

#### NUMERICAL PRECISION

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#### Let's talk about precision

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## Know your friend / enemy

- IEEE floating-point in a nutshell:
  - ± 1.0100100111011 2<sup>±010011101</sup> mantissa
- Consequences:
  - Base 2 → Even good old 0.1 isn't exact!
  - Precision relative to exponent / order of magnitude
  - Unbounded loss of accuracy on subtract / add
  - Very small / large numbers need care

# **Comparing FP numbers**

- How to tell if val ≈ ref?
  - FP precision is relative  $\rightarrow$  Relative comparison often best
  - Typical algorithm looks like  $|val ref| < tol \cdot |ref|$
  - Nice side-effect: tolerance is (mostly) data-agnostic
- Two limits of relative comparisons
  - Orders of magnitude may matter (e.g. spatial tolerances)
  - Breaks down when reference is close to zero

# Some choice can be good

- When in doubt, start with relative comparisons
- If they prove inadequate, consider other algs...
  - Absolute comparisons : |val ref| < tol
  - « Small enough » : |val| < tol</p>
  - « Close or small » : relative unless val & ref are both small
  - L2 norm of difference of matrices vs ref matrix, etc.

# Too much choice will kill you

- FP test assertions currently used in ACTS :
  - BOOST\_CHECK\_CLOSE(val, ref, tol)
  - BOOST\_CHECK\_CLOSE\_FRACTION(val, ref, tol)
  - BOOST\_CHECK\_SMALL(val, tol)
  - BOOST\_TEST(val == ref[, tol]
  - BOOST\_CHECK(val.isApprox(ref[, tol]))
  - checkCloseXyz(val, ref)
  - STL container element-wise comparison (test-specific)

## **Consistency matters**

- The previous assertions disagree on many things:
  - Are relative tolerances given as fractions? Percentages?
  - Can I compare floats with integers? Doubles?
  - Does it work with scalars? Eigen types? STL containers?
  - Is there a default tolerance? A hidden global one?
  - What happens when a value/reference is near zero?
  - Are matrices compared element-wise or by L2 norm?
  - How good is the error reporting?

# Trying to improve upon this

- Key goal: Assertions should be easy to understand
  - Follow typical & shared conventions
  - Inputs are explicit (nothing global, nothing hardcoded)
  - Simple, general-purpose and predictable logic
- Some flexibility on comparison algs, input types
- Report errors as clearly as possible
- My attempt at resolving this: acts-core!490

#### One remaining problem

/root/acts-core/Tests/Integration/PropagationTestHelper.hpp(527):
error: in "covariance\_transport\_disc\_disc\_/\_45":
check Acts::Test::checkCloseOrSmall((calculated\_cov),
 (obtained\_cov), (reltol), (1e-4)) has failed. [...]

The	failure	occured during	a matrix cor	mparison, wher	e the value was
	35447.7	31.4111	-1.80979	59.5127	0.291849
	31.4111	25761.4	53.0901	1.53086	-8.93186
	-1.80979	53.0901	0.112616	3.72723e-06	-0.0356915
	59.5127	1.53086	3.72723e-06	0.1	-1.98435e-11
	0.291849	-8.93186	-0.0356915	-1.98435e-11	0.1
and the reference was					
	35448	20.9458	-1.8162	59.5128	0.291879
	20.9458	25864.9	53.1939	1.52074	-8.93245
	-1.8162	53.1939	0.112616	1.91157e-06	-0.0356914
	59.5128	1.52074	1.91157e-06	0.1	0
	0.291879	-8.93245	-0.0356914	Θ	0.1

#### Help wanted!

- Seeing this now because we used isApprox() before
  - isApprox() based on L2 norm: ||val-ref|| < tol · ||ref||</p>
  - Comparison dominated by large diagonal terms
  - But... does L2 norm make sense for covariance?
- Question: how should I handle this issue?
  - Is this difference physically significant?
  - Should I consider it to be a propagator bug?

# Beyond that: single precision experiment

- Step 1: Evaluate SP tolerances of ACTS code 
   → OK
  - acts-core!491: Making tests pass under Verrou emulation
  - Affected by previous issue, otherwise looking good...
- - Need review from someone who knows the physics!
- Step 3: Fix the unacceptable part → TBD
  - Look out for easy « precision bottlenecks »
  - Move what we can to SP, keep rest in double precision

