### Solving ordinary differential equations in C++

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## Outline

- Introduction
- 2 Tutorial
- Technical details
- 4 Discussion

# The interface problem in C/C++

- Many frameworks exist to do numerical computations.
- Data has to be stored in containers or collections.
- **GSL**: gsl\_vector, gsl\_matrix
- NR: pointers with Fortran-style indexing
- Blitz++, MTL4, boost::ublas
- QT: QVector, wxWidgets: wxArray, MFC: CArray

But: All books on C++ recommend the use of the STL containers std::vector, std::list,...

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GoF Design Pattern: Adaptor, also known as Wrapper

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#### Theoretical solution of the interface mess

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#### Alternative

Generic, container independent algorithms

## Example

```
void CrankNicolsonEvolution::prepareVector(gsl vector complex* phi) {
 gsl vector complex* phi temp = gsl vector complex alloc(dim);
 //we need a copy of phi for this
 gsl vector complex memcpv(phi temp, phi):
 for( int i=1; i < dim-1; i++ ) {
   //phi n = phi n - i*dt/2 * (phi n-1 + phi n+1 + pot[n]*phi n)
   gsl vector complex set(phi, i, gsl complex add(
               gsl vector complex get(phi temp, i),
               gsl complex mul imag(
                  gsl complex add(
                     gsl complex add( gsl vector complex get(phi temp, i-1),
                                   gsl_vector_complex_get(phi_temp, i+1)),
                     gsl complex mul real( gsl vector complex get(phi temp, i),
                                           potential[i] )),
                  -dt/2.0))):
  if(periodic) {
   //periodic boundaries: i=0
   gsl vector complex_set(phi, 0, gsl_complex_add(
               gsl vector complex get(phi temp, 0),
               gsl complex mul imag(
                  gsl complex add(
                     gsl complex add( gsl vector complex get(phi temp, dim-1),
                                      gsl vector complex get(phi temp, 1)),
                     gsl complex mul real( gsl vector complex get(phi temp, θ),
                                           potential[0] )).
                  -dt/2.0))):
   //periodic boundaries: i=dim-1
   gsl vector complex set(phi, dim-1, gsl complex add(
               gsl vector complex get(phi temp, dim-1),
               gsl complex mul imag(
                  gsl complex add(
                     gsl complex add( gsl vector complex get(phi temp, dim-2),
                                      gsl vector complex get(phi temp, 0)),
                     gsl complex mul real(gsl vector complex get(phi temp, dim-1),
                                           potential[dim-1] )),
                  -dt/2.0))):
  } else {
```

# Portability of your algorithm

How to run your algorithm?

- Single machine, single CPU
- Single machine, multiple CPU's (OpenMP, threads, ...)
- Multiple machines (MPI)
- GPU (Cuda, Thrust, OpenCL)

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Which data types are used by your algorithm?

- Build-in data types double, complex<double>
- Arbitrary precision types GMP, MPFR
- Vectorial data types float2d, float3d

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#### Theoretical solution

GoF Design Pattern: Strategy, also known as Policy

Gamma, Holm, Johnson, Vlissides: Design Patterns, Elements of Reuseable Object-Oriented Software, 1998.

# Numerical integration of ODEs

Find a numerical solution of an ODE an its initial value problem

$$\dot{x}=f(x,t)$$
,  $x(t=0)=x_0$ 

Example: Explicit Euler

$$x(t + \Delta t) = x(t) + \Delta t \ f(x(t), t) + \mathcal{O}(\Delta t^2)$$

General scheme of order s

$$x(t)\mapsto x(t+\Delta t)$$
 , or  $x(t+\Delta t)=\mathcal{F}_t x(t)+\mathcal{O}(\Delta t^{s+1})$ 

Solving ordinary differential equations in C++

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#### Modern C++

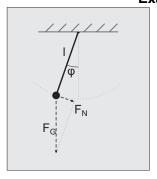
- Generic programming, functional programming
- Heavy use of the C++ template system
- Fast, easy-to-use and extendable.
- Container independent
- Portable

Introduction

2 Tutorial

- Technical details
- Discussion

### **Example – Pendulum**



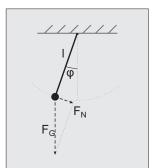
Pendulum - Newtons law ma = F

Acceleration 
$$a = I\ddot{\varphi}$$

Force

 $F = F_N = -mg \sin \varphi$ Result in an ode for the angle  $\ddot{\varphi} = -g/I\sin\varphi$ 

#### Example - Pendulum

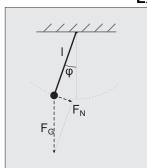


$$\ddot{arphi} = -g/I \sin arphi$$
Small angle  $\sin arphi pprox \phi$ 
Harmonic oscillator
 $\ddot{arphi} = -g/I arphi$ 

An analytic solution is known  $\varphi = A\cos\omega t + B\sin\omega t$ 

Amplitude *A* and *B* must be determined from initial conditions: 
$$\varphi(t=0) = \varphi_0, \ \dot{\varphi}(t=0) = \dot{\varphi}_0$$
 $B = \varphi_0, \ A = \dot{\varphi}_0/\omega$ 

#### **Example – Pendulum**

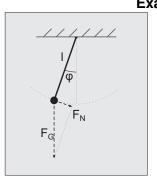


Full equation  $\ddot{\varphi}=g/I\sin\varphi$  has also analytic solution Jacobi elliptic function Lets enhance the ODE, add friction and external driving

 $\ddot{\varphi} = g/I \sin \varphi - \mu \dot{\varphi} + \varepsilon \sin \omega t$ 

No analytic solution is known. We need to solve this equation numerically.

### **Example - Pendulum**



 $\ddot{\varphi} = g/I \sin \varphi - \mu \dot{\varphi} + \varepsilon \sin \omega t$ Create a first order ODE  $x_1 = \varphi, x_2 = \dot{\varphi}$   $\dot{x_1} = x_2, \dot{x_2} = -g/I \sin x_1 - \mu x_2 + \mu x_3 + \mu x_4 + \mu x_5 + \mu x_5$ 

 $x_1 - x_2$ ,  $x_2 - g/r\sin x_1 - \mu x_2 + \varepsilon \sin \omega t$  $x_1$  and  $x_2$  are the state space variables.

Introduction

2 Tutorial

Technical details

Discussion

Introduction

2 Tutorial

Technical details

Discussion



# First example – Lorenz system

```
#include <iostream>
#include <boost/array.hpp>
#include <boost/numeric/odeint.hpp>
using namespace std;
using namespace boost::numeric::odeint;
const double sigma = 10.0;
const double R = 28.0;
const double b = 8.0 / 3.0;
typedef boost::array< double , 3 > state_type;
void lorenz ( const state type &x , state type &dxdt , double t )
    dxdt[0] = sigma * (x[1] - x[0]);
    dxdt[1] = R * x[0] - x[1] - x[0] * x[2];
    dxdt[2] = -b * x[2] + x[0] * x[1];
void write lorenz (const state type &x , const double t )
    cout << t << '\t' << x[0] << '\t' << x[1] << '\t' << x[2] << endl;
```

- The r.h.s. of the ODE is a simple function
- Observer

```
runge_kutta4< state_type > stepper;
```

```
runge_kutta4< state_type > stepper;
```

```
controlled_runge_kutta< runge_kutta_cash_karp54< state_type > > stepper;
```

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runge_kutta4< state_type > stepper;
```

```
controlled_runge_kutta< runge_kutta_cash_karp54< state_type > > stepper;
```

```
dense_output_runge_kutta< controlled_runge_kutta<
    runge_kutta_dopri5< state_type > > > stepper;
```

```
runge_kutta4< state_type > stepper;
```

```
controlled_runge_kutta< runge_kutta_cash_karp54< state_type > > stepper;
```

```
dense_output_runge_kutta< controlled_runge_kutta<
   runge_kutta_dopri5< state_type > > > stepper;
```

```
runge_kutta_dopri5< state_type > stepper;
make_dense_output( 1.0e-6 , 1.0e-6 , stepper ); // incomplete
```

#### Different steppers:

```
runge_kutta4< state_type > stepper;
```

```
controlled_runge_kutta< runge_kutta_cash_karp54< state_type > > stepper;
```

```
dense_output_runge_kutta< controlled_runge_kutta<
    runge_kutta_dopri5< state_type > > > stepper;
```

```
runge_kutta_dopri5< state_type > stepper;
make_dense_output( 1.0e-6 , 1.0e-6 , stepper ); // incomplete
```

### All together:

# Second example – Fermi-Pasta-Ulam lattice

$$\dot{q}_k = p_k$$

$$\dot{p}_k = -q_k^2 + \Delta q_k + \beta \{ (q_{k+1} - q_k)^3 - (q_k - q_{k-1})^3 \}$$

$$\Delta q_k = q_{k+1} - 2q_k + q_{k-1}$$

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$$\Delta q_k = q_{k+1} - 2q_k + q_{k-1}$$

#### State type consists of coordinates *q* and momentas *p*

```
typedef std::vector<double> vector_type;
vector_type q( 256 ) , p( 256 );
// initialize q.p
std::pair< state_type , state_type > state = std::make_pair( q , p );
```

# Second example – Fermi-Pasta-Ulam lattice

$$\dot{q}_{k} = p_{k} 
\dot{p}_{k} = -q_{k}^{2} + \Delta q_{k} + \beta \{ (q_{k+1} - q_{k})^{3} - (q_{k} - q_{k-1})^{3} \} 
\Delta q_{k} = q_{k+1} - 2q_{k} + q_{k-1}$$

### State type consists of coordinates *q* and momentas *p*

```
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vector_type q( 256 ) , p( 256 );
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```

### Hamiltonian system $\implies$ Symplectic solvers needed

```
symplectic_rkn_sb3a_mclachlan< vector_type > stepper;
```

## Fermi-Pasta-Ulam lattice continued

## Trivial first component $\dot{q}_k = p_k$

```
struct fpu {
   double m_beta;
   fpu(double beta) : m_beta(beta) { }

   void operator()(const vector_type &q, vector_type &dpdt) const {
        // ...
   }
};
```

## Fermi-Pasta-Ulam lattice continued

### Trivial first component $\dot{q}_k = p_k$

### All together

```
struct statistics_observer {
    void operator()( const state_type &x , double t ) const {
        // write the statistics
    }
};
int main(int argc, char **argv)
{
    vector_type q( 256 ) , p( 256 );
    // initialize q,p
    integrate_const( symplectic_rkn_sb3a_mclachlan< state_type >(), fpu(1.0),
        make_pair( q , p ), 0.0, 10.0, 0.01, statistics_observer() );
    return 0;
}
```

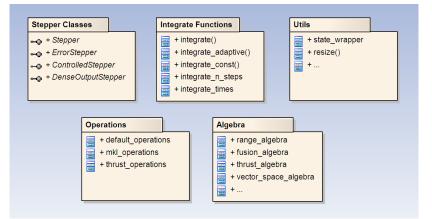
Introduction

2 Tutorial

Technical details

Discussion

### Structure of odeint



# Internals – Example Euler's method

User provides

$$y_i = f_i(x(t), t)$$

odeint provides

$$x_i(t + \Delta t) = x_i(t) + \Delta t \cdot y_i$$

(In general vector operations like  $z_i = a_1 x_{1,i} + a_2 x_{2,i} + \dots$ )

#### Instantiation

```
euler<state_type, value_type, deriv_type,
    time_type, algebra, operations > stepper;
```

All elements for container independence and portability are already included in this line!

$$y_i = f_i(x(t))$$
  
$$x_i(t + \Delta t) = x_i(t) + \Delta t \cdot y_i$$

```
euler<state_type, value_type, deriv_type,
   time_type, algebra, operations > stepper;
```

### Data types

- state\_type the type of x
- value\_type the basic numeric type, e.g. double
- deriv\_type the type of y
- time\_type the type of t,  $\Delta t$

$$y_i = f_i(x(t))$$
  
$$x_i(t + \Delta t) = x_i(t) + \Delta t \cdot y_i$$

```
euler<state_type, value_type, deriv_type,
    time_type, algebra, operations > stepper;
```

#### Algebra policies, perform the iteration

Algebra must be a class with public methods

- for\_each1(x,op) Performs  $op(x_i)$  for all i
- for\_each2(x1,x2,op) Performs  $op(x1_i,x2_i)$  for all i
- ...

$$y_i = f_i(x(t))$$
  
$$x_i(t + \Delta t) = x_i(t) + \Delta t \cdot y_i$$

```
euler<state_type, value_type, deriv_type,
    time_type, algebra, operations > stepper;
```

#### Operations do the basic computation

Operations must be a class with the public classes (functors)

- scale\_sum1 Calculates  $x = a1 \cdot y1$
- scale\_sum2 Calculates  $x = a1 \cdot y1 + a2 \cdot y2$
- ...

$$y_i = f_i(x(t))$$
  
$$x_i(t + \Delta t) = x_i(t) + \Delta t \cdot y_i$$

```
euler<state_type, value_type, deriv_type,
    time_type, algebra, operations > stepper;
```

#### All together

```
m_algebra.for_each3(xnew ,xold, y ,
    operations_type::scale_sum2<value_type,time_type>(1.0,dt));
```

## Stepper concepts

### Concepts

"... In generic programming, a concept is a description of supported operations on a type..."

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"... In generic programming, a concept is a description of supported operations on a type..."

#### odeint provides

Stepper concept

```
stepper.do_step(sys,x,t,dt);
```

ErrorStepper concept

```
stepper.do_step(sys,x,t,dt,xerr);
```

ControlledStepper concept

```
stepper.try_step(sys,x,t,dt);
```

DenseOutputStepper concept

```
stepper.do_step(sys);
stepper.calc_state(t,x);
```

# Supported methods

Method Euler Runge-Kutta 4 Runge-Kutta Cash-Karp Runge-Kutta Fehlberg Runge-Kutta Dormand-Prince	Class name euler runge_kutta4 runge_kutta_cash_karp54 runge_kutta_runge_fehlberg78 runge_kutta_dopri5	Concept SD S SE SE SED
Runge-Kutta controller	controlled_runge_kutta	C
Runge-Kutta dense output	dense_output_runge_kutta	D
Symplectic Euler	symplectic_euler	S
Symplectic RKN	symplectic_rkn_sb3a_mclachlan	S
Rosenbrock 4	rosenbrock4	ECD
Implicit Euler	implicit_euler	S
Adams-Bashforth-Moulton Bulirsch-Stoer	adams_bashforth_moulton bulirsch_stoer	S CD

S – fulfills stepper concept

E – fulfills error stepper concept

C – fulfills controlled stepper concept

D – fulfills dense output stepper concept

# Integrate functions

- integrate\_const
- integrate\_adaptive
- integrate\_times
- integrate\_n\_steps

Perform many steps, use all features of the underlying method

# Integrate functions

- integrate\_const
- integrate\_adaptive
- integrate\_times
- integrate\_n\_steps

Perform many steps, use all features of the underlying method

An additional observer can be called

```
integrate_const(stepper, sys, x, t_start,
t_end, dt, obs);
```

### More internals

- Header-only, no linking → powerful compiler optimization
- Memory allocation is managed internally
- No virtual inheritance, no virtual functions are called
- Different container types are supported, for example
  - STL containers (vector, list, map, tr1::array)
  - MTL4 matrix types, blitz++ arrays, Boost.Ublas matrix types
  - thrust::device\_vector
  - Fancy types, like Boost.Units
  - ANY type you like
- Explicit Runge-Kutta-steppers are implemented with a new template-metaprogramming method
- Different operations and algebras are supported
  - MKL
  - Thrust
  - gsl

### **ODEs on GPUs**

Graphical processing units (GPUs) are able to perform up to 10<sup>6</sup> operations at once in parallel

#### Frameworks

- CUDA from NVIDIA
- OpenCL
- Thrust a STL-like library for CUDA and OpenMP

#### Applications:

- Parameter studies
- Large systems, like ensembles or one- or two dimensional lattices
- Discretizations of PDEs

### odeint supports CUDA, through Thrust

## Example: Parameter study of the Lorenz system

```
typedef thrust::device vector<double> state type;
typedef runge kutta4<state type ,value type ,state type ,value type ,
     thrust algebra .thrust operations > stepper type;
struct lorenz system {
    lorenz_system(size_t N ,const state_type &beta)
    : m N(N) , m beta(beta) {}
    void operator()( const state_type &x , state_type &dxdt , double t ){
        // ..
    size t m N;
    const state type &m beta;
};
int main ( int arc , char* argv[] )
    const size t N = 1024:
    vector<value type> beta host(N);
    for( size_t i=0 ; i<N ; ++i )
        beta host[i] = 56.0 + \text{value type}(i) * (56.0) / \text{value type}(N - 1)
              );
    state type beta = beta_host;
    state_type x(3 * N, 10.0);
    integrate const( stepper type() , lorenz(N,beta) , x , 0.0 , 10.0 , 0.01
         );
    return 0;
```

### Conclusion

- odeint provides a fast, flexible and easy-to-use C++ library for numerical integration of ODEs.
- Its container independence is a large advantage over existing libraries.
- Portable
- Generic programming is the main programming technique.

### Outlook

- Submission to the boost libraries
- Dynamical system classes for easy implementation of interacting dynamical systems
- More methods: implicit methods and multistep methods.
- Implementation of the Taylor series method

```
taylor_fixed_order< 25 , 3 > taylor_type stepper;

stepper.do_step(
   fusion::make_vector
   (
       sigma * ( arg2 - arg1 ) ,
       R * arg1 - arg2 - arg1 * arg3 ,
       arg1 * arg2 - b * arg3
   ) , x , t , dt );
```

### Resources

#### Download and documentation

odeint.com

### An article about the used techniques exists at

http://www.codeproject.com/KB/recipes/odeint-v2.aspx

### Development

https://github.com/headmyshoulder/odeint-v2

#### Contributions and feedback

are highly welcome