Legolas++: automatic vectorisation for tensor-based algorithms

Software Development Workshop for Enterprise, HPC and Al Tue 10th & Wed 11th December 2019

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Purpose: Apply 1 algorithm to N problems

LEGOLAS++ Arrays

Solving N Tridiagonal systems

Under the hood

Performances!



Legolas++ arrays

Purpose: Apply 1 algorithm to N problems

LEGOLAS++ Arrays

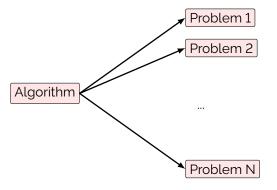
Solving N Tridiagonal systems

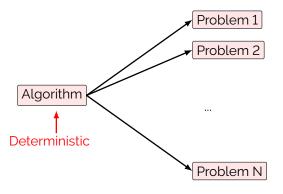
Under the hood

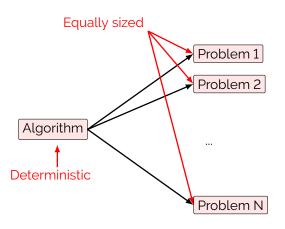
Performances



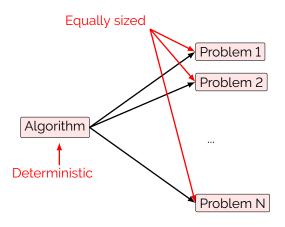
Algorithm







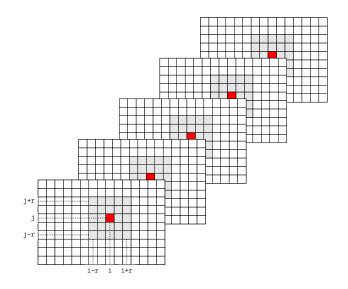




→ Vectorised and parallel Implementation



Example: Convolution Filter



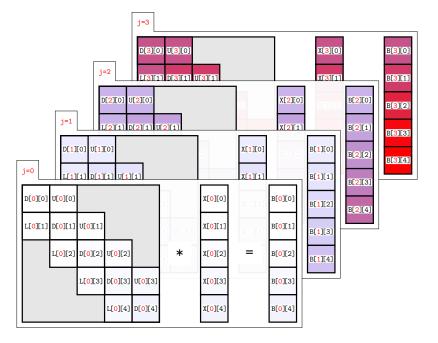


Running Example: Tridiagonal Linear Systems

Thomas Algorithm for TX = B:

| D[0] | U[0] | | | | | X[0] | | B[0] |
|------|------|------|------|------|---|------|---|------|
| L[1] | D[1] | U[1] | | | | X[1] | | B[1] |
| | L[2] | D[2] | U[2] | | * | X[2] | = | B[2] |
| | | L[3] | D[3] | U[3] | | x[3] | | B[3] |
| | | | L[4] | D[4] | | X[4] | | B[4] |

 \rightarrow algo : Thomas(L,D,U,X,B)



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LEGOLAS++ Arrays

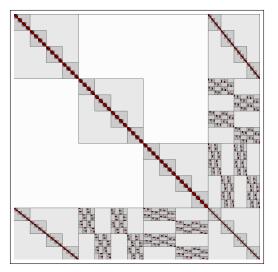
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LEGOLAS++ library

LEGOLAS++: building blocks for linear algebra solvers. Highly structured sparse linear algebra.



LEGOLAS++ Array<T,D> Template class

Rectangular N-Dimensional Arrays (Tensors):

```
//3D Array containing 100 double elts
Legolas::Array < double ,3 > X3D(10,5,2);
```



LEGOLAS++ Array<T,D> Template class

Rectangular N-Dimensional Arrays (Tensors):

```
//3D Array containing 100 double elts
Legolas::Array < double ,3 > X3D(10,5,2);

Legolas::Array < double ,2 > X2D = X3D[0];

for (int k=0; k<10; k++)
  for (int j=0; j<5; j++)
   for (int i=0; i<2; i++)
      X3D[k][j][i]=2.*X3D[k][j][i]+1.;</pre>
```

LEGOLAS++ Parallel Expression Template

```
//3D Array containing 100 double elts
 Legolas:: Array < double, 3 > X (10,5,2);
 Legolas:: Array < double, 3 > Y(10,5,2);
 Legolas::Array <double, 3> Z(10,5,2);
 for (int k=0; k<10; k++)</pre>
  for (int j=0; j<5; j++)
   for (int i=0; i<2; i++)
   Y[k][j][i]+=2.*X[k][j][i]+Z[k][j][i];
Y+=2.*X+Z; //Expression Template
Y+=2.*X+Z || seq; // Sequential
Y+=2.*X+Z || par; // MultiThreaded
```

LEGOLAS++ Array<T,D> Nested Types

| Array <t,d> Nested Types</t,d> | | | | | | | |
|--------------------------------|---|--|--|--|--|--|--|
| Scalar | Array <t,d>::Scalar = T</t,d> | | | | | | |
| Element | <pre>Element Array<t,d>::Element</t,d></pre> | | | | | | |
| | $$ Array <t,d-1> if $D>1$</t,d-1> | | | | | | |
| | $\hat{=}$ T& if $D=1$ | | | | | | |

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LEGOLAS++ Arrays

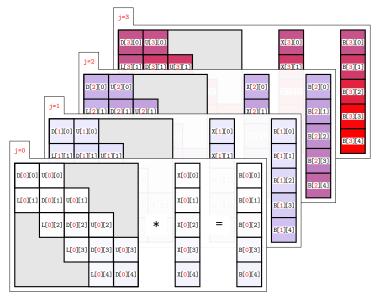
Solving N Tridiagonal systems

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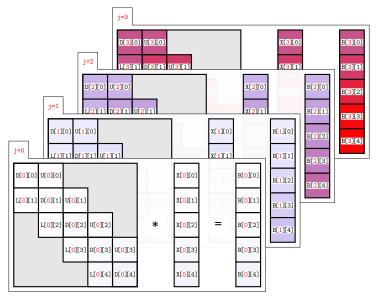


Legolas::Array<float,2> L,D,U,X,B (L,D,U)[j][i]



→ algo : Thomas

Legolas::Array<float,2> L,D,U,X,B (L,D,U)[j][i]



 \rightarrow algo: Thomas(L[j],D[j],U[j],X[j],B[j])



Solving N Tridiagonal systems

```
struct MultiThomasSolver{
  template <class A2D>
  void operator()(int begin, int end,
                   A2D D2D, A2D U2D, A2D L2D
                   A2D B2D, A2D X2D) const{
    using Element = typename A2D:: Element;
    using Scalar = typename A2D::Scalar;
    Element S(X2D[0].shape());
    //Loop on tridiagonal systems
    for (int j=begin ; j<end ; j++){</pre>
      auto D=D2D[j]; auto U=U2D[j]; auto L=L2D[j];
      auto B=B2D[i]; auto X=X2D[i];
      //Thomas Algorithm starts here
        //forward...
//backward...
}}:
```

Thomas Algorithm : TX=B with T=(L,D,U)

```
Scalar s=D[0], one=1.0, sm1=one/s;
const int size=X.size():
//forward
X[0] = B[0] * sm1;
for(int i=1; i<size; i++){</pre>
  S[i]=U[i-1]*sm1;
  s=D[i]-L[i]*S[i]:
  X[i]=B[i]-L[i]*X[i-1];
  sm1=one/s;
  X[i] *= sm1;
//backward
for (int i=(size-2);i>=0; i--){
  X[i] -= S[i+1] * X[i+1]:
};
```

LEGOLAS++ map

```
int main () {
  size_t ni=200;//System size
  size_t nj=800; // Number of systems
  using A2D=Legolas::Array<float,2>;
  A2D u(nj,ni),l(nj,ni),d(nj,ni);
  A2D X(nj,ni),B(nj,ni);
  //.. Arrays initialization
  . . .
  //Solve all systems (sequential)
  Legolas::map(MultiThomasSolver(),d,u,1,B,X);
```

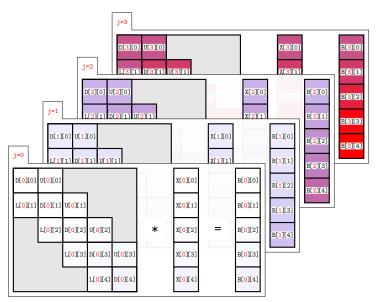
SIMD Parallelism

SIMD : Single Instruction Multiple Data \rightarrow CPU (MMX,SSE,AVX,AVX2,AVX512,...)



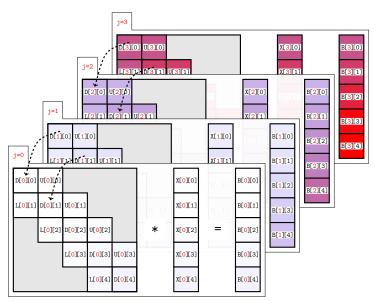
 \rightarrow OK for contiguous data!

Vectorisation with SIMD width=2





Vectorisation with SIMD width=2





Vectorisation with SIMD width=2





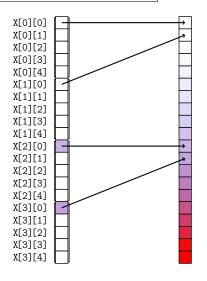
LEGOLAS++: Data Layout Interleaving

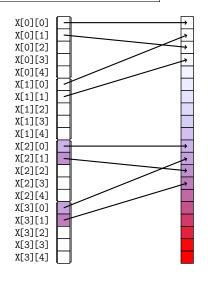
The LEGOLAS++ Array template class:

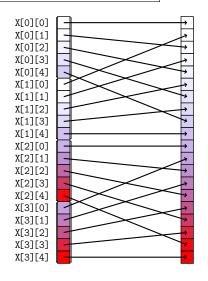
accepts two extra parameters P and DP that control multi-dimensional array data layout.
P specifies the *packing factor* that must be applied to the dimension dimension DP.

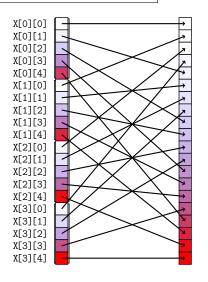
| X[0][0] | \bigcap | | |
|---------|-----------|--|--|
| X[0][1] | | | |
| X[0][2] | | | |
| X[0][3] | | | |
| X[0][4] | | | |
| X[1][0] | П | | |
| X[1][1] | | | |
| X[1][2] | | | |
| X[1][3] | | | |
| X[1][4] | | | |
| X[2][0] | | | |
| X[2][1] | | | |
| X[2][2] | | | |
| X[2][3] | | | |
| X[2][4] | | | |
| X[3][0] | Ш | | |
| X[3][1] | | | |
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| X[3][4] | Ш | | |











LEGOLAS++ for N tridiagonal systems

```
int main () {
  size_t ni=200;//System size
  size_t nj=800; // Number of systems
  using A2D=Legolas::Array<float,2>;
  A2D u(nj,ni),l(nj,ni),d(nj,ni);
  A2D X(nj,ni),B(nj,ni);
  //.. Arrays initialization
  //Solve all systems (sequential)
  Legolas::map(MultiThomasSolver(),d,u,1,B,X);
}
```

LEGOLAS++ for *N* tridiagonal systems

```
int main () {
  size_t ni=200;//System size
  size_t nj=800; // Number of systems
 //using A2D=Legolas::Array<float,2>;
 using A2D=Legolas::Array<float,2,4,2>;
 A2D u(nj,ni),l(nj,ni),d(nj,ni);
 A2D X(nj,ni),B(nj,ni);
 //.. Arrays initialization
 //Automatic SSE Vectorization
 Legolas::map(MultiThomasSolver(),d,u,1,B,X);
}
```

LEGOLAS++ for N tridiagonal systems

```
int main () {
  size_t ni=200;//System size
  size_t nj=800; // Number of systems
 //using A2D=Legolas::Array<float,2>;
 //using A2D=Legolas::Array<float,2,4,2>;
 using A2D=Legolas::Array<float,2,8,2>;
 A2D u(nj,ni),l(nj,ni),d(nj,ni);
 A2D X(nj,ni),B(nj,ni);
 //.. Arrays initialization
 //Automatic AVX Vectorization
 Legolas::map(MultiThomasSolver(),d,u,1,B,X);
}
```

LEGOLAS++ for N tridiagonal systems

```
int main () {
  size_t ni=200;//System size
  size_t nj=800;//Number of systems
 //using A2D=Legolas::Array<float,2>;
 //using A2D=Legolas::Array<float,2,4,2>;
 using A2D=Legolas::Array<float,2,8,2>;
 A2D u(nj,ni),l(nj,ni),d(nj,ni);
 A2D X(nj,ni),B(nj,ni);
 //.. Arrays initialization
 //Automatic AVX Vectorization+parallelization
 //Legolas::map(MultiThomasSolver(),d,u,1,B,X);
 Legolas::parmap(MultiThomasSolver(),d,u,1,B,X);
}
```



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LEGOLAS++



LEGOLAS++

controls Data Layout Interleaving : Legolas::Array<T,D,P,DP>

FGOLAS++

- controls Data Layout Interleaving : Legolas::Array<T,D,P,DP>
- uses Eigen for SIMD operations: http://eigen.tuxfamily.org

LEGOLAS++

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- controls Data Layout Interleaving : Legolas::Array<T,D,P,DP>
- uses Eigen for SIMD operations: http://eigen.tuxfamily.org
- uses Intel TBB for multithreading: https://www.threadingbuildingblocks.org
- ightharpoonup header-only library (<2K lines)!

Eigen Blocks (fixed size arrays)

```
typedef Eigen::Array<float,p,1> Block;
```

Eigen overloads all the usual arithmetic operators, functions and expressions.

For example:

```
Block a,b,c,d;
a+=b-c;    //_mm_add_ps
a=b*c/d;    //_mm_mul_ps,_mm_div_ps
```

are transformed, at compile time, into the corresponding explicit intrinsic SIMD functions without any performance penalty compared to hand-coded assembly or intrinsic formulation.

► Let us define X as:

Legolas::Array<float,2,4,2> X(nj,ni);

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```
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```

▶ and Block as:

```
using Block=Eigen::Array<float,4,1>;
```

Let us define X as:

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Legolas::Array<float,2,4,2> X(nj,ni);
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▶ and Block as:

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using Block=Eigen::Array<float,4,1>;
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Then X can be reinterpreted as non-interleaved array of Blocks:

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 - ► Legolas::Array<Block,2>

► Let us define X as:

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```

▶ and Block as:

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using Block=Eigen::Array<float,4,1>;
```

- Then X can be reinterpreted as non-interleaved array of Blocks:
 - ► Legolas::Array<Block,2>
- ► LEGOLAS++ getPackedView() method returns this view:



► Let us define X as:

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```

and Block as:

```
using Block=Eigen::Array<float,4,1>;
```

- Then X can be reinterpreted as non-interleaved array of Blocks:
 - Legolas::Array<Block,2>
- LEGOLAS++ getPackedView() method returns this view :
 - Legolas::Array<Block,2> Xp=X.getPackedView();

LEGOLAS++ map

```
template <class ALGO, typename... ARRAYS>
void map(ALGO a, ARRAYS... rest){
  int s=getArraysSize(rest...);//e.g 5
  int p=getPackSize(rest...); //e.g 2
  int sp=getPackedArraysSize(rest...);// 5/2=2
  // Vectorized part:
  // algo a in [0,sp=s/p[ -> [0,2[
  a(0, sp, rest.getPackedView()...);
  // Scalar remainder
 // scalar algo a a in [sp*p,s[ ->[4,5[
  a(sp*p,s,rest...);
```



```
struct MultiThomasSolver{
 template <class A2D>
 void operator()(int begin, int end,
                   A2D D2D, A2D U2D, A2D L2D
                   A2D B2D, A2D X2D) const{
    using Element = typename A2D:: Element;
    using Scalar =typename A2D::Scalar;
    Element S(X2D[0].shape());
    Scalar one (1.0), s, sm1;
    //Loop on tridiagonal systems
    for (int j=begin ; j<end ; j++){</pre>
      auto D=D2D[j]; auto U=U2D[j]; auto L=L2D[j];
      auto B=B2D[j]; auto X=X2D[j];
      const int size=X.size();
      X[0]=B[0]*sm1:
      for(int i=1; i<size; i++){</pre>
        S[i]=U[i-1]*sm1;
```

```
struct MultiThomasSolver{
 template <class A2D>
 void operator()(int begin, int end,
                  A2D D2D, A2D U2D, A2D L2D
                   A2D B2D, A2D X2D) const{
    using Element=typename A2D::Element;//Array<Block,1>
    using Scalar = typename A2D::Scalar;
    Element S(X2D[0].shape());
    Scalar one (1.0), s, sm1;
    //Loop on tridiagonal systems
    for (int j=begin ; j<end ; j++){</pre>
      auto D=D2D[j]; auto U=U2D[j]; auto L=L2D[j];
      auto B=B2D[j]; auto X=X2D[j];
      const int size=X.size();
      X[0]=B[0]*sm1:
      for(int i=1; i<size; i++){</pre>
        S[i]=U[i-1]*sm1;
```



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struct MultiThomasSolver{
 template <class A2D>
 void operator()(int begin, int end,
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    using Element=typename A2D::Element;//Array<Block,1>
    using Scalar = typename A2D::Scalar; //Block
    Element S(X2D[0].shape());
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    //Loop on tridiagonal systems
    for (int j=begin ; j<end ; j++){</pre>
      auto D=D2D[j]; auto U=U2D[j]; auto L=L2D[j];
      auto B=B2D[j]; auto X=X2D[j];
      const int size=X.size();
      X[0]=B[0]*sm1:
      for(int i=1; i<size; i++){</pre>
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```



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  template <class A2D>
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                   A2D B2D, A2D X2D) const{
    using Element=typename A2D::Element;//Array<Block,1>
    using Scalar = typename A2D::Scalar; //Block
    Element S(X2D[0].shape());
    Scalar one (1.0), s, sm1;
    //Loop on tridiagonal systems
    for (int j=begin ; j<end ; j++){</pre>
      auto D=D2D[j]; auto U=U2D[j]; auto L=L2D[j];
      auto B=B2D[j]; auto X=X2D[j];
      const int size=X.size();
      X[0] = B[0] * sm1; // _mm_mul_ps
      for(int i=1; i<size; i++){</pre>
        S[i]=U[i-1]*sm1;//_mm_mul_ps
        . . .
```



LEGOLAS++ parmap

```
template <class ALGO,typename... ARRAYS>
void parmap(int begin, int end,
ALGO algo, ARRAYS... rest){

  tbb::parallel_for(
  tbb::blocked_range<int>(begin,end),
  [=](tbb::blocked_range<int> r){
    algo(r.begin(),r.end(),rest...);},
  tbb::auto_partitioner());
}
```

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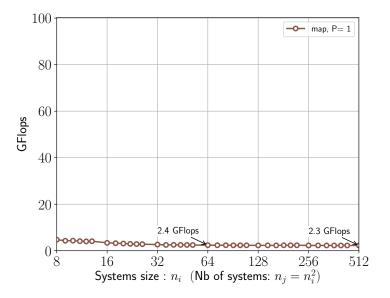
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Solving N Tridiagonal systems

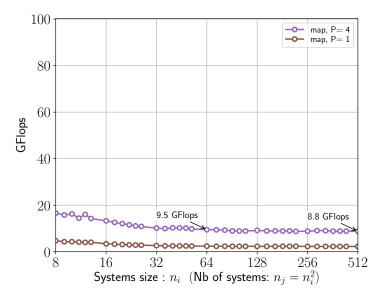
Under the hood

Performances!

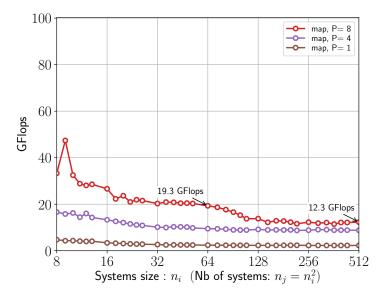




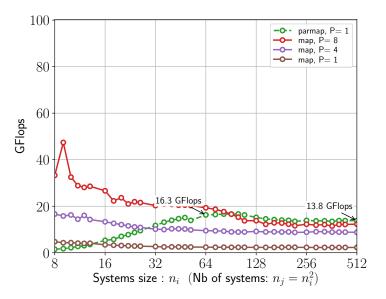




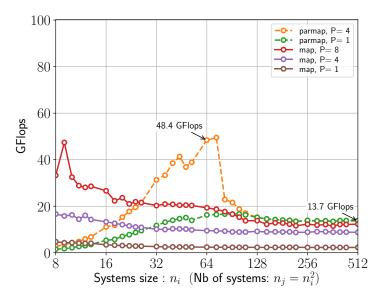




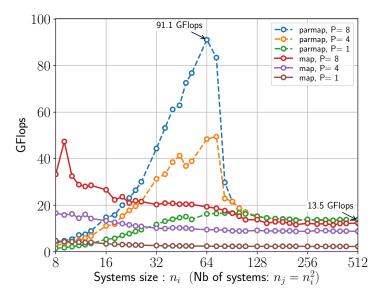






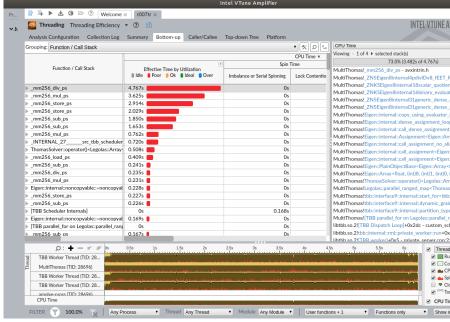








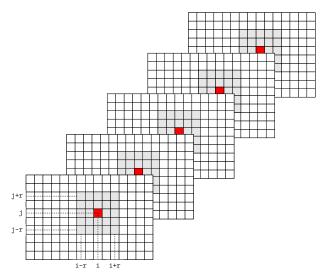
Intel VTune Amplifier

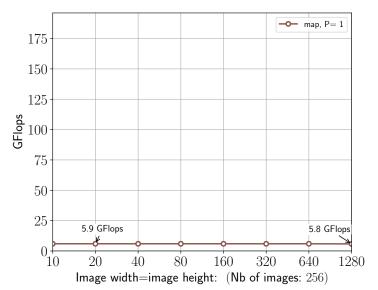


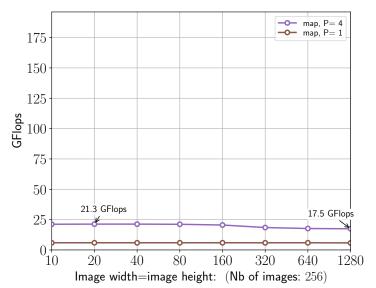


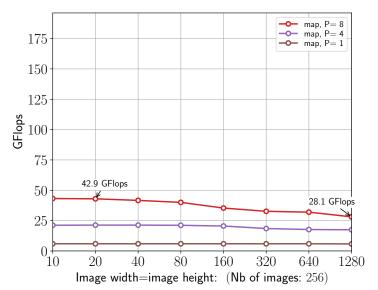
Example: Convolution Image Filter

convolution($\{(s_0, t_0), (s_1, t_1), \dots, (s_n, t_n)\}$)

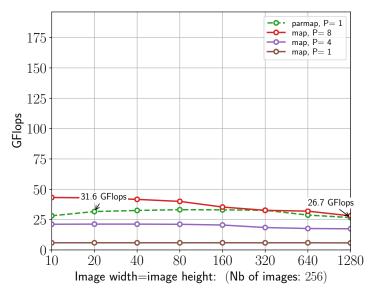


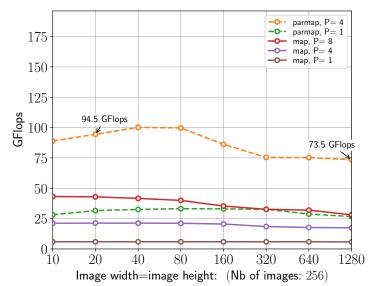




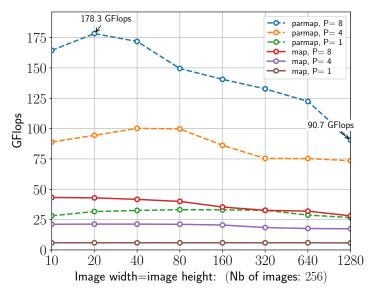














Conclusion

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 - multiple applications of a deterministic algorithm



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 - multiple applications of a deterministic algorithm
 - to equally sized problems,
- ► Legolas::Array<T,D,P,DP> enables for automatic
 - vectorisation via Data Layout Interleaving (+Eigen)
 - parallelisation (Intel TBB)

- Considering the
 - multiple applications of a deterministic algorithm
 - to equally sized problems,
- ► Legolas::Array<T,D,P,DP> enables for automatic
 - vectorisation via Data Layout Interleaving (+Eigen)
 - parallelisation (Intel TBB)
 - from a single user expression of algorithms.



Thank you for your attention!







 Applied Mathematics (Simulation, Mathematical Optimisation...);



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- ➤ Software Development (Julia, Rust, Go, C, C++, Java...);



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- Software Development (Julia, Rust, Go, C, C++, Java...);
- Hardware Architectures (SMP Clusters, GPU, DSPs, IoT...).



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The success of scientific and technical software depends on the **optimal combination of these 3 areas in strong interaction**.



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- Hardware Architectures (SMP Clusters, GPU, DSPs, IoT...).

The success of scientific and technical software depends on the **optimal combination of these 3 areas in strong interaction**.

Training

- ▶ HPC & //ism
- Julia, Go
- Numerical Quality



- Applied Mathematics (Simulation, Mathematical Optimisation...);
- Software Development (Julia, Rust, Go, C, C++, Java...);
- ► Hardware Architectures (SMP Clusters, GPU, DSPs, IoT...).

The success of scientific and technical software depends on the **optimal combination of these 3 areas in strong interaction**.

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- Architecture
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→ More information on triscale-innov.com

Legolas++ arrays

```
int main(int argc, char** argv) {
 typedef float RealType;
 typedef Legolas:: Array < Real Type, 4,8,4 > A4D;
 const int nz = 200; // nimages
 const int nc = 4; //nchannels(rgba)
 A4D sourceImages(nz, nc, ny, nx);
 A4D targetImages(nz, nc, ny, nx);
 //init source images
 typedef Convolution < GaussianKernel <5>> MyFilter;
 Legolas::map(MyFilter(), sourceImages, targetImages);
 return 0;
```

```
template <class KERNEL>
struct Convolution {
 template <class A4D>
 void operator()(int begin, int end,
 const A4D source, A4D target) const {
   typedef typename A4D::RealType Scalar;
   for (int k = begin; k < end; k++) {
     auto source3D = source[k]:
     auto target3D = target[k];
     const int nv = source3D[0].size();
     const int nx = source3D[0][0].size():
     const int r = KERNEL::radius:
     const Scalar zero(0.0):
     Scalar sumR(zero).sumG(zero).sumB(zero);
     for (int i = 0: i < nv: i++) {
       for (int i = 0: i < nx: i++) {
          sumR = zero: sumG = zero: sumB = zero:
         for (int si = -r: si <= r: si++) {
            int jp = j + sj; if (jp < 0 \mid \mid jp >= ny) jp = j - sj;
            for (int si = -r: si <= r: si++) {
             int ip = i + si: if (ip < 0 || ip >= nx) ip = i - si:
             sumR += kmat[si + r][si + r] * source3D[0][ip][ip];
             sumG += kmat[si + r][si + r] * source3D[1][ip][ip];
             sumB += kmat[sj + r][si + r] * source3D[2][jp][ip];
          }
          target3D[0][i][i] = sumR;
          target3D[1][i][i] = sumG;
         target3D[2][j][i] = sumB;
       }}}};
```

Vectorize (and parallelize) multiple applications of deterministic algorithms to regular set of datasets.



.egolas++ array

Vectorize (and parallelize) multiple applications of deterministic algorithms to regular set of datasets.

Let $V = \{v_1, v_2, \dots, v_n\}$ be a dataset consisting of n multidim-array variables.

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- Let a(V) be a *deterministic* algorithm to be applied to a dataset V:

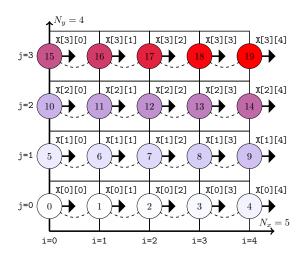
- Let $V = \{v_1, v_2, \dots, v_n\}$ be a dataset consisting of n multidim-array variables.
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 - The sequence of operations does not depend on the dataset.

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- Let $W = \{V_1, V_2, \dots, V_m\}$ be a set of m datasets. W is regular if:
 - $\qquad \forall j,j' \in [1,m]^2, \ \forall i \in [1,n], \quad \mathsf{size}(v_{i,j}) = \mathsf{size}(v_{i,j'})$

PDE on 2D Cartesian mesh

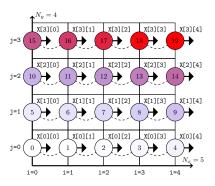
$$-X_{i-1,j} + 2X_{i,j} - X_{i+1,j} = hB_{i,j}$$

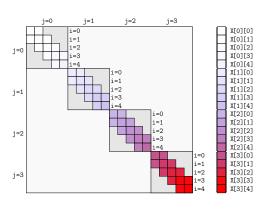




Two-level block matrix structure

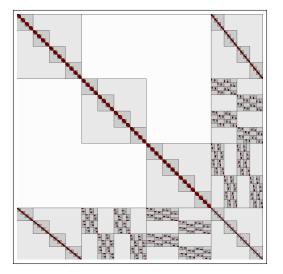
Diagonal<Tridiagonal>





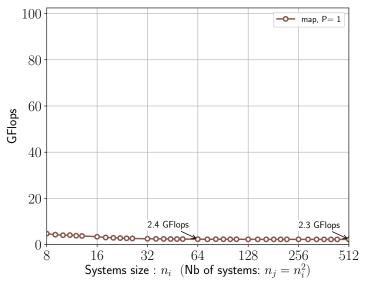
LEGOLAS++ template library

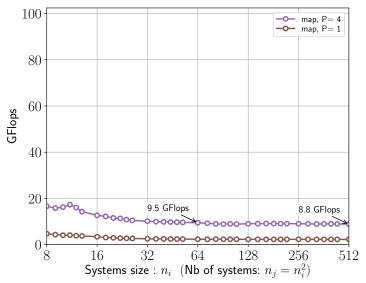
LEGOLAS++: building blocks for linear algebra solvers. Highly structured sparse linear algebra.

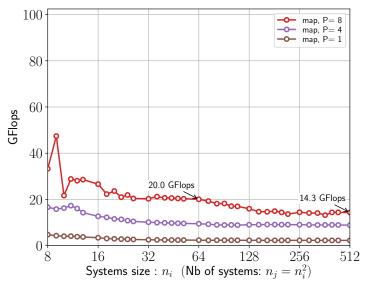


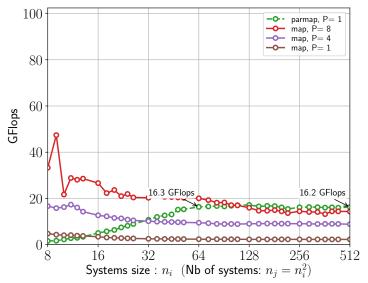
LEGOLAS++ for N tridiagonal systems

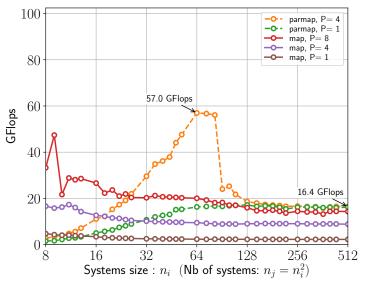
```
struct MultiThomasSolver{
  template <class A2D>
  void operator()(int begin, int end,
  A2D D2D, A2D U2D, A2D L2D, A2D B2D, A2D X2D) const{
    typedef typename A2D::Element Element;
    typedef typename A2D::Scalar Scalar;
    Element S(X2D[0].shape());
    Scalar one(1.0),s,sm1;
    //Loop over the tridiag systems
    for (int j=begin ; j<end ; j++){
      auto D=D2D[i]:auto U=U2D[i]: auto L=L2D[i]:
      auto B=B2D[j]; auto X=X2D[j];
      //Thomas Algorithm starts here
      s=D[0]; sm1=one/s;
      const int size=X.size():
      //forward
      X[0]=B[0]*sm1:
      for(int i=1; i<size; i++){</pre>
        S[i]=U[i-1]*sm1:
        s=D[i]-L[i]*S[i];
        X[i]=B[i]-L[i]*X[i-1]:
        sm1=one/s:
        X[i] *= sm1;
      //backward
      for (int i=(size-2):i>=0 : i--){
        X[i] -= S[i+1] * X[i+1];
}}};
```

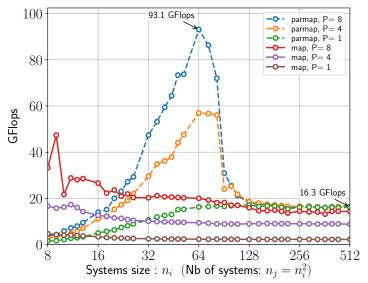




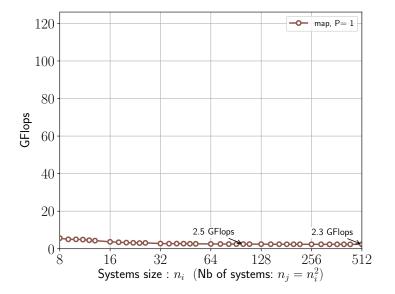




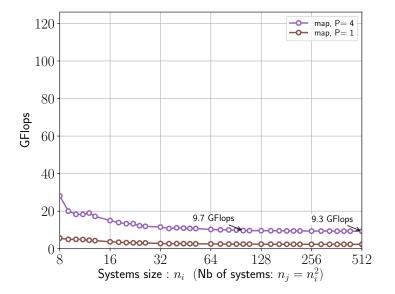


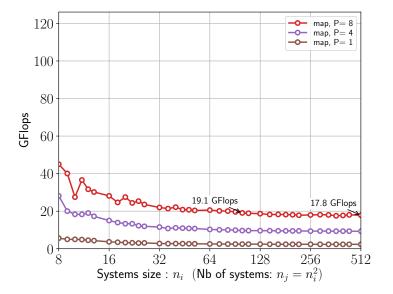


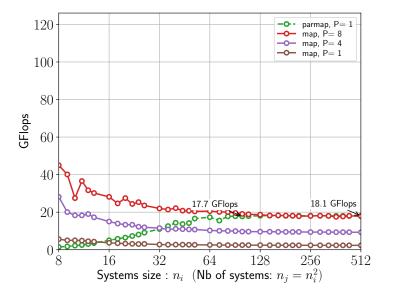
LEGOLAS++ Laplacian (Skylake 4-core,4GHz,AVX2)

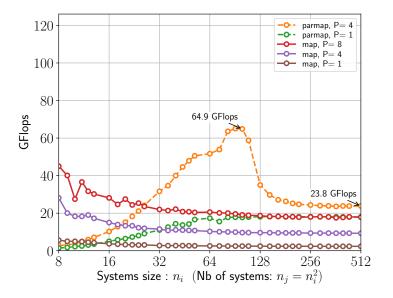


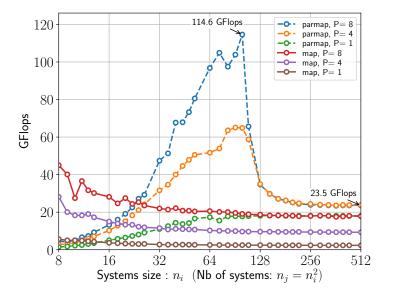
LEGOLAS++ Laplacian (Skylake 4-core,4GHz,AVX2)











LEGOLAS++ Array<T,D> Basic API

| Array <t,d> Nested Types</t,d> | |
|--------------------------------|--|
| Scalar | Array <t,d>::Scalar = T</t,d> |
| Element | Array <t,d>::Element</t,d> |
| | $\hat{=}$ Array <t,d-1> if $D>1$</t,d-1> |
| | $\hat{=}$ T& if $D=1$ |
| Shape | int tuple for array sizes |

| Array <t,d> Methods</t,d> | |
|---------------------------|--------------------------------|
| Array(int n1,,nD) | Ctor (variadic) |
| Element operator[] | Element accessor |
| int size() | returns the number of elements |
| Shape shape() | returns (n1,,nD) |



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A2D= Array<float,2>

```
struct MultiThomasSolver{
 template <class A2D>
 void operator()(int begin, int end,
                   A2D D2D, A2D U2D, A2D L2D
                   A2D B2D, A2D X2D) const{
    typedef typename A2D::Element Element;
    typedef typename A2D::Scalar Scalar;
    Element S(X2D[0].shape());
    Scalar one (1.0), s, sm1;
    //Loop on tridiagonal systems
    for (int j=begin ; j<end ; j++){</pre>
      auto D=D2D[j]; auto U=U2D[j]; auto L=L2D[j];
      auto B=B2D[j]; auto X=X2D[j];
      const int size=X.size();
      X[0]=B[0]*sm1:
      for(int i=1; i<size; i++){</pre>
        S[i]=U[i-1]*sm1;
        . . .
```

A2D= Array<float,2>

```
struct MultiThomasSolver{
 template <class A2D>
 void operator()(int begin, int end,
                  A2D D2D, A2D U2D, A2D L2D
                   A2D B2D, A2D X2D) const{
    typedef typename A2D::Element Element;//Array<float,1>
    typedef typename A2D::Scalar Scalar;
    Element S(X2D[0].shape());
    Scalar one (1.0), s, sm1;
    //Loop on tridiagonal systems
    for (int j=begin ; j<end ; j++){</pre>
      auto D=D2D[j]; auto U=U2D[j]; auto L=L2D[j];
      auto B=B2D[j]; auto X=X2D[j];
      const int size=X.size();
      X[0]=B[0]*sm1:
      for(int i=1; i<size; i++){</pre>
        S[i]=U[i-1]*sm1;
```

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```
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 template <class A2D>
 void operator()(int begin, int end,
                  A2D D2D, A2D U2D, A2D L2D
                   A2D B2D, A2D X2D) const{
    typedef typename A2D::Element Element;//Array<float,1>
    typedef typename A2D::Scalar Scalar; //float
    Element S(X2D[0].shape());
    Scalar one (1.0), s, sm1;
    //Loop on tridiagonal systems
    for (int j=begin ; j<end ; j++){</pre>
      auto D=D2D[j]; auto U=U2D[j]; auto L=L2D[j];
      auto B=B2D[j]; auto X=X2D[j];
      const int size=X.size();
      X[0] = B[0] * sm1;
      for(int i=1; i<size; i++){</pre>
        S[i]=U[i-1]*sm1;
```

LEGOLAS++ for *N* tridiagonal systems

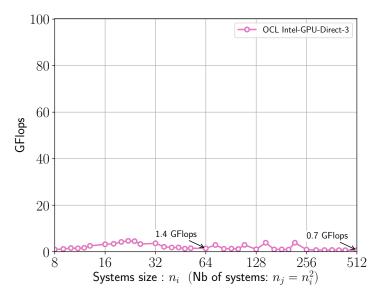
```
int main () {
  size_t ni=200;//System size
  size_t nj=800; // Number of systems
  typedef Legolas::Array<float,2> A2D;
  A2D u(nj,ni),l(nj,ni),d(nj,ni);
  A2D X(nj,ni),B(nj,ni);
  //.. Arrays initialization
  auto multiThomasSolver=MultiThomasSolver();
  //Solve all systems (sequentially)
  multiThomasSolver(0,nj,d,u,1,B,X);
```

LEGOLAS++ Parallel Expression Template

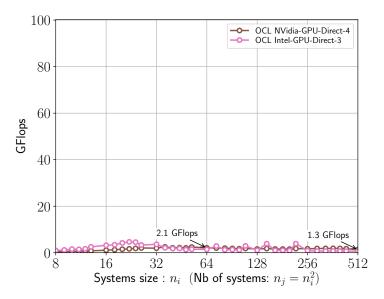
```
//3D Array containing 100 double elts
 Legolas:: Array < double, 3 > X (10,5,2);
 Legolas:: Array < double, 3 > Y(10,5,2);
 Legolas::Array <double, 3> Z(10,5,2);
 for (int k=0; k<10; k++)</pre>
  for (int j=0; j<5; j++)
   for (int i=0; j<2; i++)
   Y[k][j][i]+=2.*X[k][j][i]+Z[k][j][i];
Y+=2.*X+Z; //Expression Template
Y+=2.*X+Z || sec; // Sequential
Y+=2.*X+Z || par; // MultiThreaded
```

LEGOLAS++ OpenCL Thomas (Direct)

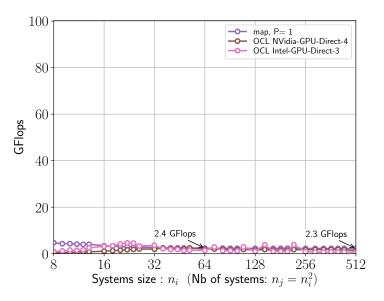
```
__kernel void thomasAlgorithm(
#define SYSTEM_SIZE %(system_size)d
__global float *L,__global float *D,
__global float *U,__global float *X,
__global float *B){
//o: system offset
int o=get_global_id(0)*SYSTEM_SIZE;
float s=D[o+0];
float sm1=1.0/s;
float S1[SYSTEM SIZE]:
//forward
X[o+0] = B[o+0] * sm1;
for (int i=1 ; i<SYSTEM_SIZE ; i++){</pre>
    S1[i]=U[o+i-1]*sm1:
    s=D[o+i]-L[o+i]*Sl[i];
    X[o+i]=B[o+i]-L[o+i]*X[o+i-1];
    sm1=1./s:
    X[o+i]*=sm1:
//backward
for (int i=(SYSTEM SIZE-2):i>=0 : i--){
    X[o+i]-=Sl[i+1]*X[o+i+1]:
}}
```



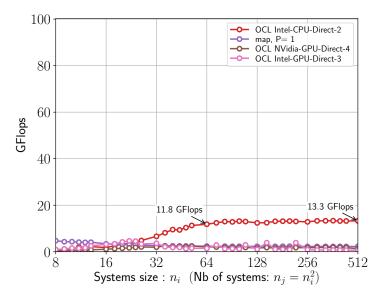




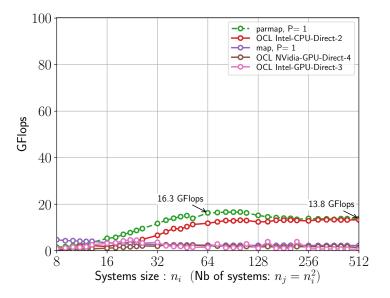




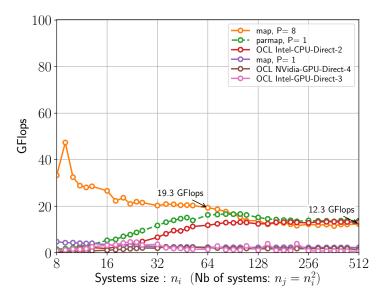




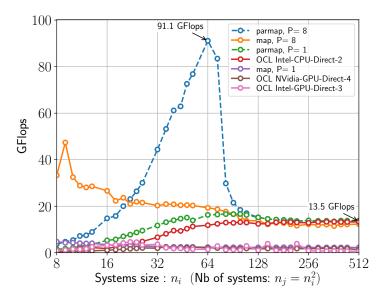












LEGOLAS++ OpenCL Thomas (Transposed)

```
kernel void thomasAlgorithm(
#define SYSTEM_SIZE %(system_size)d
__global float *L,__global float *D,__global ↔
   float *U.
__global float *X,__global float *B, unsigned ↔
   n systems)
//o: system offset
int gid=get_global_id(0);
float s=D[gid];
float sm1=1.0/s:
//forward
X[gid]=B[gid]*sm1;
//local workspace
float S1[SYSTEM_SIZE];
for (int i=1 : i<SYSTEM SIZE : i++) {
  int I=gid+n_systems*i;
  int Im1=gid+n systems*(i-1);
  Sl[i]=U[Im1]*sm1:
  s=D[I]-L[I]*S1[i];
  X[I]=B[I]-L[I]*X[Im1]:
  sm1=1./s;
  X[I]*=sm1;
}
//backward
for (int i=(SYSTEM_SIZE-2);i>=0 ; i--){
  int I=gid+n systems*i:
  int Ip1=gid+n_systems*(i+1);
  X[I]-=S1[i+1]*X[Ip1]:
11
```

