

Fluka and Geant4

Using Fluka for CALICE

- ▶ Motivation
- ▶ Method
- ▶ Initial results
- ▶ Future

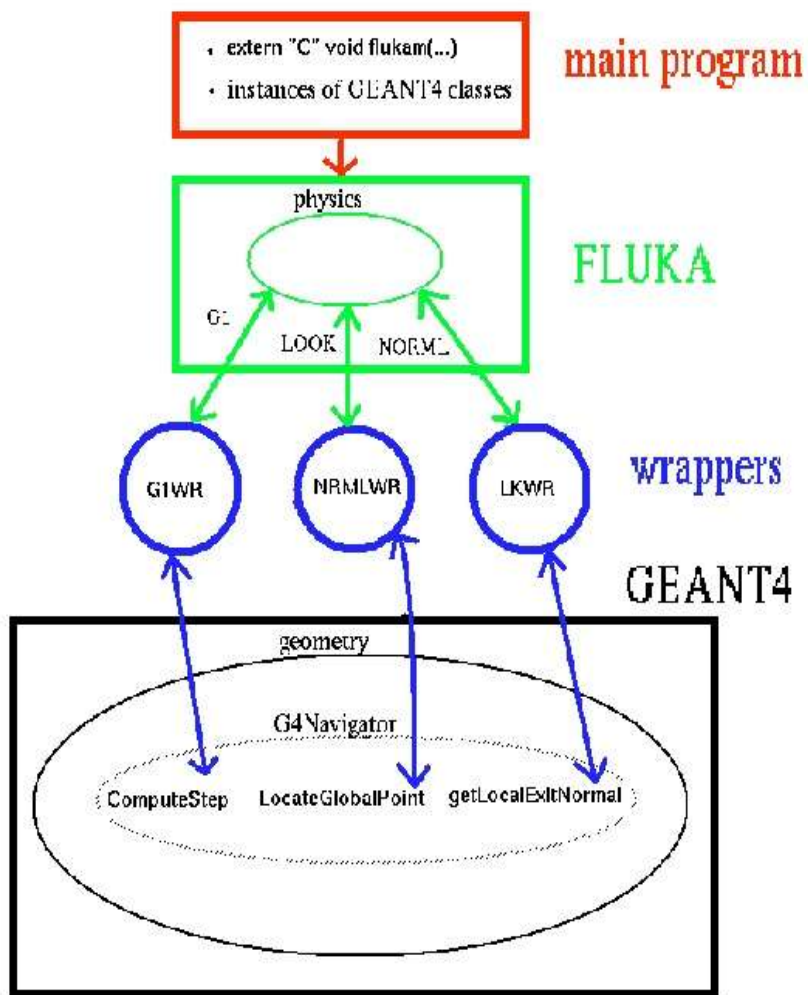
David Ward (Cambridge) for

Nigel Watson (CCRLC-RAL & Birmingham)

Motivation

- Systematic comparison of Mokka and Fluka physics modelling of CALICE test beam
- Particularly interesting for hadronic interactions
 - ▶ See G3/G4 studies (DRW, George Mavromanolakis)
- Wish to...
 - ▶ Test new Mokka detector models
 - ▶ Investigate full TDR type geometry
 - ▶ Avoid coding each geometry directly in Fluka
 - ⇒ error prone, may introduce non-physics differences
- Issues
 - ▶ Fluka geometry defined by data cards
 - ▶ Only limited geometrical structures supported
 - ▶ Repeated structures at 1 level only

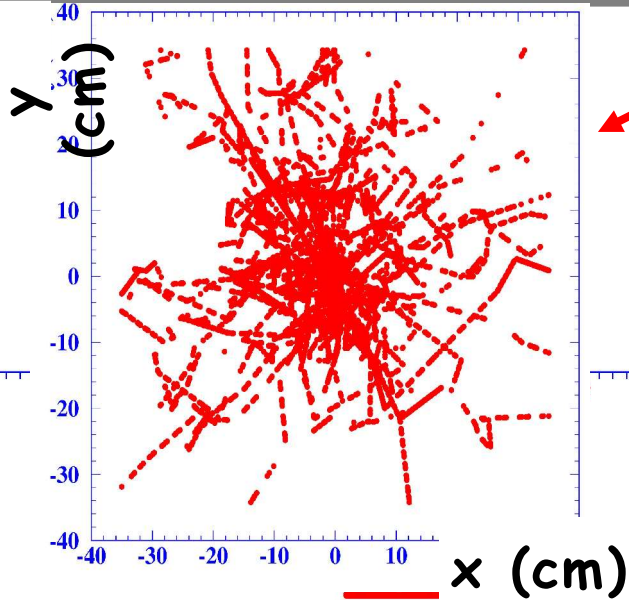
"Flugg" Package (P. Sala et al)



- Geometry & physics decoupled in G4 and Fluka
- Wrappers for f77/C++
- Fluka authors' comparisons of G4 with Flugg (FLUKa+G4 Geometry)
 - ▶ Simple detectors, identical results
 - ▶ Complex T36 calorimeter: 81 layers Pb (10mm)-scint.(2.5mm) Consistent results
- My first test
 - ▶ Use T36 calorimeter as above

[From ATL-SOFT-98-039]

Flugg Issues

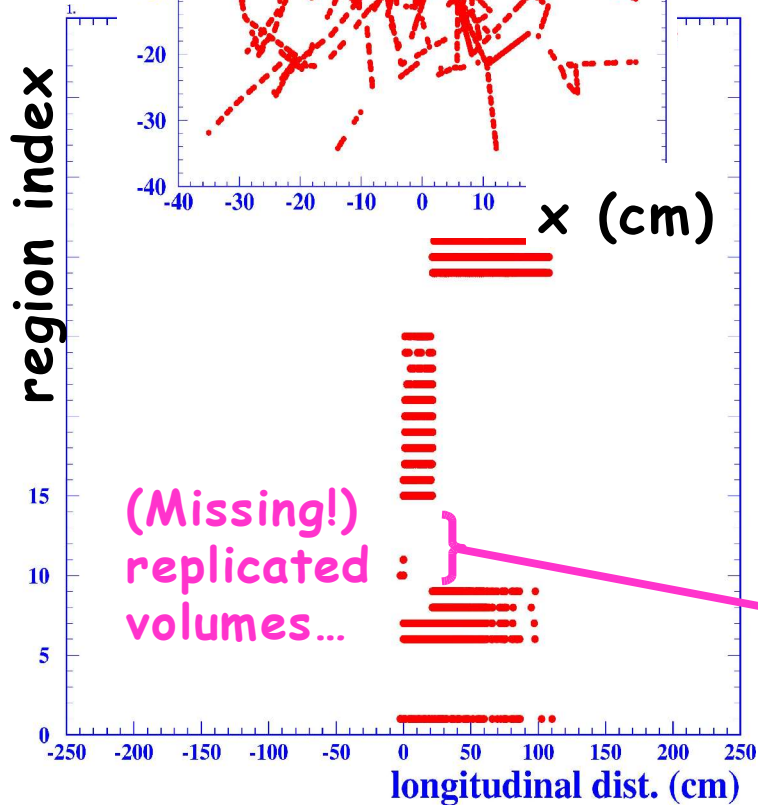


■ Transverse response of T36 calo. to 10 GeV π^- in flugg

■ User control available:

▶ at every tracking step, via rudimentary drawing routine (slow)

▶ at every energy deposition event



Note

■ For G4 replicated or parametrised volumes (correspond to Fluka "lattice volumes")

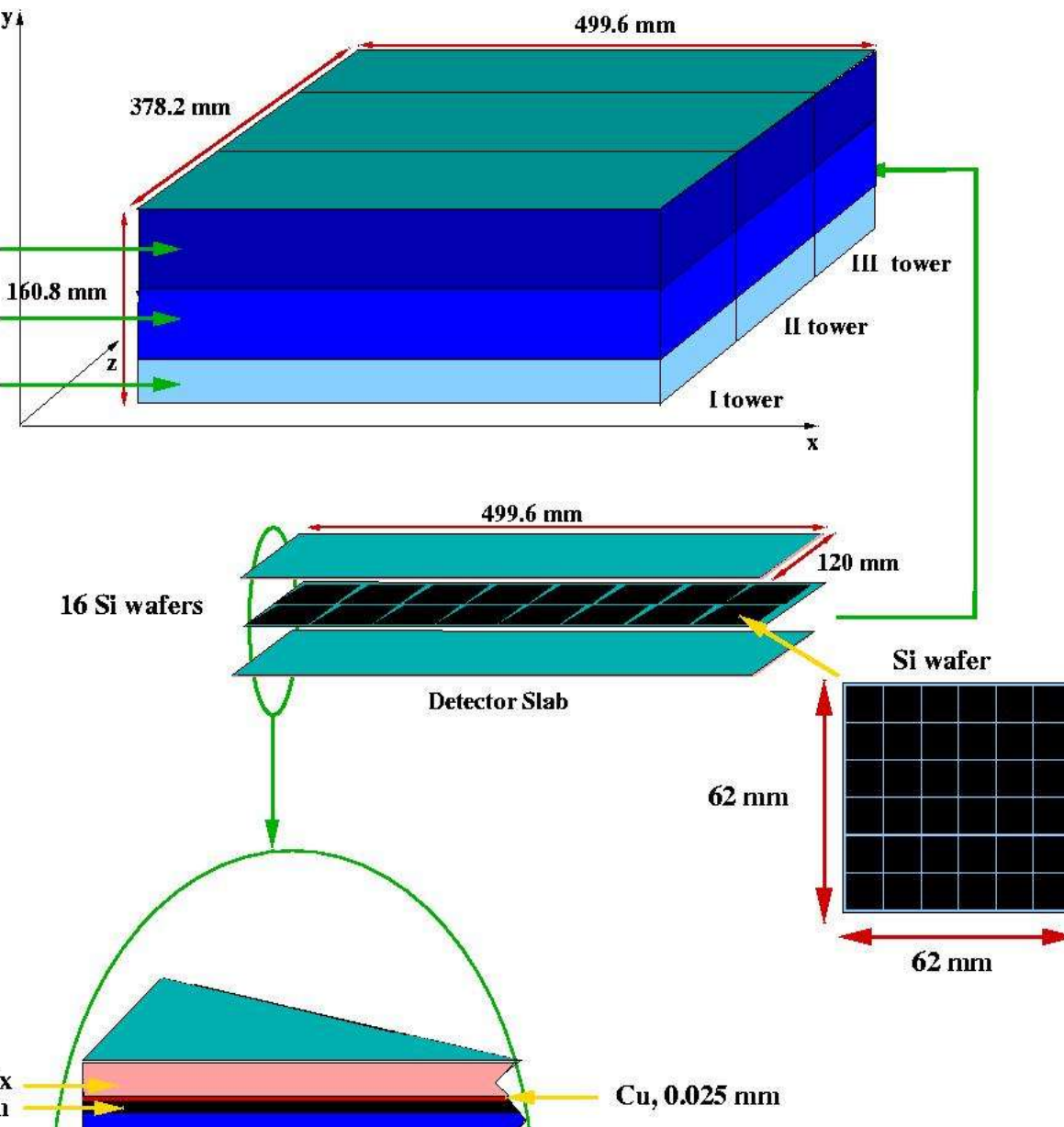
▶ Region index is degenerate

▶ Boundary crossings sometimes

4 not detected

Montpellier, 14-Nov-

Volume Ambiguity



- fluka 'sees' 3x32 Si volumes
- id for wafers degenerate
 - ▶ in z (x3 towers)
 - ▶ in y within a stack of 5 detector slabs (10 Si layers)

Current Status

- Mokka running within flugg/Fluka framework
 - ▶ Using Mokka-01-05 + Geant4.5.0.p01 + clhep1.8.0 + gcc3.2
 - ▶ Flugg05 (Jan. 2003)
 - ▶ Fluka 2002.4 (May 2003)
- Procedure: start from Mokka release and **delete**:
 - ▶ all classes **except** for detector construction, detector parametrisation, magnetic field construction
 - ▶ corresponding #include, variable, class definitions in .cc/.hh
 - ▶ anything related to G4RunManager, DetectorMessenger
 - ▶ code where SensitiveDetector is set
 - ▶ interactive code, visualisation, etc.
- Validation
 - ▶ Minimal debugging tools in flugg, e.g. P55 prototype geometry

Flugg Operation

Two pass operation

■ One-time initialisation

- ▶ Read G4 geometry/material definitions
- ▶ Generate fluka input cards
 - ⇒ Material/compound definitions
 - ⇒ Material to volume assignments

■ Subsequent runs with a given geometry model

- ▶ Use generated Fluka cards
- ▶ Tracking within G4 geometry
- ▶ Physics processes from Fluka

First pass, G4 → Fluka conversion

Connecting to the database models00

Building sub_detector P66WNominal, geometry db P66WNominal, driver proto01:

Ecal prototype driver with W ideal thickness (reference)

Connecting to the database P66WNominal

proto01: proto size is (499.600000,160.800000,3

proto01: placing prototype at (0.000000,236.0000

Sub_detector P66WNominal DONE!

Building sub_detector SinglehcalFeRPC1, geometry

Single module Hcal Fe & RPC as prototype

Connecting to the database SinglehcalFeRPC1

The sensitive model in Hcal chambers is RPC1

Iron is the radiator material being placed.

Sub_detector SinglehcalFeRPC1 DONE!

tRadlen() = 89867.3 mm

Styropor->GetRadlen() = 17518.3 mm

C->GetRadlen() = 188.496 mm

CGAGeometryManager starting the detector constr

Asking for the model ProtoEcalHcalRPC:

Building Proto release 01

total_W_layers = 30

MixDensite = 2.15747 g/cm3

Mix->GetRadlen() = 75.0202 mm

Proto done.

generates fluka input deck

Building Hcal...

Detector construction done.

* G4PhysicalVolumeStore (0x401b5288) has 2424 volume

* Storing information...

+ Tungsten: dens. = 19.3g/cm3, nElem = 1

Stored as TUNGSTEN

+ TungstenModified: dens. = 11g/cm3, nEl

Stored as TUNGST02

+ Copper: dens. = 8.96g/cm3, nElem = 1

Stored as COPPER

+ Silicium: dens. = 2.33g/cm3, nElem = 1

Stored as SILICIUM

+ SiVXD: dens. = 8.72g/cm3, nElem = 1

Stored as SIVXD

+ Iron: dens. = 7.87g/cm3, nElem = 1

Stored as IRON

+ Aluminum: dens. = 2.7g/cm3, nElem = 1

Stored as TETRAFLU

Stored as RPCGAS1

Stored as GRAPHITE

+ Mix: dens. = 2.15747g/cm3, nElem = 9

Stored as MIX

----- ... -----

* Printing FLUKA materials...

* Printing FLUKA compounds...

* G4PhysicalVolumeStore (0x401b5288) has 2424 volume

* Printing ASSIGNMAT...

* Printing Magnetic Field...

No field found...

*** Entering UsrIni.f!! ***

*** Entering HistIn.f!! ***

Operation

C4 volume index

material to region assignments

material definitions

& etc

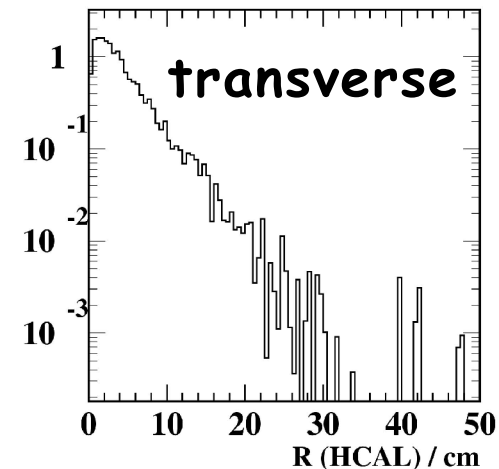
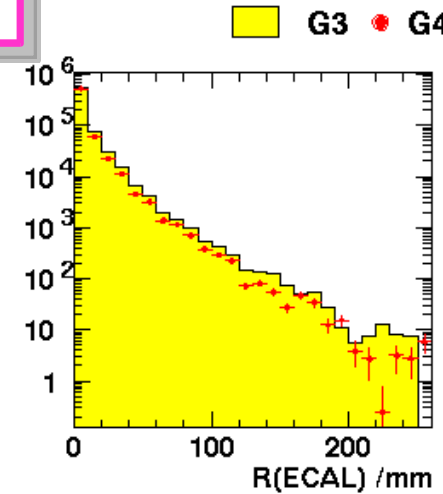
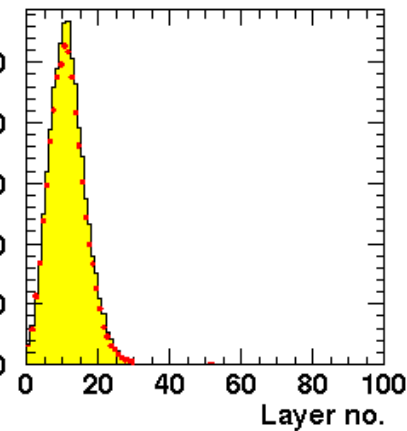
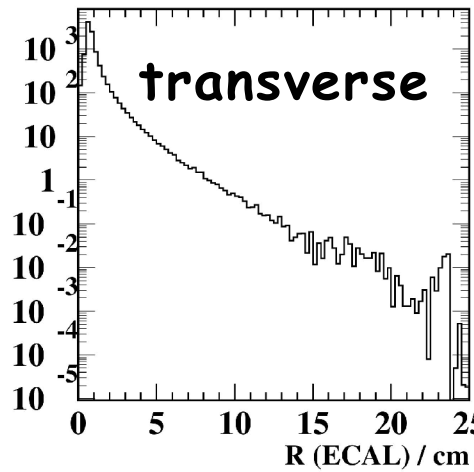
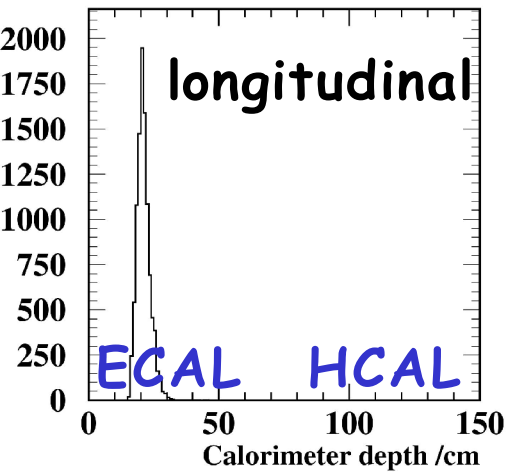
& etc

C4 volume index		material definitions				material to region assignments	
ID	Region	Material	Vol	Density	Mass	Region	Assignment
1	WorldPhysical						
2	SensWafferPhys	MATERIAL TUNGSTEN	74.0	183.840	1.930e+01	3.0	ASSIGNMAT 12.0
3	DeadWBlock	MATERIAL TUNGST02	74.0	183.840	1.100e+01	4.0	ASSIGNMAT 6.0
4	DeadWBlock	LOW-MAT TUNGSTEN	4.0				ASSIGNMAT 3.0
5	DeadWBlock	MATERIAL COPPER	29.0	63.546	8.960e+00	5.0	ASSIGNMAT 3.0
6	DeadWBlock	MATERIAL SILICIUM	14.0	28.090	2.330e+00	6.0	ASSIGNMAT 21.0
7	DeadWBlock	MATERIAL SIVXD	14.0	28.090	8.720e+00	7.0	ASSIGNMAT 21.0
9	DeadWBlock	MATERIAL IRON	26.0	55.850	7.870e+00	8.0	ASSIGNMAT 1.0
10	SlabWBlock	MATERIAL ALUMINUM	13.0	26.980	2.700e+00	9.0	
11	SlabWBlock	MATERIAL BERYLLIU	4.0	9.012	1.848e+00	10.0	
		MATERIAL ARGON	18.0	39.950	1.780e-03	11.0	
2417	EndCapChamberPhy	MATERIAL			1.290e-03	12.0	AIR
2418	EndCapChamberPhys	NITROGEN	7.0	14.010	9.990e-01	13.0	
2419	EndCapChamberPhys	OXIGEN	8.0	16.000	9.990e-01	14.0	
2420	EndCapChamberPhys	BEAM_			1.000e-05	15.0	
2421	EndCapChamberPhys	QUARTZ			2.200e+00	16.0	
		MATERIAL	14.0	28.090	9.990e-01	17.0	

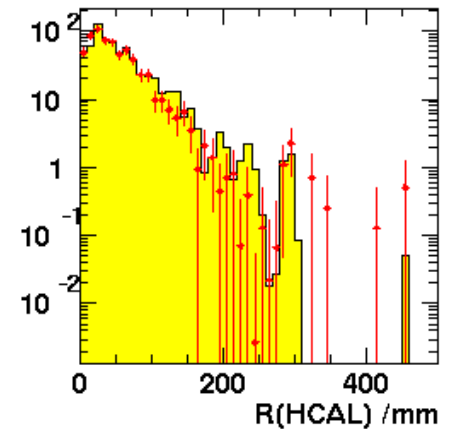
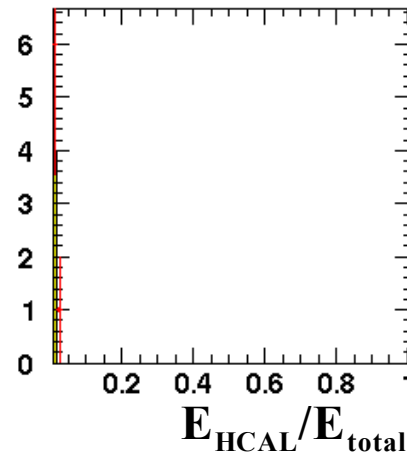
Fluka with G3/G4

ProtoEcalHcalRPC, 5 GeV e⁻

5 GeV e⁻



In these plots, fluka has energy deposited in all material, not just active layers



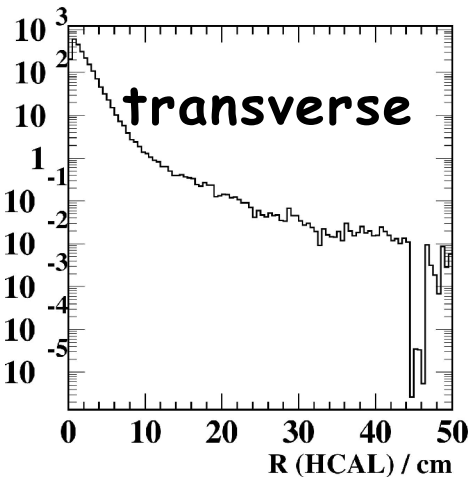
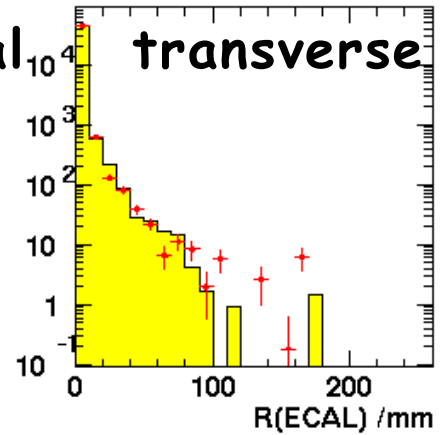
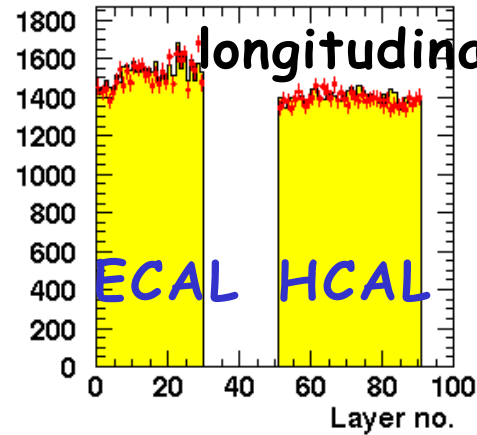
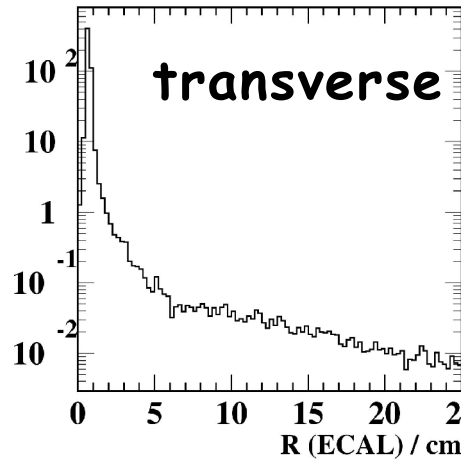
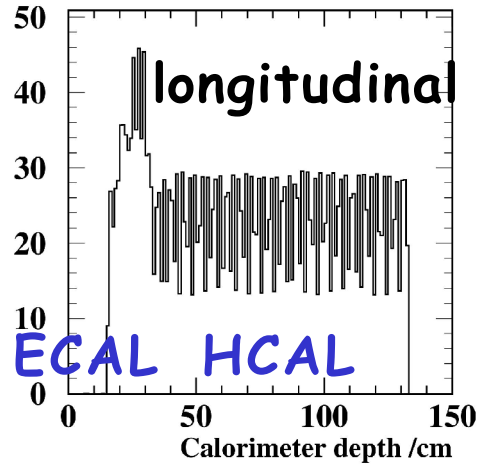
[G3/G4 plots from DRW]

Fluka with G3/G4

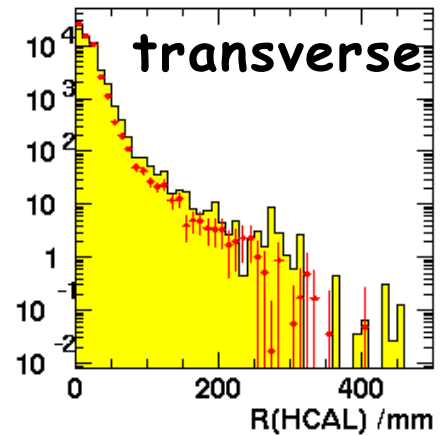
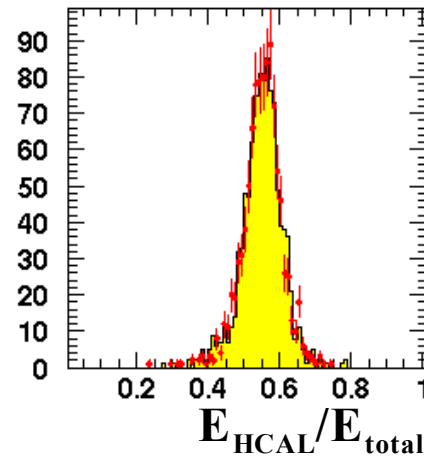
ProtoEcalHcalRPC, 5 GeV μ^-

5 GeV μ^-

■ G3 ● G4



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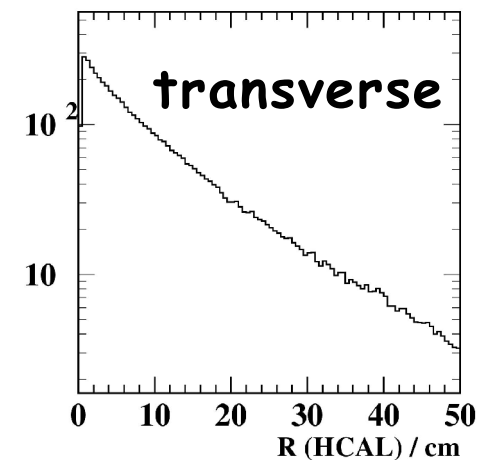
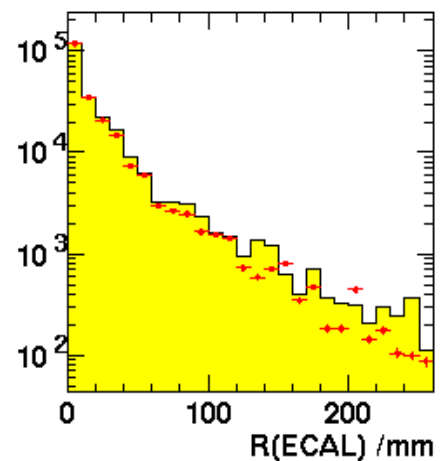
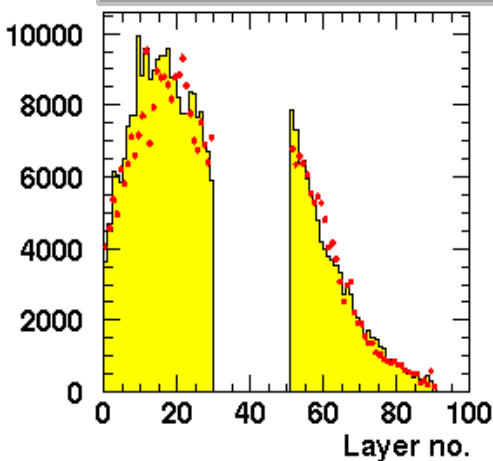
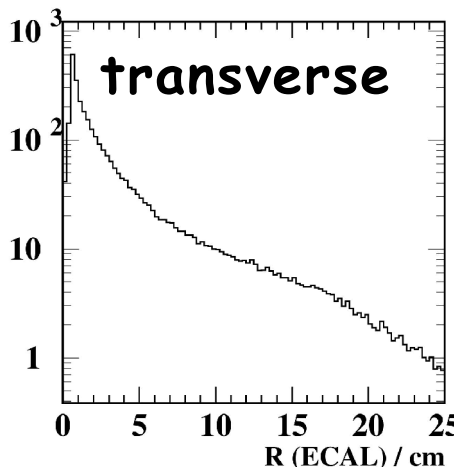
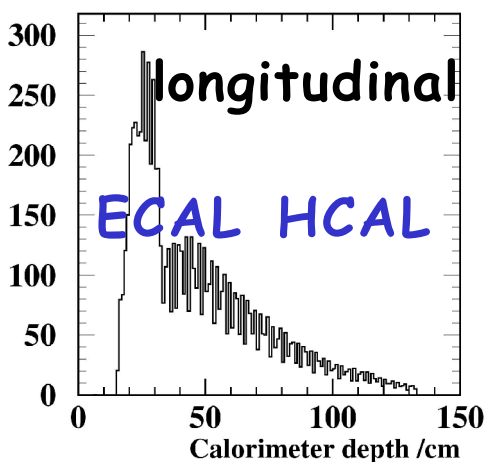
[G3/G4 plots from DRW]

Fluka with G3/G4

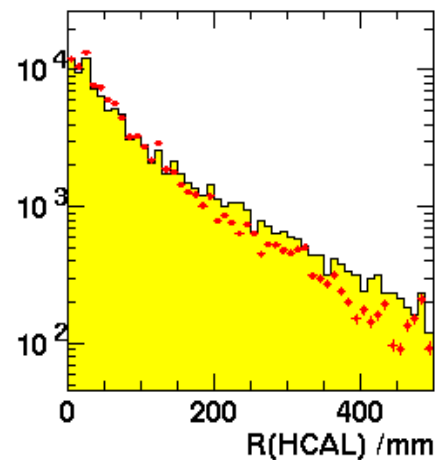
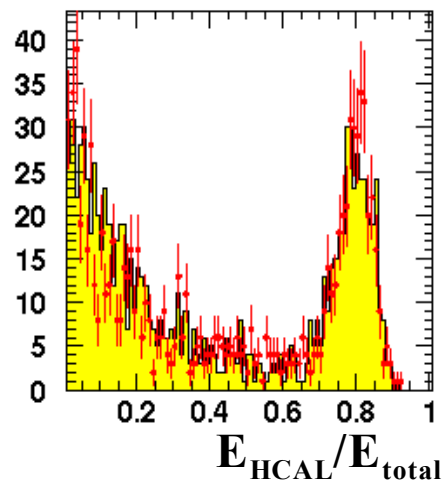
ProtoEcalHcalRPC, 5 GeV π^-

5 GeV π^-

■ G3 ● G4



In these plots, fluka has energy deposited in all material, not just active layers



Ongoing Work

- Restrict study to energy deposited in active layers
- Improve reliability for larger samples
 - ▶ ~understood technical issue
- Review energy thresholds/step size in Fluka
 - ▶ default min. K.E. > 100 keV
 - ▶ neutrons, 19.6 MeV
 - ▶ energy e/γ > 500 keV (??)
 - ▶ low energy neutron cross-sections
- Compare systematically with G3/G4 results,
 - ▶ Same initial conditions
 - ▶ Thresholds, mip normalisation, etc.
 - ▶ Adopt same output format as DRW/GM

Summary

Identified “easy” way of comparing G4/Fluka

- ▶ Alternative to deprecated G-Fluka
- ▶ Preferable to “standalone” Fluka as more efficient for variations in geometry

Integration with Mokka geometry classes

- ▶ Need to feed changes back to Mokka developers

Will be useful input when devising test beam programme/strategy