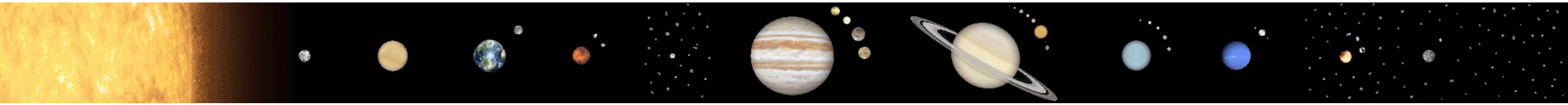
PSE Molekulardynamik
Sheet 1: First steps towards a molecular dynamics' simulation



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Task 1 & 2: Set up

- ParaView (5.9)
- CMake (3.27.4)
- Doxygen (1.10.0)
- Clang (16.0.6)
- Make (4.3)
- Graphviz (2.42.2)
- Libxerces (3.2.4)
- Clang tidy

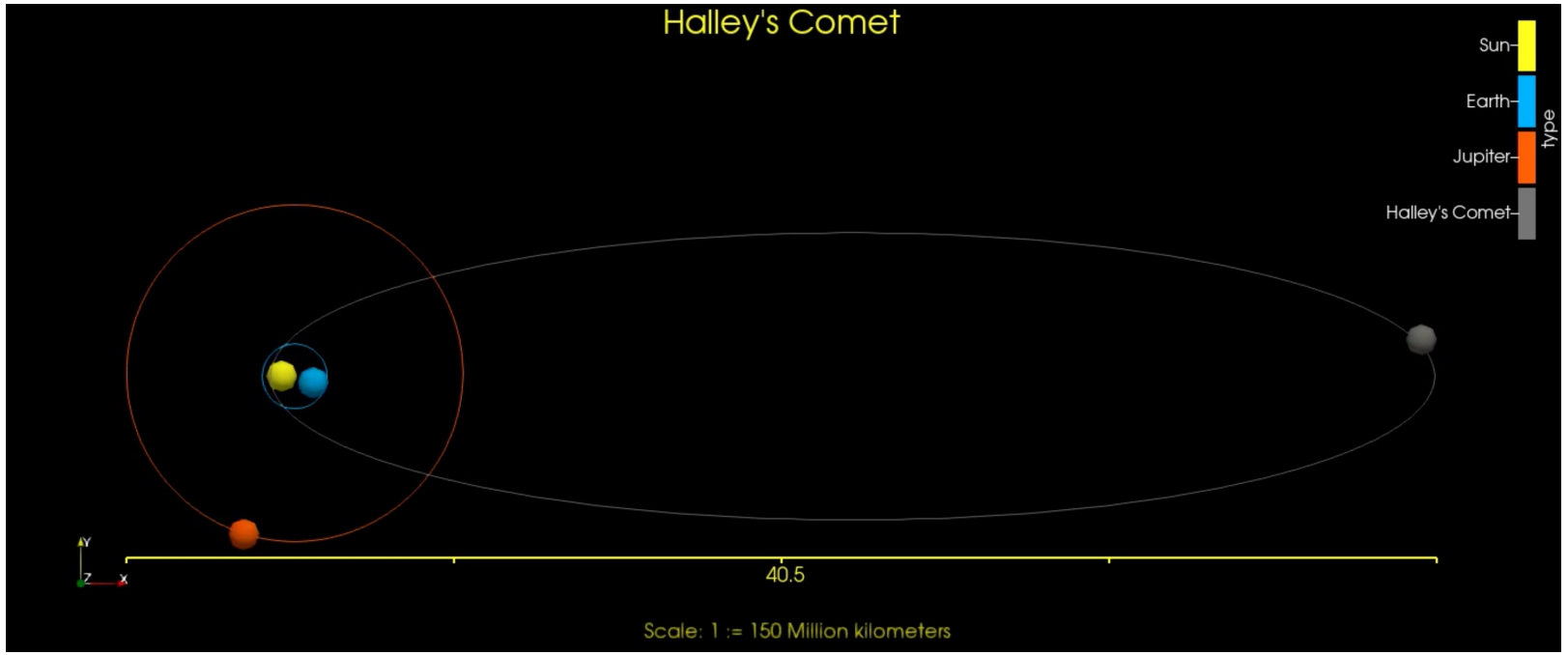
Task 3: Completion of program frame

- Implementation of **Force calculation** was straight forward
  - used L2 Norm from ArrayUtils.h
- Checked implementation by **running a simulation** in ParaView
  - adjusted output file format from .xyt to .vtu
- Boost for **argument parsing** in the command line

Task 4: Halley’s Comet

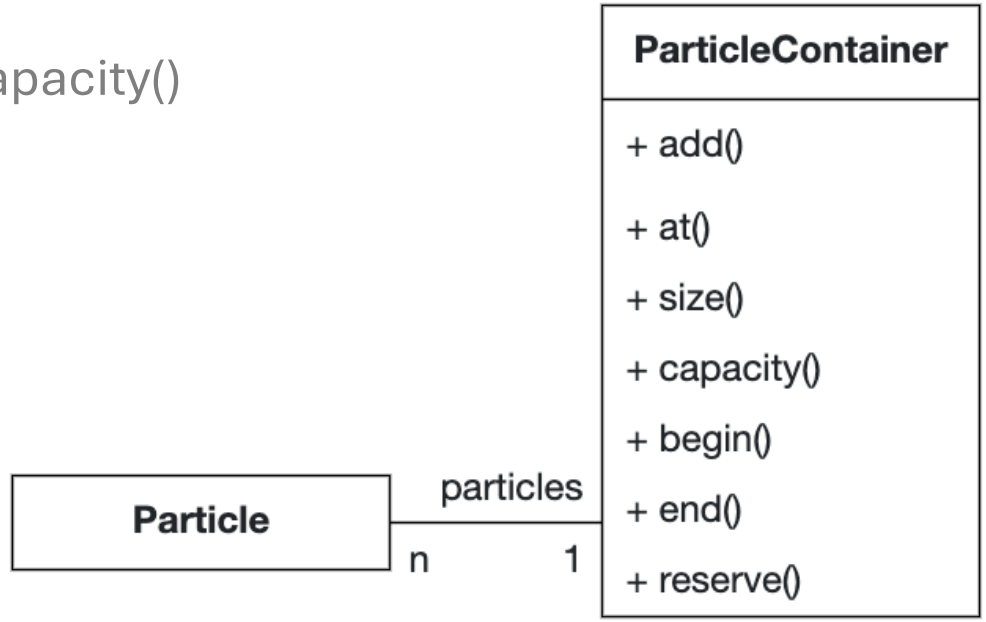
Identification of celestial bodies:

- **Sun:** most mass
- **Comet:** least mass, unconventional trajectory
- **Earth and Jupiter:** earth is closer to sun than Jupiter



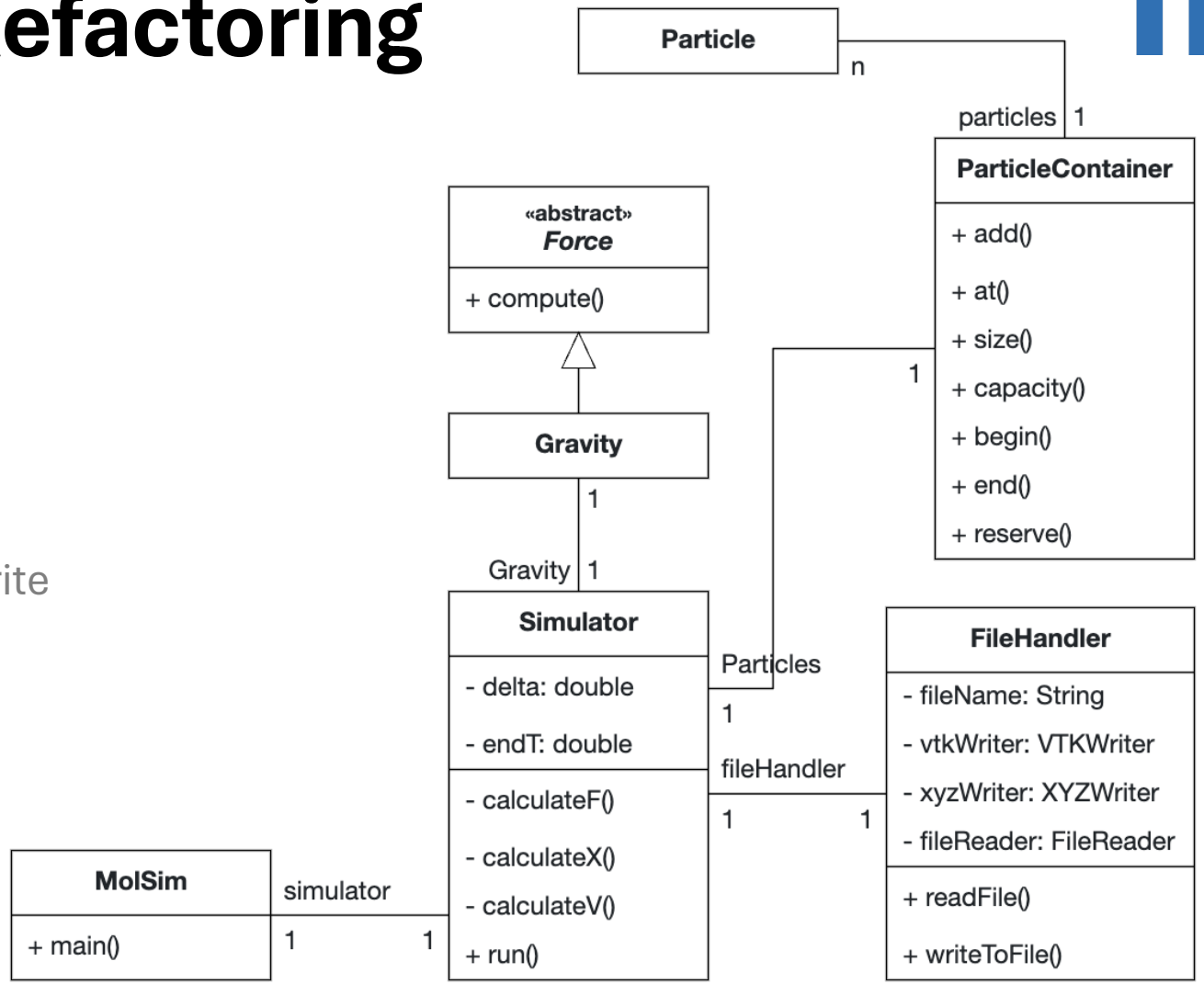
Task 5: Particle Container

- Store particles using **std::vector**
  - objects stored **consecutively** for better **performance**
  - but still **dynamic** on the other hand
  - Implementation of add(), at(), size(), capacity()
- **Iterator** pattern
  - Iterate over pairs with **nested for loop**
  - Implementation of begin(), end()



Task 5: Further Refactoring

- **Strategy pattern** for forces acting on Particles
  - abstract class Force.h as **parent** for all forces to come
  - **gravity** being the first (implementation of task 3)
- Interface for **file handling**
  - **FileHandler** encapsulates write and read classes
- **Simulator** class instead of comp-lex main method



Thank you for listening!