# Optimizing the Performance of Multi-threaded Linear Algebra Libraries, a Task Granularity based Approach

Scientific Computing Around Louisiana (SCALA)

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#### **Outline**

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Future Work

# **Objective**

#### **Objective**

A compile-time and runtime solution to optimize the performance of a linear algebra library based on

- Machine architecture
- Number of cores to run the program on
- Expression to be evaluated
  - Type of operations
  - Number of matrices involved
  - Matrix sizes

# \_\_\_\_

Introduction

#### Introduction

- Current programming models would not be able to keep up with the advances toward exascale computing
  - More complex machine architectures, deeper memory hierarchies, heterogeneous nodes, complicated networks
- AMT(Asynchronous Many-Task) model and runtime systems
  - Examples: HPX, Charm++, Legion

#### Introduction

- Performance of HPC applications heavily rely on the linear algebra library they are using.
- Linear algebra libraries
  - BLAS(Basic Linear Algebra Subprograms) are the fundamental routines for basic vector and matrix operations.
  - Examples: ScaLAPACK, ATLAS, SPIRAL
- Our motivation:
  - Phylanx, a platform to run your python code in parallel and distributed with machine learning as the target application.

# Background

#### **Background: HPX**

- HPX is a general purpose C++ runtime system for parallel and distributed applications of any scale.
- Fine-grained parallelism instead of heavyweight threads.
- Four major factors for performance degradation: SLOW
  - Starvation
  - Latency
  - Overheads
  - Waiting for contention resolution

## **Background: Blaze C++ Library**



Blaze is a high performance C++ linear algebra library based on Smart Expression Templates.

- Expression Templates:
  - Creates a parse tree of the expression at compile time and postpone the actual evaluation to when the expression is assigned to a target
- Smart:
  - Integration with highly optimized compute kernels
  - Selecting optimal evaluation method automatically for compound expressions

#### Background: Blaze, Parallelization

Depending on the operation and the size of operands, the assignment could be parallelized through four different backends

• HPX, OpenMP, C++ threads, Boost

In the current implementation, the work is equally divided between the cores at compile time.

Parallel for loop



Figure 1: An example of how C=A+B is performed in parallel in Blaze with 4 cores

#### **Background: Task Granularity**

Grain size: The amount of work performed by one task

- What causes performance degradation?
  - Overheads
  - Starvation

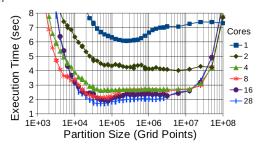


Figure 2: The effect of task size on execution time for Stencil application 1

 $<sup>^{1}</sup>$ Grubel, Patricia, et al. "The performance implication of task size for applications on the hpx runtime system." 2015 IEEE International Conference on Cluster Computing. IEEE, 2015.

## Background: Modeling Performance based on number of cores

· Amdahl's Law

$$S(p) = \frac{p}{1 + \sigma(p-1)}$$

Universal Scalability Law(USL)

$$S(p) = \frac{p}{1 + \sigma(p-1) + \kappa p(p-1)}$$

 Models the effects of linear speedup, contention delay, and coherency delay due to crosstalk

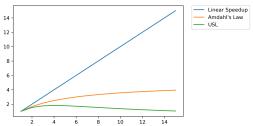


Figure 3: An example of speedup based on Amdahl's law and USL compared to the ideal linear speedup where  $\sigma=0.2$  and  $\kappa=0.05$ .

# Method

#### Method: Objective

Dynamically divide the work among the cores based on number of cores, matrix size, complexity of the operation, machine architecture. For this purpose two parameters have been introduced:

- block\_size: at each loop iteration the assignment is performed on one block
- chunk\_size: the number of loop iterations included in one task

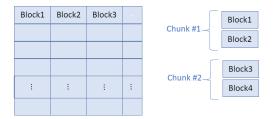


Figure 4: An example of blocking a matrix and creating chunks for chunk\_size = 2

#### Method: Data Collection

#### • Starting from DMATDMATADD benchmark: C = A + B

Category	Configuration
Matrix sizes	200, 230, 264, 300, 396, 455, 523, 600, 690, 793, 912, 1048, 1200, 1380, 1587
Number of cores	1, 2, 3, 4, 5, 6, 7, 8
Number of rows in the block	4, 8, 12, 16, 20, 32
Number of columns in the block	64, 128, 256, 512, 1024
Chunk size	Between 1 and total number of blocks (logarithmic increase)

Table 1: List of different values used for each variable for running the *DMATDMATADD* benchmark

#### Method: Data Analysis

 For simplicity we look at each matrix size individually, one number of core at a time

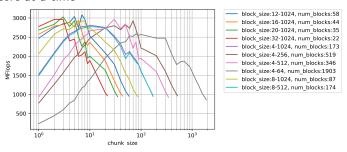
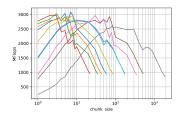


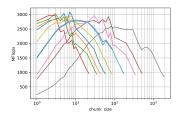
Figure 5: The results obtained from running DMATDMATADD benchmark for matrix sizes  $690 \times 690$  with different combinations of block size and chunk size on 4 cores

#### Method: Observation



- For each selected block size, there is a range of chunk sizes that gives us the best performance.
- Except for some uncommon cases, no matter which block size we choose, we are able to achieve the maximum performance if we select the right chunk size.

#### Method: Observation

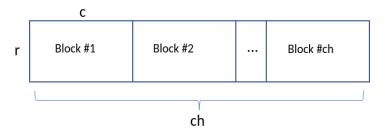


- For each selected block size, there is a range of chunk sizes that gives us the best performance.
- Except for some uncommon cases, no matter which block size we choose, we are able to achieve the maximum performance if we select the right chunk size.
- Instead of looking at block size and chunk size individually, look at grain size.

#### Method: Throughput vs. Grain Size

Grain size: The number of floating point operations performed by one thread

- Grain size represents the complexity of the expression
- For DMATDMATADD, with  $block\_size = r \times c$  and  $chunk\_size = ch$  $Grain\_size = r \times c \times ch$



#### Method: Throughput vs. Grain Size

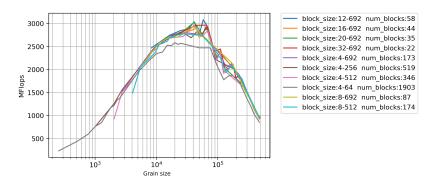


Figure 6: The results obtained from running DMATDMATADD benchmark through Blazemark for matrix size  $690 \times 690$  on 4 cores.

#### Method: Throughput vs. Grain Size

The range of grain size for maximum performance

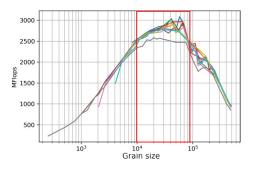


Figure 7: The results obtained from running DMATDMATADD benchmark through Blazemark for matrix size  $690 \times 690$  on 4 cores.

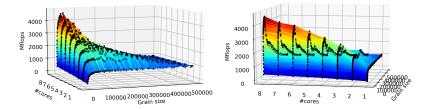


Figure 8: The results obtained from running DMATDMATADD benchmark through Blazemark for matrix size  $690 \times 690$  based on grain size and number of cores.

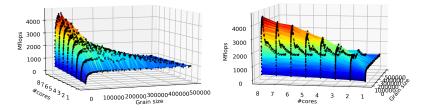


Figure 8: The results obtained from running DMATDMATADD benchmark through Blazemark for matrix size  $690 \times 690$  based on grain size and number of cores.

Can we model the relationship between the **throughput** and the **grain** size and the **number of cores**?

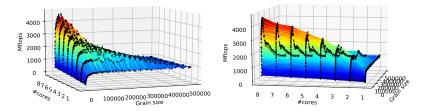


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Can we model the relationship between the **throughput** and the **grain** size and the **number of cores**?

 First we try to model the relationship between throughput and grain size.

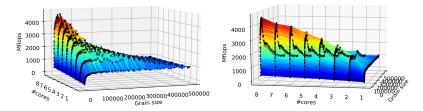


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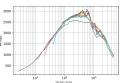
- First we try to model the relationship between throughput and grain size.
  - Polynomial Model

#### Method: Polynomial Model

In order to simplify the process and eliminate the effect of different possible factors, we started with limiting the problem to a fixed matrix size.

 Used a second order polynomial to model the relationship between the throughput and the grain size when number of cores is fixed.

$$P = ag^2 + bg + c$$



Divide the data into training(60%) and test(40%)

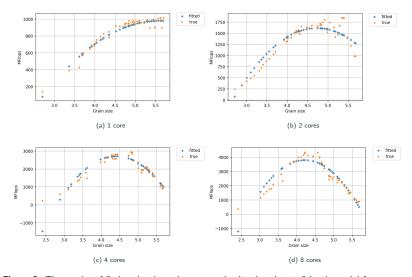
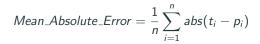


Figure 9: The results of fitting the throughput vs grain size data into a 2d polynomial for DMATDMATADD benchmark for matrix size  $690 \times 690$  with different number of cores on the test data.



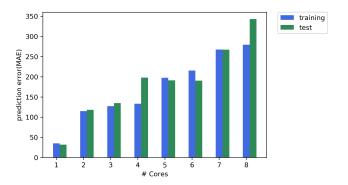


Figure 10: The training and test error for fitting data obtained from the DMATDMATADD benchmark for matrix size  $690 \times 690$  against different number of cores cores.

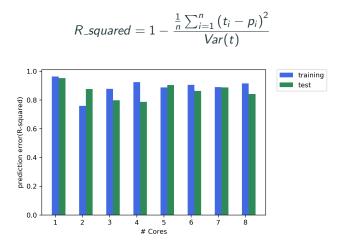
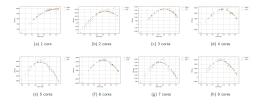
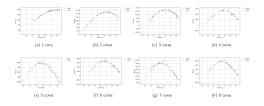


Figure 11: The training and test error for fitting data obtained from the DMATDMATADD benchmark for matrix size  $690 \times 690$  against different number of cores cores.



• We have developed a model for each number of cores:1, 2,.., 8  $P = ag^2 + bg + c$ 



- We have developed a model for each number of cores:1, 2,.., 8  $P = ag^2 + bg + c$
- Can we somehow integrate number of cores into the model?

• For  $P = ag^2 + bg + c$ , how do a, b, and c change with the number of cores?

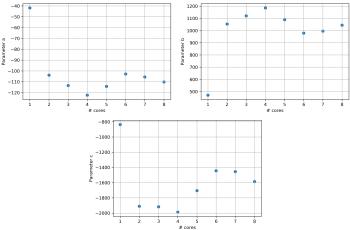


Figure 12: The parameters of the polynomial fit from the DMATDMATADD benchmark for matrix size  $1587 \times 1587$  against different number of cores.

Model the relationship with a 3rd degree polynomial

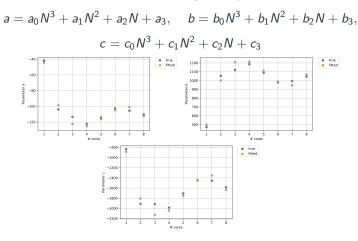


Figure 13: Fitting the parameters of the polynomial function with a 3rd degree polynomial from the DMATDMATADD benchmark for matrix size  $1587 \times 1587$  against different number of cores.

The final model:

$$P = a_{11}g^2N^3 + a_{10}g^2N^2 + ... + a_1N + a_0$$

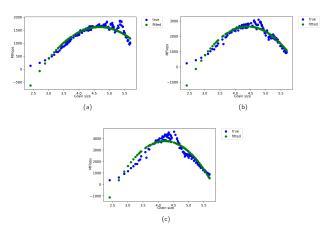


Figure 14: matrix size  $690 \times 690$  for (a) 2 core, (b) 4 cores, (c) 8 cores.

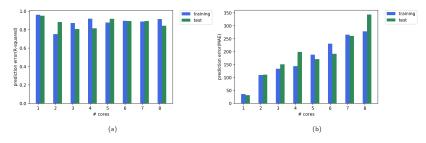


Figure 15: All the data points are included in calculation of error,(a) R-squared error (b) Mean Absolute Error(MAE) .

# Method: Finding the Grain Size Range for Maximum Performance

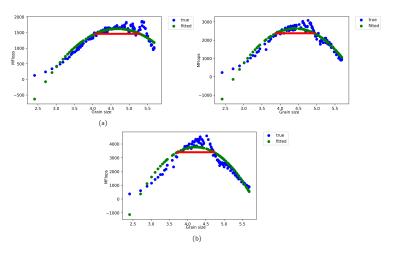
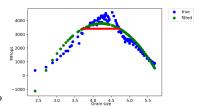


Figure 16: The range of grain size (shown as the red line) that leads to a performance within 10% of the maximum performance for (a) 2 cores, (b) 4 cores and (b) 8 cores.

# Method: Finding the Grain Size Range for Maximum Performance



How do we use the calculated range?

- ullet Select a reasonable block size, e.g.  $4 \times 256$
- Find the range of chunk size that results in the calculated range of grain size, Grain\_size=r × c × ch

# Method: Finding the Grain Size Range for Maximum Performance

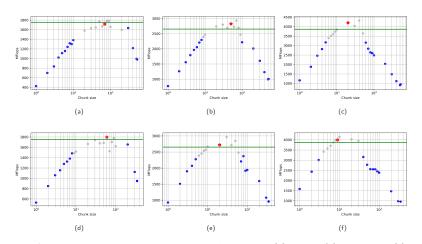


Figure 17: matrix size  $690 \times 690$  with block size of  $4 \times 256$  on (a) 2 cores, (b) 4 cores, and (c) 8 cores, and block size of  $4 \times 512$  on (d) 2 cores, (e) 4 cores, and (f) 8 cores.

## Method: Polynomial Model, Wrap up

### Strength:

• Simple model, can easily find the maximum.

#### Weakness:

• It is not physical.

### Setup: Blazemark

Blazemark is a benchmark suite provided by Blaze to compare the performance of Blaze with other linear algebra libraries.

```
Dense Vector/Dense Vector Addition:
 C-like implementation [MFlop/s]:
    100
                1115.44
                206.317
    10000000
 Classic operator overloading [MFlop/s]:
    100
                415.703
                112.557
    10000000
 Blaze [MFlop/s]:
    100
                2602.56
    10000000
                292.569
 Boost uBLAS [MFlop/s]:
    100
                1056.75
    10000000
                208.639
 Blitz++ [MFlop/s]:
    100
                1011.1
    10000000
                207.855
 GMM++ [MFlop/s]:
    100
                1115.42
    10000000
                207.699
 Armadillo [MFlop/s]:
    100
                1095.86
    10000000
                208.658
 MTL [MFlop/s]:
                1018.47
    100
                209.065
    10000000
 Eigen [MFlop/s]:
    100
                2173.48
    10000000
                209.899
```

```
N=100, steps=55116257
  C-like
             = 2.33322
                        (4.94123)
  Classic
             = 6.26062
                        (13.2586)
                        (2.11777)
  Blaze
              = 1
  Boost uBLAS = 2,4628
                        (5.21565)
                        (5.4511)
  Blitz++
             = 2.57398
  GMM++
                        (4.94129)
             = 2.33325
 Armadillo
             = 2.3749
                        (5.0295)
  MTI
             = 2.55537
                        (5.41168)
  Eigen
              = 1.19742
                        (2.53585)
N=10000000, steps=8
 C-like
             = 1.41805
                        (0.387753)
 Classic
             = 2.5993
                        (0.710753)
                        (0.27344)
  Blaze
             = 1
  Boost uBLAS = 1.40227
                        (0.383437)
  Blitz++
             = 1.40756
                        (0.384884)
  GMM++
              = 1.40862
                        (0.385172)
  Armadillo
             = 1.40215
                        (0.383403)
             = 1.39941
                        (0.382656)
  MTL
 Eigen
             = 1.39386
                        (0.381136)
```

Figure 18: An example of results obtained from Blazemark

## **Setup: Configuration**

Category	Specification
CPU	2 x Intel(R) Xeon(R) CPU E5-2450 0 @ 2.10GHz
RAM	48 GB
Number of Cores	16
Hyperthreading	Off

Table 2: Specifications of the Marvin node from Rostam cluster at CCT.

Library	Version
HPX	1.3.0
Blaze	3.5

**Table 3:** Specifications of the libraries used to run our experiments.

# Related Work

### **Related Work**

- Liu et al. estimated the optimal number of cores to run the program on based on cache specific traces.
- Khatami et al. used logistic regression to find the best chunk size based on some static and dynamic features of the loop.
- Thoman et al. proposed a compile-time and runtime solution, using an effort estimation function set the chunk size.

### **Our Contributions**

- We propose a novel analytic model to represent how the execution time is expected to change based on grain size.
- To our knowledge, there has not been a work to create a 3D model of the throughput, grain size, and number of cores.
- We are proposing a method to apply the developed model to a linear algebra library, in a way specific to our application, and the machine architecture.

# Future Work

### **Future Work**

- Generalizing the current model to different matrix sizes
- Looking into more complex expressions
- Finding an analytical model to explain the observed behavior between execution time, grain size, and number of cores.

Thank you!

### **Appendix**

### **BLAS** operations

 $\textbf{Figure 19:} \ \, \texttt{https://web.stanford.edu/class/ee392o/nlas-foils.pdf}$ 

## **Appendix**

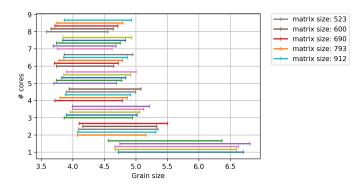


Figure 20: The range of grain size within 10% of the maximum performance of the fitted polynomial function for *DMATDMATADD* benchmark for different number of cores for matrix size  $523 \times 523$  to  $912 \times 912$ .