



The OpenMP reduction minimal example revisited: Data Sharing

Example (OpenMP reduction minimal example)

```
#include <omp.h>
#include <stdio.h>
#include <stdlib.h>
int main (int argc, char *argv[]) {
int i, n;
float a[100], b[100], sum;
/* Some initializations */
n = 100;
for (i=0; i < n; i++)</pre>
 a[i] = b[i] = i * 1.0;
sum = 0.0;
#pragma omp parallel for reduction (+: sum)
  for (i=0; i < n; i++)
    sum = sum + (a[i] * b[i]);
printf(" ____Sum = %f\n", sum);
```



The main properties of the reduction are

- $lue{}$ accumulation of data via a binary operator (here +)
- intrinsically sequential operation causing a race condition in multi-thread based implementations (since every iteration step depends on the result of its predecessor.)

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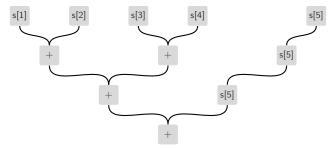


Figure: Tree reduction basic idea.



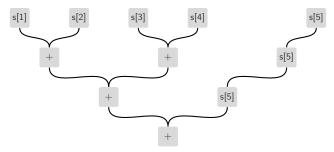


Figure: Tree reduction basic idea.

- ideally the number of elements is a power of 2
- best splitting of the actual data depends on the hardware used



Practical tree reduction on multiple cores

Example (Another approach for the dot example)

Consider the setting as before $a, b \in \mathbb{R}^{100}$. Further we have four equal cores. How do we compute the accumulation in parallel?



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1. Task pool approach: define a task pool and feed it with n/2 = 50 work packages accumulating 2 elements into 1. When these are done, schedule the next 25 and so on by further binary accumulation of 2 intermediate results per work package.



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Consider the setting as before a, $b \in \mathbb{R}^{100}$. Further we have four equal cores. How do we compute the accumulation in parallel? Basically 2 choices

- 1. Task pool approach: define a task pool and feed it with n/2 = 50 work packages accumulating 2 elements into 1. When these are done, schedule the next 25 and so on by further binary accumulation of 2 intermediate results per work package.
- 2. **#Processors=#Threads approach:** Divide the work by the number of threads, i.e. on our 4 cores each gets 25 subsequent indices to sum up. The reduction is then performed on the results of the threads.



Repetition blocked algorithms

Algorithm 1: Gaussian elimination – row-by-row-version

```
Input: A \in \mathbb{R}^{n \times n} allowing LU decomposition

Output: A overwritten by L, U

1 for k = 1: n - 1 do

2 A(k+1:n,k) = A(k+1:n,b)/A(k,k);

3 for i = k+1:n do

4 for j = k+1:n do

5 A(i,j) = A(i,j) - A(i,k)A(k,j);
```



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Observation:

- Innermost loop performs rank-1 update on the A(k+1:n,k+1:n) submatrix in the lower right,
- i.e. a BLAS level 2 operation.



Repetition blocked algorithms

Algorithm 2: Gaussian elimination – Outer product formulation

```
Input: A \in \mathbb{R}^{n \times n} allowing LU decomposition Output: L, U \in \mathbb{R}^{n \times n} such that A = LU stored in A stored in A 1 for k = 1 : n - 1 do 2 | rows= k + 1 : n:
```

3
$$A(rows, k) = A(rows, k)/A(k, k);$$

$$A(rows, rows) = A(rows, rows) - A(rows, k)A(k, rows);$$



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1 **for** k = 1 : n - 1 **do**

rows= k + 1 : n;

 $3 \quad | \quad A(rows, k) = A(rows, k)/A(k, k);$

4 | A(rows, rows) = A(rows, rows) - A(rows, k)A(k, rows);

Idea of the blocked version

- Replace the rank-1 update by a rank-r update ,
- Thus replace the $\mathcal{O}(n^2)$ / $\mathcal{O}(n^2)$ operation per data ratio the more desirable $\mathcal{O}(n^3)$ / $\mathcal{O}(n^2)$ ratio,
- Therefore exploit the fast local caches of modern CPUs more optimally.



Repetition blocked algorithms

Algorithm 3: Gaussian elimination – Block outer product formulation

```
Input: A \in \mathbb{R}^{n \times n} allowing LU decomposition, r prescribed block size
  Output: A = LU with L, U stored in A
1 k = 1:
2 while k < n do
      \ell = \min(n, k+r-1);
    Compute A(k : \ell, k : \ell) = \tilde{L}\tilde{U} via Algorithm 7;
    Solve \tilde{L}Z = A(k : \ell, \ell + 1 : n) and store Z in A;
     Solve W\tilde{U} = A(\ell + 1 : n, k : \ell) and store W in A;
      Perform the rank-r update:
      A(\ell+1:n,\ell+1:n) = A(\ell+1:n,\ell+1:n) - WZ;
    k = \ell + 1;
```



 $k = \ell + 1$:

Dense Linear Systems of Equations

Repetition blocked algorithms

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```
Input: A \in \mathbb{R}^{n \times n} allowing LU decomposition, r prescribed block size Output: A = LU with L, U stored in A

1 k = 1;

2 while k \le n do

3 | \ell = \min(n, k + r - 1);

4 | Compute A(k : \ell, k : \ell) = \tilde{L}\tilde{U} via Algorithm 7;

5 | Solve \tilde{L}Z = A(k : \ell, \ell + 1 : n) and store Z in A;

6 | Solve W\tilde{U} = A(\ell + 1 : n, k : \ell) and store W in A;

7 | Perform the rank-r update:

A(\ell + 1 : n, \ell + 1 : n) = A(\ell + 1 : n, \ell + 1 : n) - WZ;
```

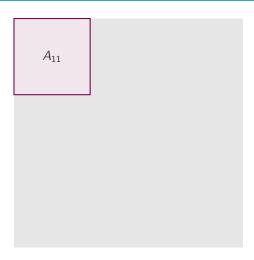
The block size r can be further exploited in the computation of W and Z and the rank-r update. It is used to optimize the data portions for the cache.



Repetition blocked algorithms

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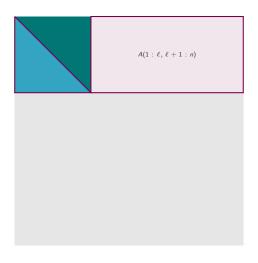




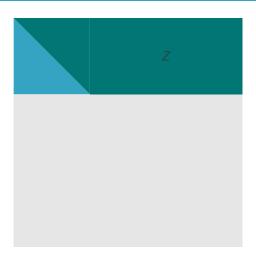






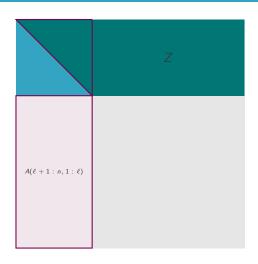






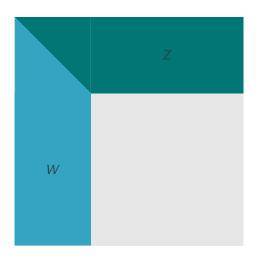


Repetition blocked algorithms

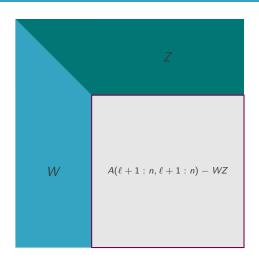


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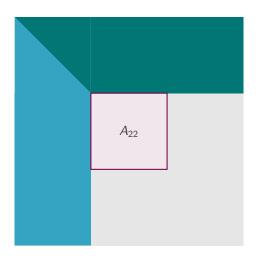












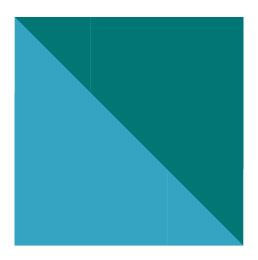














Fork-Join parallel implementation for multicore machines

We have basically two ways to implement naive parallel versions of the block outer product elimination in Algorithm 6.

Threaded BLAS available

- Compute line 4 with the sequential version of the LU
- Exploite the threaded BLAS for the block operations in lines 5–7

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Netlib BLAS

- Compute line 4 with the sequential version of the LU
- Employ OpenMP/PThreads to perform the BLAS calls for the block operations in lines 5–7 in parallel.



Fork-Join parallel implementation for multicore machines

Both these approaches fall into the class of parallel codes described by the following paradigm.

Definition (Fork-Join Parallelism)

An algorithm that performs certain parts sequentially between others that are executed in parallel is *called fork-join-parallel*.

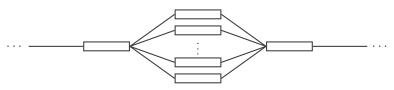


Figure: A sketch of the fork-join execution model.



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Dense Linear Systems of Equations

Fork-Join parallel implementation for multicore machines

Advantages

- Easy to achieve.
- Many threaded BLAS implementations available.
- Basically usable from any user code that requires linear system solves.

Disadvantages

- Very naive implementation.
- Sequential fraction limits the speedup (Amdahl's law).
- Therefore, only useful for small numbers of cores.



DAG scheduling of block operations aiming at manycore systems

Definition (Directed Acyclic Graph (DAG))

A directed acyclic graph is a graph where

- all edges have one distinct direction,
- directions are such that no cycles are possible for any path in the graph.

Where is the connection to parallel mathematical algorithms?

• Consider every node in the graph a task in the computation.



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- Every task requires a certain number of previous tasks to have finished.
- Also none of the previous tasks depend on the later ones.
- Thus, the dependencies give us the directions and cycles can not appear by construction.



 $\mathsf{D}\mathsf{A}\mathsf{G}$ scheduling of block operations aiming at manycore systems





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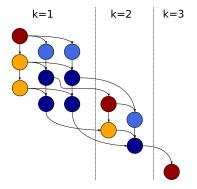


Figure: Dependency graph of Algorithm 6 for a 3×3 block subdivision.



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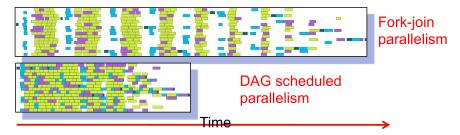


Figure: The superiority of DAG scheduling of tasks over fork-join parallelism.