help on default parameters:
`./testing_dgemm -h`

Summary

- Source code: affine loops with “tiled” kernels
 - Quark or Generic interface for declaring tasks
- `tools/q2j/q2j -anti zmyop.c > dplasma/lib/zmyop.jdf`
- `tools/dague-compiler/daguepp -o zmyop < zmyop.jdf`
- Write wrapper and `main()`
 - `dplasma/lib/zmyop_wrapper.c`
 - `dplasma/testing/tesing_zmyop.c`
- Update `*/CMakeLists.txt`, `dplasma_z.h`, `subs.py`