# OpenCL Day 2: Loop Reasoning, Optimizing Temporal and Spatial Locality

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#### Content of Day2 Lecture:

Dependence Analysis

Problem Statement, Dependency Definition
Dependency Matrix, Loop Parallelism and Interchange
Loop Distribution, Array Expansion/Privatization
Exercise: Putting it all together

Optimizing Spatial Locality (Coalesced Memory)
Motivating Example on A Contrived Program
Uncoalesced Matrix Transposition Kernel
Coding Exercise: Entirely-Coalesced Transposition
Program Optimized By Transposition + Coding Exercise
Fully Fused Program + Coding Exercise

Optimizing Temporal Locality
Naive Matrix-Matrix Multiplication
Block-Tiled Matrix Multiplication + Coding Exercise
Register-Tiled Matrix Multiplication + Coding Exercise

#### **Problem Statement**

### Three Loop Examples

Iterations are ordered *lexicographically*, in the order they occur in the sequential execution:  $\vec{k} = (i=2, j=4) < \vec{l} = (i=3, j=3)$ .

- Which of the three loop nests is amenable to parallelization?
- ► Loop interchange is one of the most simple and useful code transformations, e.g., used to enhance locality of reference, and even to "create" parallelism.
- ► In which loop nest is it safe to interchange the loops?

### **Definition of a Dependency**

#### Load-Store Classification of Dependencies

**Def. Loop Dependence:** There is a dependence from statement S1 to S2 in a loop nest *iff* there exist iterations  $\vec{k} \leq \vec{l}$  and an execution path from statement S1 to statement S2 such that:

- 1. S1 accesses memory location M on iteration  $\vec{k}$ , and
- 2. S2 accesses memory location M on iteration  $\vec{l}$ , and
- 3. one of these accesses is a write.

We say that S1 is the source and S2 is the sink of the dependence, because S1 executes before S2 in the sequential program execution. Dependence depicted with an arrow pointing from source to sink.

## **Definition of a Dependency**

Load-Store Classification of Dependencies

**Def. Loop Dependence:** There is a dependence from statement *S*1 to *S*2 in a loop nest *iff* there exist iterations  $\vec{k} \leq \vec{l}$  and an execution path from statement *S*1 to statement *S*2 such that:

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We say that S1 is the source and S2 is the sink of the dependence, because S1 executes before S2 in the sequential program execution. Dependence depicted with an arrow pointing from source to sink.

We are most interested in cross iteration dependencies, i.e.,  $\vec{k} < \vec{l}$ .

### **Loop-Nest Dependencies**

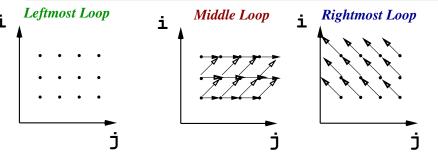
Lexicographic ordering, e.g.,  $\vec{k} = (i=2, j=4) < \vec{l} = (i=3, j=3)$ .

#### Three Loop Examples

## **Loop-Nest Dependencies**

Lexicographic ordering, e.g.,  $\vec{k} = (i=2, j=4) < \vec{l} = (i=3, j=3)$ .

#### Three Loop Examples



How can I summarize this information?

# **Aggregate Dependencies via Direction Vectors**

### Three Loop Examples

Dependencies depicted via an edge *from* the stmt that executes first in the loop nest (*source*), *to* the one that executes later (*sink*).

**Def. Dependence Direction:** Assume there exists a dependence from *S*1 in iteration  $\vec{k}$  to *S*2 in  $\vec{l}$  ( $\vec{k} \le \vec{l}$ ). The direction vector is:

- 1.  $\vec{D}(\vec{k}, \vec{l})_m =$  "<" if  $\vec{k}_m < \vec{l}_m$ , 2.  $\vec{D}(\vec{k}, \vec{l})_m =$  "=" if  $\vec{k}_m = \vec{l}_m$ ,
- 3.  $\vec{D}(\vec{k}, \vec{l})_m = ">" if <math>\vec{k}_m > \vec{l}_m$ .
- 4.  $\vec{D}(\vec{k}, \vec{l})_m =$  "\*" otherwise.

If the source is a write and the sink a read then RAW dependency, if the source is a read then WAR, if both are writes then WAW.

### **How to Compute the Direction Vectors?**

- ► For any two statements *S*1 and *S*2 that may access the same array *A* (and one of the accesses is a write),
- ▶ in two symbolic iterations  $l^1 \equiv (i_1^1, \dots i_m^1)$  and  $l^2 = (i_1^2, \dots i_m^2)$  (assuming  $l^1 \leq l^2$ )
- ▶ on indices  $A[e_1^1, \ldots, e_n^1]$  and  $A[e_1^2, \ldots, e_n^2]$ , respectivelly,
- ► then the direction vectors may be derived from the equations

$$\left\{egin{array}{l} e_1^1=e_1^2\ \cdots\ e_n^1=e_n^2 \end{array}
ight.$$

(The system of equations models the definition of a dependency: both accesses need to refer to the same memory location!)

```
Direction Vectors/Matrix for Three Loops

DO i = 1, N
```

DO j = 1, N S1 A[j,i]=A[j,i].. ENDDO ENDDO

### Direction Vectors/Matrix for Three Loops

Direction matrix:  $S1 \rightarrow S1$ :  $\lceil = , = \rceil$ 

 $S1 \rightarrow S1$ :  $\lceil =, = \rceil$ 

### Direction Vectors/Matrix for Three Loops

```
D0 i = 2, N
 DO i = 1, N
                                         D0 i = 2, N
ENDDO
                      FNDDO
                                            ENDDO
 ENDDO
                     ENDDO
                   S1\rightarrowS1: (j1,i1)=(j2-1,i2) For S1\rightarrowS1:
For S1 \rightarrow S1:
   (j1,i1)=(j2,i2)
                         i1 = i2 \& j1 < j2
   i1 = i2 \& j1 = j2
                   S2 \rightarrow S2: (j1,i1)=(j2-1,i2-1)
                        i1 < i2 & j1 < j2
Direction matrix:
                   S1→S1: [=,<]
```

S2→S2: 「<,<7

### Direction Vectors/Matrix for Three Loops

```
D0 i = 2, N
  DO i = 1, N
                                                         D0 i = 2, N
                             D0 i = 2, N
DO j = 1, N
S1 A[j,i]=A[j,i]. S1 A[j,i]=A[j-1,i]. S1 A[i,j]=A[i-1,j+1].
                                                           ENDDO
    ENDDO
                              ENDDO
                                                          ENDDO
  ENDDO
                            ENDDO
                         S1 \rightarrow S1: (j1,i1)=(j2-1,i2) For S1 \rightarrow S1:
For S1 \rightarrow S1:
                                   i1 = i2 & j1 < j2 (i1,j1) = (i2-1,j2+1)
    (j1,i1)=(j2,i2)
    i1 = i2 & j1 = j2 \\ S2 \rightarrow S2: (j1, i1) = (j2-1, i2-1)
                                                          i1 < i2 & j1 > j2
                                   i1 < i2 \& j1 < j^2Direction matrix:
Direction matrix:
                          S1→S1: [=,<]
S1 \rightarrow S1: \lceil =, = \rceil
                                                      S1→S1: [<,>]
                          S2→S2: 「<,<7
```

**Th. Parallelism:** A loop in a loop nest is parallel *iff* for each entry in its direction column, the entry is either = or there exists an outer loop whose corresponding direction is <.

A direction vector cannot have > as the first non-= symbol, as that would mean that I depend on something in the future.

## Parallelism and Loop Interchange

```
Direction Vectors/Matrix for Three Loops
                            D0 i = 2, N
  DO i = 1, N
                                                          D0 i = 2, N
                            D0 i = 2, N
    DO j = 1, N
                                                         D0 j = 1, N
S1 A[j,i]=A[j-1,i]... S1 A[j,j]=A[j-1,i]... S1 A[i,j]=... S1 A[i,j]=... S1 A[i,j]=...
    ENDDO
                                                                AΓi-1,j+17
                              ENDDO
                                                          ENDDO ENDDO
  ENDDO
                            ENDDO
For S1 \rightarrow S1:
                                                         For S1 \rightarrow S1:
                         S1 \rightarrow S1: (j1,i1)=(j2-1,i2)
    (j1,i1)=(j2,i2)
                                                            (i1,j1) = (i1-1,j2+1)
                                  i1 = i2 \& j1 < j2
    i1 = i2 \& j1 = j2
                                                           i1 < i2 & j1 > j2
                         S2 \rightarrow S2: (j1,i1)=(j2-1,i2-1)
                                   i1 < i2 & j1 < j2
                                                         Direction matrix:
Direction matrix:
                         S1→S1: [=,<]
S1 \rightarrow S1: \lceil =, = \rceil
                                                         S1→S1: [<,>]
                         S2→S2: 「<,<7
```

**Th. Loop Interchange:** Interchanging two loops in a loop nest is legal *iff* interchanging the columns of the direction matrix in the same way *does NOT result* in a > direction as the leftmost non-= direction in a row.

## Parallelism and Loop Interchange

```
Direction Vectors/Matrix for Three Loops
 DO i = 1. N
S1 A[j,i]=A[j,i]... S1 A[j,i]=A[j-1,i]... S1 A[i,j]=A[i-1,j+1]...
   ENDDO
                     S2 B\lceil j, i \rceil = B\lceil j-1, i-1 \rceil \dots ENDDO
                    ENDDO ENDDO
 ENDDO
                                            ENDDO
                     For S1\rightarrowS1: j1 = j2-1 For S1\rightarrowS1: i1 = i2-1
For S1 \rightarrow S1: j1 = j2
           i1 = i2
                             i1 = i2 j1 = j2+1
                    (i2,j2)-(i1,j1)=[=,<] (i2,j2)-(i1,j1)=
(i2,j2)-(i1,j1)=
                   For S2 \rightarrow S2: j1 = j2-1 [<,>]
\Gamma = , = 1
                                i1 = i2-1
```

(i2,j2)-(i1,j1)=[<,<]

Interchange is safe for the 1st and 2nd nests, but not for the 3rd! e.g., [=,<]  $\rightarrow$  [<,=] (for the second loop nest) [<,<]

## Parallelism and Loop Interchange

```
Direction Vectors/Matrix for Three Loops
 DO i = 2, N
                                    DO j = 1, N
S1 A[j,i]=A[j,i]... S1 A[j,i]=A[j-1,i]... S1 A[i,j]=A[i-1,j+1]...
    ENDDO
                      S2 B\lceil j, i \rceil = B\lceil j-1, i-1 \rceil \dots ENDDO
                     ENDDO ENDDO
 ENDDO
                                              ENDDO
                      For S1 \rightarrow S1: j1 = j2-1 For S1 \rightarrow S1: i1 = i2-1
For S1 \rightarrow S1: j1 = j2
           i1 = i2
                                i1 = i2 j1 = j2+1
                     (i2,j2)-(i1,j1)=[=,<] (i2,j2)-(i1,j1)=
(i2,j2)-(i1,j1)=
                      For S2 \rightarrow S2: j1 = j2-1 [<,>]
\Gamma = , = 1
```

(i2, j2)-(i1, j1)=[<,<]

i1 = i2-1

```
Interchange is safe for the 1st and 2nd nests, but not for the 3rd! e.g., [=,<] \rightarrow [<,=] (for the second loop nest) [<,<]
```

After interchange, loop *j* of the second loop nest is parallel.

Corollary: A parallel loop can be always interchanged inwards.

## **Dependency Graph and Loop Distribution**

**Def. Dependency Graph:** edges from the source of the dependency, i.e., early iteration, to the sink, i.e., later iteration.

**Th. Loop Distribution:** Statements that are in a dependence cycle remain in one (sequential) loop. The others are distributed to separate loops in graph order; if no cycle then parallel loops.

#### Vectorization Example

```
DO i = 3, N

S1 A[i] = B[i-2] ...

S2 B[i] = B[i-1] ...

ENDDO

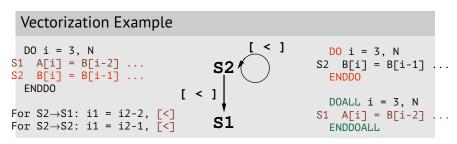
For S2→S1: i1 = i2-2, [<]

For S2→S2: i1 = i2-1, [<]
```

## **Dependency Graph and Loop Distribution**

**Def. Dependency Graph:** edges from the source of the dependency, i.e., early iteration, to the sink, i.e., later iteration.

**Th. Loop Distribution:** Statements that are in a dependence cycle remain in one (sequential) loop. The others are distributed to separate loops in graph order; if no cycle then parallel loops.



**Corollary:** It is always legal to distribute a parallel loop; but requires array expansion for local variables or if output dependencies are present.

## **Loop Distribution May Require Array Expansion**

```
float tmp;
float tmp;
for(i=2; i<N; i++) {
    tmp = 2*B[i-2];
    A[i] = tmp;
    B[i] = tmp+B[i-1]
}
float tmp[N];
for(int i=2; i<N; i++) {
    tmp[i] = 2*B[i-2];
    B[i] = tmp[i]+B[i-1];
}
forall(int i=2; i<N; i++) {
    A[i] = tmp[i];
}</pre>
```

No matter where tmp is declared (inside or outside the loop) it needs to be expanded into an array in order to do loop distribution.

If tmp is declared outside the loop then requires **privatization**,

## **Loop Distribution May Require Array Expansion**

```
float tmp;
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for(i=2; i<N; i++) {
    tmp = 2*B[i-2];
    A[i] = tmp;
    B[i] = tmp+B[i-1]
}
float tmp[N];
for(int i=2; i<N; i++) {
    tmp[i] = 2*B[i-2];
    B[i] = tmp[i]+B[i-1];
}
forall(int i=2; i<N; i++) {
    A[i] = tmp[i];
}</pre>
```

No matter where tmp is declared (inside or outside the loop) it needs to be expanded into an array in order to do loop distribution.

If tmp is declared outside the loop then requires **privatization**, because it actually causes frequent waw dependencies. However its value is written before being used within the same iteration. Hence it is semantically equivalent to a locally declared variable, which will remove the output (waw) dependency.

Distribution requires array expansion of the scalar tmp.

## False Dependencies (WAR/WAW)

- ► Cross-Iteration Anti Dependencies (WAR) correspond to a read from the array as it was before the loop ⇒ can be eliminated by reading from a copy of the array.
- ► Cross-Iteration WAW Dependencies (WAW): If they correspond to the case in which every read from a scalar or array location is covered by a previous same-iteration write ⇒ can be eliminated privatization (renaming), which semantically moves the declaration of the variable (scalar or array) inside the loop.
- ► Direction-vectors reasoning is limited to relatively simple loop nests, e.g., difficult to reason about privatization in such a way.

### **Example: Eliminating WAW Dependencies**

```
float A[M];
for(i=0; i<N; i++){
  for(int j=0, j<M; j++)
        A[j] = (4*i+4*j) % M;
  for(int k=0; k<N; k++)
        X[i][k] = X[i][k-1] *
        A[A[(2*i+k)%M]%M];
}</pre>
```

### **Example: Eliminating WAW Dependencies**

```
// The write to A[j] causes many WAWs,

float A[M];
for(i=0; i<N; i++){
   for(int j=0, j<M; j++)
        A[j] = (4*i+4*j) % M;
   for(int k=0; k<N; k++)
        X[i][k] = X[i][k-1] *
        A[A[(2*i+k))M],M];
```

### **Example: Eliminating WAW Dependencies**

```
// The write to A[j] causes many WAWs,
                                 // but A is fully written in loop j
float AFM7;
                                 // float AFN7FM7;
for(i=0; i<N; i++){</pre>
                                 forall(int i=0; i<N; i++){</pre>
                                  float AFM7;
  for(int j=0, j<M; j++)
      A[j] = (4*i+4*j) \% M; for(int j=0, j<M; j++)
  for(int k=0; k<N; k++)
                                       A[j] = (4*i+4*j) % M; // A[i][j]
      X[i][k] = X[i][k-1] *
                                  for(int k=0; k<N; k++)
         A\Gamma A\Gamma (2*i+k)%M7%M7;
                                       X[i][k]=X[i][k-1] *
                                          A\Gamma A\Gamma (2*i+k)%M7%M7;
                                          //AΓi][AΓi,...]]
```

For OpenCL programming you would likely use array expansion (float A[N, M]) because the size M is likely unknown at compile time.

# Putting Everything Together

```
float X[M][N][N]; /* ... compute X ... */
float A[N][N];
for(int i = 0; i < M; i++) {
  //initialize A
  for(int k = 0; k < N; k++) {
    for(int j = 0; j < N; j++) {
        float x = X[i][j][k];
        A[k][j] = x * x;
  // convergence loop
  for(int t = 0; t < T; t++) {
    for(int j = 0; j < N; j++) {
      for(int k = 0; k < N; k++) {
        A[k][j] = fsqrt(A[k][j] * A[k-1][j]);
        X[i][k][j] += fsqrt(B[i][k-1][j-1]);
```

**Exercise:** Assume the code above runs on a dataset having M = 128 and N = 512 and, less important, say T = 256.

- (1) Is this code suitable for GPU execution?
- (2) Transform the code to one suitable for GPU execution.

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#### Dependence Analysis

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## **Motivation: Coalesced Accesses to Global Memory**

Which are the parallel/sequential loops?

```
real A[height][width];
real B[height][width];
// Non-Coalesced Access
for(i=0; i<height; i++) {
  real accum = 0.0;

  for(j=0; j<width; j++) {
    real tmpA = A[i][j];
    accum = tmpA*tmpA - accum;
    B[i][j] = accum;
}
}</pre>
```

## **Motivation: Coalesced Accesses to Global Memory**

Which are the parallel/sequential loops?

```
real A[height][width];
                                 A' = transpose(A);
real B[height][width];
                                 // Coalesced Accesses
// Non-Coalesced Access
                                 for(i=0; i<height; i++) {</pre>
for(i=0; i<height; i++) {</pre>
                                   real accum = 0.0;
  real accum = 0.0:
                                   for(int j=0; j<width; j++) {</pre>
                                      real tmpA = A'[j][i];
  for(j=0; j<width; j++) {</pre>
    real tmpA = A[i][j];
                                      accum = tmpA*tmpA - accum;
    accum = tmpA*tmpA - accum;
                                      B'[i][i] = accum;
    B[i][j] = accum;
                                   = transpose(B');
```

- ► Loop i is parallel;
- ▶ loop j is sequential because of the RAW cross-iteration dependencies on accum.
- ► The transformed program performs about 3× the number of global-memory accesses than the original.

# **Motivation: Coalesced Accesses to Global Memory**

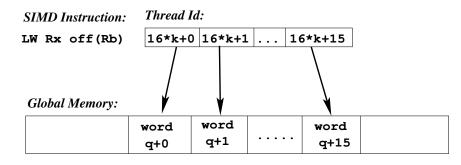
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for(i=0; i<height; i++) {</pre>
                                    real accum = 0.0;
  real accum = 0.0:
                                    for(int j=0; j<width; j++) {</pre>
                                      real tmpA = A'[j][i];
  for(j=0; j<width; j++) {</pre>
    real tmpA = A[i][j];
                                      accum = tmpA*tmpA - accum;
    accum = tmpA*tmpA - accum;
                                      B'[i][i] = accum;
    B[i][j] = accum;
                                   = transpose(B');
```

- ▶ Loop i is parallel;
- ▶ loop j is sequential because of the RAW cross-iteration dependencies on accum.
- ► The transformed program performs about 3× the number of global-memory accesses than the original.
- ▶ But it is significantly faster than the original (on GPUs).

### **Motivation: What Are Coalesced Accesses?**

**Coalesced access**: a (quarter) wave accesses in a SIMD instruction consecutive words in global-memory.



### **Uncoalesced Transposition Kernel**

- ► Semantically we chunk the matrix into T×T blocks;
- ► Each block is processed by one workgroup, e.g., T=16.

- ▶ Local workgroup:  $T \times T$ ;
- ► Global(1): [width/T] \* T; Global 1: [height/T] \* T.

### **Uncoalesced Transposition Kernel**

- ightharpoonup Semantically we chunk the matrix into TimesT blocks;
- ► Each block is processed by one workgroup, e.g., T=16.

```
// Initial
                               // Blocked
for(i=0; i< height; i++)  for(ii=0; ii< height; ii+=T)  {
 for(j=0; j<width; j++) {</pre>
                                 for(jj=0; jj<width; jj+=T) {</pre>
    B[j*height+i] =
                                   for(i=ii; i<min(ii+T,height); i++) {</pre>
           AΓi*width+il;
                                     for(j=jj; j<min(jj+T,width); j++) {</pre>
                                       B[j*height+i] = A[i*width+j];
                               } } } }
   ▶ Local workgroup: T \times T;
   ► Global(1): [width/T] * T; Global 1: [height/T] * T.
 kernel void naiveTransp( | global real* A, | global real* B
                            , int height, int width ) {
     unsigned int i = get global id(1);
     unsigned int j = get_global_id(0);
     if( (j >= width) || (i >= height) ) return;
     B[j*height + i] = A[i*width + j]; // uncoalesced access!
```

## **Exercise 1: Fill In Coalesced Transposition Kernel**

```
// verifies that workgroup size is: TILE*TILE
__kernel __attribute__((reqd_work_group_size(TILE, TILE, 1)))
kernel void coalsTransp( global real* A,
       global real* B, uint32 t height, uint32 t width ) {
  local float tile[TILE][TILE];
                                                    // local memory
  int gidx = get_global_id(0); int gidy = get_global_id(1);
  int lidx = get local id(0); int lidy = get local id(1);
  // 1. read current element from global to local memory.
       make sure to NOT access arrays out of bounds!
  // 2. current thread loads from local memory the transposed element
       at workgroup level, and writes it in the proper position in B.
  // Useful functions: get group id(0), get local size(0); or 1.
  // 3. Insert necessary workgroup-level synchronization
       barrier(CLK LOCAL MEM FENCE);
```

Transposing a (blocked) matrix  $\equiv$  moving each block to the transposed position + transposing each block!

## **Exercise 1: Fill In Coalesced Transposition Kernel**

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 int gidx = get_global_id(0); int gidy = get_global_id(1);
 int lidx = get local id(0); int lidy = get local id(1);
 // 1. read current element from global to local memory.
       make sure to NOT access arrays out of bounds!
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       at workgroup level, and writes it in the proper position in B.
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 // 3. Insert necessary workgroup-level synchronization
       barrier(CLK LOCAL MEM FENCE);
```

Transposing a (blocked) matrix  $\equiv$  moving each block to the transposed position + transposing each block!

- ▶ Implement in Day2-Exercises/Transp/kernels.cl
- Are the read and write accesses to global memory coalesced?
- ► Why/Where should you place the barrier?

# **Exercise 2: Coalescing By Transposition**

```
real A[height][width];
real B[height][width];
// Non-Coalesced Access
for(i=0; i<height; i++) {
  real accum = 0.0;

  for(j=0; j<width; j++) {
    real tmpA = A[i][j];
    accum = tmpA*tmpA - accum;
    B[i][j] = accum;
} }</pre>
```

# **Exercise 2: Coalescing By Transposition**

```
real A[height][width];
                                 A' = transpose(A);
real B[height][width];
                                 // Coalesced Accesses
// Non-Coalesced Access
                                 for(i=0; i<height; i++) {</pre>
for(i=0; i<height; i++) {</pre>
                                    real accum = 0.0:
  real accum = 0.0;
                                    for(int j=0; j<width; j++) {</pre>
  for(j=0; j<width; j++) {</pre>
                                      real tmpA = A'[j][i];
    real tmpA = A[i][j];
                                      accum = tmpA*tmpA - accum;
    accum = tmpA*tmpA - accum;
                                      B'[j][i] = accum;
    B[i][j] = accum;
} }
                                      transpose(B');
```

- ► Open File "Day2-Exercises/Transp/coalescing.c", function named "runGPUcoalsProgram".
- ► Fill the correct sizes for the "coalsProgrm" kernel call.
- ► Open File "Day2-Exercises/Transp/kernels.cl" and implement kernel named "coalsProgrm".
- ► Fix bugs until it validates! :-)
- ► What is the speedup with respect to the naive program?
- ▶ What fraction of the memcpy bandwidth is achieved?

#### **Exercise 3: Fused Program With Coalesced Accesses**

- ▶ Open File "Day2-Exercises/Transp/kernels.cl" and implement the missing pieces from kernel named "optimProgrm";
- ► The idea is to have coalesced accesses "fused" into the kernel by each thread processing CHUNK elements at a time by:
  - collectively copying CHUNK \* WORKGROUP\_SIZE input elements from global to local memory in coalesced fashion;
  - each thread now processes CHUNK elements from its row;
  - collectively copying CHUNK \* WORKGROUP\_SIZE result elements from local to global memory in coalesced fashion;
  - ► repeat in a loop (inside the kernel) until all row's elements have been processed!
- ► Fix bugs until it validates! :-)
- ► What is the speedup with respect to the transposition-optimized program?
- ▶ What fraction of the memcpy bandwidth is achieved?

#### Content of Day2 Lecture:

#### Dependence Analysis

Problem Statement, Dependency Definition
Dependency Matrix, Loop Parallelism and Interchange
Loop Distribution, Array Expansion/Privatization
Exercise: Putting it all together

Optimizing Spatial Locality (Coalesced Memory)

Motivating Example on A Contrived Program

Uncoalesced Matrix Transposition Kernel

Coding Exercise: Entirely-Coalesced Transposition

Program Optimized By Transposition + Coding Exercise

Fully Fused Program + Coding Exercise

### Optimizing Temporal Locality

Naive Matrix-Matrix Multiplication
Block-Tiled Matrix Multiplication + Coding Exercise
Register-Tiled Matrix Multiplication + Coding Exercise

## Matrix Multiplication: Loop Strip Mining

Accesses to A and B invariant to loops i and  $j \Rightarrow$ The idea is to apply Block Tiling to optimize temporal locality!

#### **Loop Strip Mining**

Strip Mining is always safe: the transformed loop executes the same instructions in the same order as the original loop!

Strip mining all loops by some tile T result in:

### **Loop Strip Mining**

float tmp = 0.0;

```
for(int ii = 0; ii < N; ii += T) {
\Rightarrow for(int i = ii < N; ii += T) 
for(int i = 0; i < N; i++) {
    loop_body(i)
                                  for(int i = ii; i < MIN(ii+T,N); i++) {
                                     loop_body(i)
 Strip Mining is always safe: the transformed loop executes the
 same instructions in the same order as the original loop!
 Strip mining all loops by some tile T result in:
                                         // Parallel
 for (int ii = 0; ii < M; ii += T) {
    for (int i = ii; i < MIN(ii+T,M); i++) { // Parallel
      for (int jj = 0; jj < N; jj += T) { // Parallel
        for (int j = jj; j < MIN(jj+T,N); j++) { // Parallel
```

for(int kk = 0; kk < U; k += T) { // Reduction
for(int k=kk; k<MIN(kk+T,U); k++) {// Reduction</pre>

tmp += A[i,k]\*B[k,j];

#### Block Tiling: Strip Mining + Interchange

Now interchange the loops if index i and jj. Why is this safe?

```
for (int ii = 0; ii < M; ii += T) {
                                           // Global 1
  for (int jj = 0; jj < N; jj += T) {
                                             // Global 0
    for (int i = ii; i < MIN(ii+T,M); i++) { // Local 1
     for (int j = jj; j < MIN(jj+T,N); j++) { // Local 0
         float tmp = 0.0:
         for(int kk = 0; kk < U; k += T) { // Sequential
           //What slices of A and B are used in the loop below?
           for(int k=kk; k<MIN(kk+T,U); k++) { // Sequential</pre>
               tmp += A[i,k]*B[k,j];
         C[i,j] = tmp;
} } } }
```

#### **Block Tiling: Strip Mining + Interchange**

Now interchange the loops if index i and jj. Why is this safe?

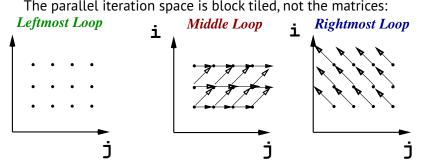
```
for (int ii = 0; ii < M; ii += T) {
                                               // Global 1
  for (int jj = 0; jj < N; jj += T) {
                                             // Global 0
    for (int i = ii; i < MIN(ii+T,M); i++) { // Local 1
      for (int j = jj; j < MIN(jj+T,N); j++) { // Local 0
          float tmp = 0.0:
          for(int kk = 0; kk < U; k += T) { // Sequential
            //What slices of A and B are used in the loop below?
            for(int k=kk; k<MIN(kk+T,U); k++) { // Sequential</pre>
                tmp += A[i,k]*B[k,j];
          C[i,j] = tmp;
} } } }
A[ii : MIN(ii+T,M)][kk : MIN(kk+T,U)]
B\lceil kk : MIN(kk+T,U)\rceil\lceil jj : MIN(jj+T,N)\rceil
```

Each array slice has  $T \times T$  elements and there are  $T \times T$  threads in a workgroup; so each thread can copy one element to local memory. Thus, we cut the number of global-memory accesses by a T factor.

- ➤ Open File "Day2-Exercises/MMMult/mmm.c", function named "runGPUverMMM".
- ► Place there the correct sizes for the matrix multiplication kernels. That would make the naive version work correctly.
- ▶ Open File "Day2-Exercises/MMMult/mmm.cl" and implement kernel named "blockMMM".
- Fix bugs until it validates! :-)

- ▶ Open File "Day2-Exercises/MMMult/mmm.c", function named "runGPUverMMM".
- Place there the correct sizes for the matrix multiplication kernels. That would make the naive version work correctly.
- ▶ Open File "Day2-Exercises/MMMult/mmm.cl" and implement kernel named "blockMMM".
- Fix bugs until it validates! :-)

The parallel iteration space is block tiled, not the matrices:



```
Copy A[ii:MIN(ii+T,M)][kk:MIN(kk+T,U)] in A_loc of size [T][T]
Copy B[kk:MIN(kk+T,U)][jj:MIN(jj+T,N)] in B_loc of size [T][T]
for (int ii = 0; ii < M; ii += T) {
                                                 // Global 1
  for (int jj = 0; jj < N; jj += T) {
                                                 // Global 0
    for (int i = ii; i < MIN(ii+T,M); i++) { // Local 1
      for (int j = jj; j < MIN(jj+T,N); j++) { // Local 0
          float tmp = 0.0:
          for(int kk = 0; kk < U; k += T) { // Sequential
            // Collectively copy to local memory.
             // Loop of index k should only use local memory.
             // Add barrier here. Why?
             for(int k=kk; k<MIN(kk+T,U); k++) { // Sequential</pre>
                 tmp += A loc[i-ii,k-kk] * B loc[k-kk,j-jj];
          } } // Add barrier here. Why?
          C[i,j] = tmp;
} } } }
```

Modify loops of index i, j and k to go from 0 to T with stride 1. Unroll loop k.

OpenCL equivalences: Global workgroup size:

```
Copy A[ii:MIN(ii+T, M)][kk:MIN(kk+T, U)] in A_loc of size [T][T]
Copy B[kk:MIN(kk+T,U)][jj:MIN(jj+T,N)] in B_loc of size [T][T]
for (int ii = 0; ii < M; ii += T) {
                                                // Global 1
  for (int jj = 0; jj < N; jj += T) {
                                                 // Global 0
    for (int i = ii; i < MIN(ii+T,M); i++) { // Local 1
      for (int j = jj; j < MIN(jj+T,N); j++) { // Local 0
          float tmp = 0.0:
          for(int kk = 0; kk < U; k += T) { // Sequential
            // Collectively copy to local memory.
             // Loop of index k should only use local memory.
             // Add barrier here. Why?
             for(int k=kk; k<MIN(kk+T,U); k++) { // Sequential</pre>
                 tmp += A loc[i-ii,k-kk] * B loc[k-kk,j-jj];
          } } // Add barrier here. Why?
          C[i,j] = tmp;
} } } }
```

Modify loops of index i, j and k to go from 0 to T with stride 1. Unroll loop k.

OpenCL equivalences:

Global workgroup size:  $(\lceil M/T \rceil * T) \times (\lceil N/T \rceil * T)$ . Local Workgroup size:  $T \times T$ .

```
Copy A[ii:MIN(ii+T,M)][kk:MIN(kk+T,U)] in A_loc of size [T][T]
Copy B[kk:MIN(kk+T,U)][jj:MIN(jj+T,N)] in B_loc of size [T][T]
for (int ii = 0; ii < M; ii += T) {
                                                      // Global 1
  for (int jj = 0; jj < N; jj += T) {
                                              // Global 0
     for (int i = ii; i < MIN(ii+T,M); i++) { // Local 1
       for (int j = jj; j < MIN(jj+T,N); j++) { // Local 0
            float tmp = 0.0:
            for(int kk = 0; kk < U; k += T) { // Sequential
              // Collectively copy to local memory.
               // Loop of index k should only use local memory.
               // Add barrier here. Why?
               for(int k=kk; k<MIN(kk+T,U); k++) { // Sequential</pre>
                    tmp += A loc[i-ii,k-kk] * B loc[k-kk,j-jj];
            } } // Add barrier here. Why?
            C[i,j] = tmp;
} } } }
Modify loops of index i, i and k to go from 0 to T with stride 1. Unroll loop k.
OpenCL equivalences:
Global workgroup size: (\lceil M/T \rceil * T) \times (\lceil N/T \rceil * T). Local Workgroup size: T \times T.
Localids: i - ii = get_local_id(1), j - jj = get_local_id(0)
Globalids: i = get_global_id(1), j = get_global_id(0)
Similar: ii = get_group_id(1)*T, jj = get_group_id(0)*T
Local memory declared with: __local float A_loc[T][T];
```

Loops of index i and j form the local workgroup. Interchange them innermost and distribute them:

```
// Global 1
for (int ii = 0: ii < M: ii += T) {
 for (int jj = 0; jj < N; jj += T) { // Global 0
   float A_loc[T][T], B_loc[T][T]; // local memory
   float tmp regTTTTT: // register memory: still represented as a scalar for a thread
   for (int i = ii; i < ii+T; i++) { // Local 1
     for (int j = jj; j < jj+T; j++) { // Local 0
          tmp\Gamma i - ii\Gamma \Gamma i - ii\Gamma = 0.0:
    } }
    for(int kk = 0; kk < U; k += T) { // Sequential
     for (int i = ii; i < ii+T; i++) { // Local
        for (int j = jj; j < jj+T; j++) { // Local
          A_{loc[i-ii][k-kk]} = (i < M \& k < U)? A[i][kk+j-jj] : 0.0; // copy to local mem
          B \log[k-kk][j-jj] = (j < N \&\& k < U)? B[kk+i-ij][j] : 0.0; // copy to local mem
      } } // needs barrier here, why?
     for (int i = ii; i < ii+T; i++) { // Local 1
       for (int j = jj; j < jj+T; j++) { // Local 0
           for(int k=kk; k < kk+T; k++) { // Sequential
                tmp[i][j] += A loc[i-ii,k-kk] * B[k-kk,j-jj]; // use local memory
      } }
     for (int i = ii; i < ii+T; i++) { // Local 1
       for (int j = jj; j < jj+T; j++) { // Local 0
          C\Gamma i.i7 = tmp:
      } } // needs barrier here, why?
```

Modify loops of index i, i and k to go from 0 to T with stride 1. Unroll loop k.

OpenCL equivalences:

Global workgroup size:

```
Loops of index i and j form the local workgroup. Interchange them innermost and distribute them:
```

```
// Global 1
for (int ii = 0: ii < M: ii += T) {
 for (int jj = 0; jj < N; jj += T) { // Global 0
    float A_loc[T][T], B_loc[T][T]; // local memory
    float tmp regTTTTT: // register memory: still represented as a scalar for a thread
    for (int i = ii; i < ii+T; i++) { // Local 1
      for (int j = jj; j < jj+T; j++) { // Local 0
          tmp\Gamma i - ii\Gamma \Gamma i - ii\Gamma = 0.0:
    } }
    for(int kk = 0; kk < U; k += T) { // Sequential
      for (int i = ii: i < ii+T: i++) { // Local
        for (int j = jj; j < jj+T; j++) { // Local
          A_{loc[i-ii][k-kk]} = (i < M \& k < U)? A[i][kk+j-jj] : 0.0; // copy to local mem
          B \log[k-kk][j-jj] = (j < N \&\& k < U)? B[kk+i-ij][j] : 0.0; // copy to local mem
      } } // needs barrier here, why?
      for (int i = ii; i < ii+T; i++) { // Local 1
        for (int j = jj; j < jj+T; j++) { // Local 0
            for(int k=kk; k < kk+T; k++) { // Sequential
                tmp[i][j] += A loc[i-ii,k-kk] * B[k-kk,j-jj]; // use local memory
      } }
      for (int i = ii; i < ii+T; i++) { // Local 1
        for (int j = jj; j < jj+T; j++) { // Local 0
          C\Gamma i.i7 = tmp:
      } } // needs barrier here, why?
Modify loops of index i, i and k to go from 0 to T with stride 1. Unroll loop k.
```

OpenCL equivalences:

Global workgroup size:  $(\lceil M/T \rceil * T) \times (\lceil N/T \rceil * T)$ . Local Workgroup size:  $T \times T$ . Local ids:

Loops of index i and j form the local workgroup. Interchange them innermost and distribute them:

```
for (int ii = 0; ii < M; ii += T) { // Global 1
  for (int jj = 0; jj < N; jj += T) { // Global 0
    float A_loc[T][T], B_loc[T][T]; // local memory
    float tmp regTTTTT: // register memory: still represented as a scalar for a thread
    for (int i = ii; i < ii+T; i++) { // Local 1
      for (int j = jj; j < jj+T; j++) { // Local 0
          tmp\Gamma i - ii\Gamma \Gamma i - ii\Gamma = 0.0:
    } }
    for(int kk = 0; kk < U; k += T) { // Sequential
      for (int i = ii: i < ii+T: i++) { // Local
        for (int j = jj; j < jj+T; j++) { // Local
          A loc[i-ii][k-kk] = (i < M && k < U)? A[i][kk+j-jj] : 0.0; // copy to local mem
          B \left[ \log \left[ k - kk \right] \right] = (j < N \& k < U)? B \left[ kk + i - ii \right] \left[ j \right] : 0.0; // copy to local mem
      } } // needs barrier here, why?
      for (int i = ii; i < ii+T; i++) { // Local 1
        for (int j = jj; j < jj+T; j++) { // Local 0
            for(int k=kk: k < kk+T: k++) { // Sequential</pre>
                 tmp[i][j] += A loc[i-ii,k-kk] * B[k-kk,j-jj]; // use local memory
      } }
      for (int i = ii; i < ii+T; i++) { // Local 1
        for (int j = jj; j < jj+T; j++) { // Local 0
          C\Gamma i.i7 = tmp:
      } } // needs barrier here, why?
Modify loops of index i, i and k to go from 0 to T with stride 1. Unroll loop k.
```

OpenCL equivalences:

Global workgroup size:  $(\lceil M/T \rceil * T) \times (\lceil N/T \rceil * T)$ . Local Workgroup size:  $T \times T$ . Local ids: i - ii = get\_local\_id(1), j - jj = get\_local\_id(0) Global ids: i = get\_global\_id(1), j = get\_global\_id(0)

Local memory declared with: \_\_local float A\_loc[T][T];

```
for (int i = 0; i < M; i++) { // Parallel | Matrices: for (int j = 0; j < N; j++) { // Parallel | \blacktriangleright A \in \mathcal{M}^{M \times U} | float c = 0.0 | for(int k = 0; k < U; k++) { // Reduction | \blacktriangleright B \in \mathcal{M}^{U \times N} | \blacktriangleright C \in \mathcal{M}^{M \times N} | \blacktriangleright C \in \mathcal{M}^{M \times N} | Loops of indices i and j are parallel.
```

We can do even better by sacrificing some parallelism in excess:

- ► stripmine parallel dimension i by a tile T and stride 1; then move it innermost and sequentialize it;
- $\blacktriangleright$  stripmine parallel dimension j by a tile T $\times$ T and stride T and
- stripmine the result by a tile T and stride 1;
- ► the two tiles of parallel dimension j form the local workgroup.

Do not forget to perform array expansion on c, which we will rename cs.

```
unsigned int ii, i, jjj, jj, j, kk, k;
forall (ii = 0; ii < M; ii += T ) {
                                    // Parallel Global 1
 forall (jjj = 0; jjj < N; jjj += T*T ) { // Parallel Global 0
   float cs[T][T][T];
   forall(jj=jjj; jj<min(jjj+T*T,N); jj+=T){ // Parallel Local 1</pre>
     forall(j=jj; j<min(jj+T,N); j++) { // Parallel Local 0</pre>
       for (i=ii; i<min(ii+T,M); i++) { // sequential</pre>
         cs\Gamma(ii-iii)/T\Gamma\Gamma i-ii\Gamma\Gamma i-ii\Gamma = 0.0:
   } } }
   for (kk = 0; kk < U; kk += T) {
                                               // sequential
     // here we will insert a collective copy to local
     // memory of the slice: A[ii : ii+T, kk : kk+T]
     for (k = kk; k < min(kk+T,U); k++)  // sequential
       forall (jj=jjj; jj<min(jjj+T*T,N); jj+=T) { // Parallel Local 1
         // please modify here to read A from local memory
            // hoist out B[k][i] outside loop of index i (invariant)
             cs[(ii-iii)/T][i-ii][i-ii] += A[i][k] * B[k][i];
   forall (jj=jjj; jj<min(jjj+T*T,N); jj+=T) {      // Parallel Local 1</pre>
     forall (j=jj; j<min(jj+T,N); j++) {
                                       // Parallel Local 0
       for (i = ii; i < min(ii+T, M); i++) {
                                                // sequential
         C[i][i] = cs[(ii-iii)/T][i-ii][i-ii];
```

- ➤ Open File "Day2-Exercises/MMMult/mmm.c", function named "runGPUverMMM".
- ► Fill the correct sizes for the "rgblkMMM" kernel.
- ► Open File "Day2-Exercises/MMMult/mmm.cl" and implement kernel named "rgblkMMM".
- ► Fix bugs until it validates! :-) Tile size is called RT and not T! OpenCL equivalences:

- ► Open File "Day2-Exercises/MMMult/mmm.c", function named "runGPUverMMM".
- ► Fill the correct sizes for the "rgblkMMM" kernel.
- ► Open File "Day2-Exercises/MMMult/mmm.cl" and implement kernel named "rgblkMMM".
- ► Fix bugs until it validates! :-) Tile size is called RT and not T!

#### OpenCL equivalences:

- ► Global worksize 1:  $\lceil M/RT \rceil$ ; Global worksize 0:  $\lceil N/RT^2 \rceil * RT^2$ ;
  - ▶ Local worksize 0: RT²; Local worksize 1: 1;
    ▶ ii = qet\_qroup\_id(1) \* RT;
  - j = jj + get\_global\_id(0);
  - locx = get\_local\_id(0) % RT; locy = get\_local\_id(0) / RT;
  - ► Loops of index i should be normalized [0..RT-1] and unrolled (#pragma unroll);
    - ► Array cs should be one dimensional: float cs[RT];
    - ► \_\_local real A\_loc[RT][RT+1]; for bank conflicts.