

# Status of event generation and Status of SGV

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ILD SW pre-meeting, Fukuoka, May, 2012

# Outline

- 1 Generator status
- 2 SGV
  - Calorimeter simulation
- 3 Summary and Outlook

# Generator status: where we are

- After receiving the final beam-parameters at Christmas, the DBD bench-mark generation is soon complete:
- $t\bar{t}h$  is done.
- Beam-strahlung background is done, both at 1 TeV and 500 GeV.
- $WW$  with it's backgrounds is done.
- $\nu\nu h$  with it's backgrounds is done, but not yet on the grid.
- $\gamma\gamma$  backgrounds (both low and high  $p_T$ ) are done, but not yet on the grid.
- “mini-jets” will be produced at SLAC by the end of the week, and on the grid soon after.
- $t\bar{t}$  at 500 GeV (and other 6 fermions) is lagging behind, but should be done in a month
- 2- and 4-fermions at 500 GeV can be done in a week, once the beam-spectrum files for the new parameters are in place.

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## File locations:

Sub-directories of:

lfn:/grid/ilc/prod/ilc/mc-dbd/generated/1000-B1b\_ws/

lfn:/grid/ilc/prod/ilc/mc-dbd/generated/xxx-TDR\_ws/

(xxx=250, 350, 500)

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## Information on samples:

<http://ilcsoft.desy.de/dbd/generated/>

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# Generator status

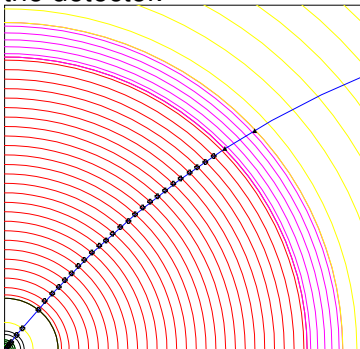
This was just the bottom line !!

See the **talk tomorrow** in the “Physics analysis” slot at **14:00** for the details.

# SGV: How tracking works

SGV is a machine to calculate covariance matrices

**Tracking:** Follow track-helix through the detector.

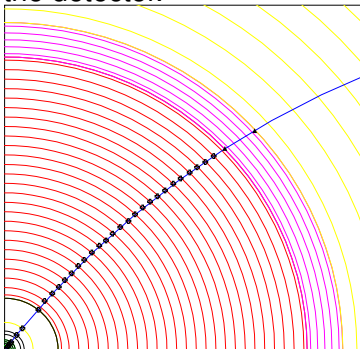


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- Smear perigee parameters (Choleski decomposition: takes all correlations into account)
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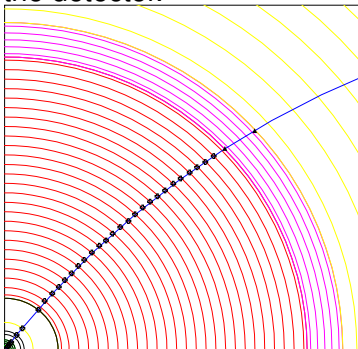


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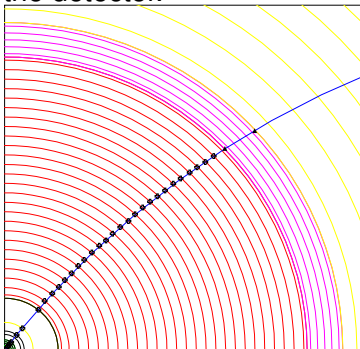


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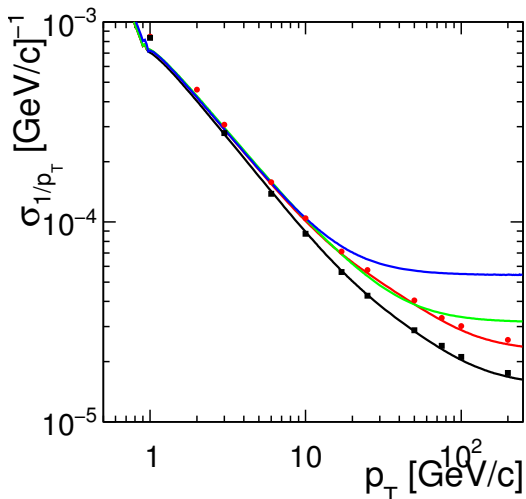
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## SGV and FullSim LDC/ILD: momentum resolution

Lines: SGV, dots: Mokka+Marlin



# SGV: How the rest works

## Calorimeters:

- Follow particle to intersection with calorimeters. Simulate:
  - Response type: MIP, EM-shower, hadronic shower, below threshold, etc.
  - Simulate response from parameters.

## Other stuff:

- EM-interactions in detector material simulated
- Plug-ins for particle identification, track-finding efficiencies,...



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# SGV: Technicalities

## Features:

- Written in **Fortran 95**.
- 20 000 lines + 10 000 lines of comments.
- Some **CERNLIB** dependence.
- Re-write of **battle-tested** f77 SGV 2-series (LEP, Tesla, LOI, ...)
- **Managed in SVN**. Install script included, much ameliorated in the HEAD version.
- Callable **PYTHIA**, **Whizard** or input from **PYJETS** or **stdhep**.
- Output of **generated event** to PYJETS or stdhep.
- **samples** subdirectory with READMEs, steering and code.
- **output LCIO DST**.

# Installing SGV

Do

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svn export https://svnsrv.desy.de/public/sgv/trunk/ sgv/
```

Then

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cd sgv ; bash install
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This will take you about **30 seconds** ...

- Study README do get the first test job done (another **30 seconds**)
- Look README in the `samples` sub-directory, to enhance the capabilities, eg.:
  - Get STDHEP installed.
  - Get CERNLIB installed in native 64bit.
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# Calorimeter simulation

## The issues:

- Clearly: Random E, shower position, shower shape.
- But also association errors:
  - Clusters might merge.
  - Clusters might split.
  - Clusters might get wrongly associated to tracks.
- Consequences:
  - If a (part of) a neutral cluster associated to track → Energy is lost.
  - If a (part of) a charged cluster not associated to any track → Energy is double-counted.
  - Other errors (split neutral cluster, charged cluster associated with wrong track ....) are of less importance.

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# Calorimeter simulation: SGV strategy

- Concentrate on what really matters:
  - True charged particles **splitting off** (a part of) their shower: **double-counting**.
  - True neutral particles **merging** (a part of) their shower with charged particles: **energy loss**.
- Don't care about neutral-neutral or charged-charged merging.
- Nor about multiple splitting/merging.
- Then: identify the **most relevant variables** available in fast simulation:
  - Cluster energy.
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# Observations

- Identify and factorise:
  - ① Probability to split
  - ② If split, probability to split off/merge the entire cluster.
  - ③ If split, but not 100 %: Form of the p.d.f. of the fraction split off.
- Depends on:
  - ① Isolation - strongly for merging, slightly for splitting - but can be treated in two energy bins with no energy dependence in the bin. %5 over-all dependence on barrel/endcap.
  - ② Energy. Is small for splitting, important for merging at low E.
  - ③ Energy and isolation (very little for splitting) via the average.
- All cases (EM/had - split/merge - Barrel/endcap) can be described by the same functional shapes.
- Functions are combinations of exponentials and lines.
- 28 parameters  $\times$  4 cases (em/had  $\times$  double-counting/loss)
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# Parametrisation in SGV

Call **ZAUPFL** in ZAUSER (instead of ZAUSHO which is for course-grained calorimeters):

- On first call: Reads and **stores parameters**
- **ZAUMKC** makes clusters from simulated showers of each particle.
- **Loop** (once) over all particles.
  - Determine shortest distance to neighbour of right type and charge in the same calorimeter. Fortran95's whole-array capabilities very useful to produce efficient code !
  - Simulate:
    - **Split shower** - combine cluster.
- **Remove** and push out any empty showers or clusters. Adjust pointers.

**Analyse** as before: Charged particles from module ZATRS, neutrals from ZACLU.

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  - **Simulate:**
    - No split or **complete** split or **partial** split
    - If partial: **fraction** split off
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**Analyse** as before: Charged particles from module ZATRS, neutrals from ZACLU.

# Parametrisation in SGV

Call **ZAUPFL** in ZAUSER (instead of ZAUSHO which is for course-grained calorimeters):

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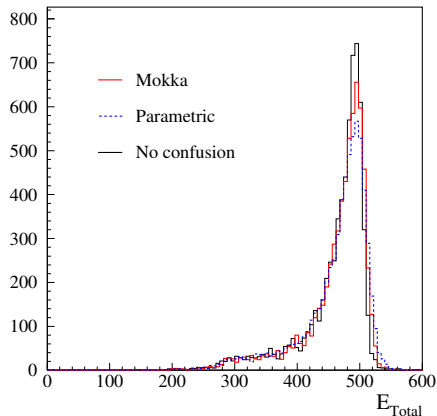
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# Checking the parametrisation

- Twiddle knobs: E vs p, overall split - probability.
- Total seen energy
- Total neutral energy
- Lost and double counted energy.

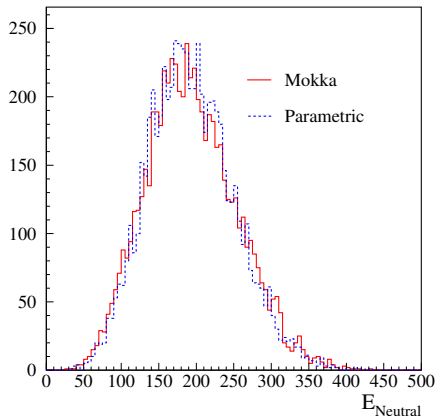
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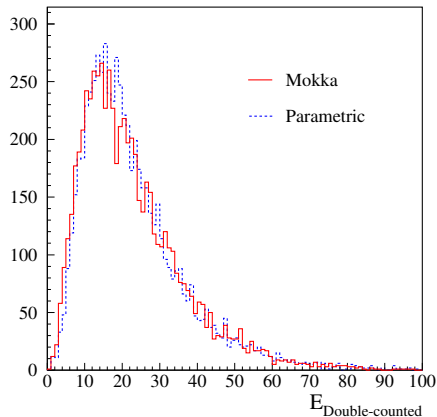
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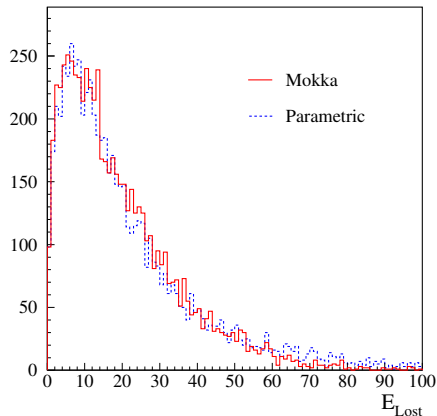
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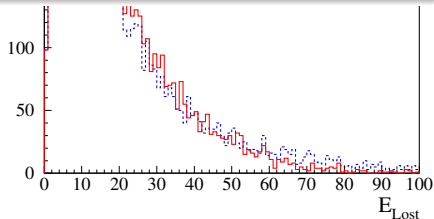
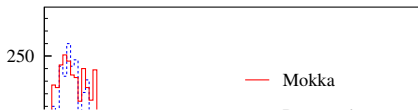


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Looks good ! Some further tests, then commit to SVN - this or next week.

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# Summary

- The status of the **SGV** program was presented.
- The particle-flow parametrisation, presented in earlier meetings, has been integrated into SGV.
- First comparisons to Mokka/Marlin with a first tentative tuning was shown.

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(On time-scale days to weeks)

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  - In the last case: output `STDHEP` of event
- Finish up `particle flow` parametrisation.
- Fix a few identified issues, then Release `SGV3.0` (no rc1).
- Add secondary vertexes to LCIO output.
- Produce **LCIO DST:s** for the DBD bench-marks: DBD analyses can start  $\approx$  now, while waiting for full-sim.

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# Thank You !

# Backup

# BACKUP SLIDES



## Use-cases at the ILC

- Used for **fastsim physics studies**, eg. arXiv:hep-ph/0510088, arXiv:hep-ph/0508247, arXiv:hep-ph/0406010, arXiv:hep-ph/9911345 and arXiv:hep-ph/9911344.
- Used for **flavour-tagging training**.
- Used for overall **detector optimisation**, see Eg. Vienna ECFA WS (2007), See Ilcagenda > Conference and Workshops > 2005 > ECFA Vienna Tracking
- **GLD/LDC merging and LOI**, see eg. Ilcagenda > Detector Design & Physics Studies > Detector Design Concepts > ILD > ILD Workshop > ILD Meeting, Cambridge > Agenda > Sub-detector Optimisation I

The latter two: Use the Covariance machine to get **analytical expressions** for performance (ie. *not* simulation)

# White paper

- Written in Fortran 95.
- CERNLIB dependence. Much reduced wrt. old F77 version, mostly by using Fortran 95's built-in matrix algebra.
- Managed in SVN. Install script included.
- Features:
  - Callable PYTHIA, Whizard.
  - Input from PYJETS or stdhep.
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  - samples subdirectory with steering and code for eg. scan single particles, create hbook ntuple with "all" information (can be converted to ROOT w/ h2root). And: output LCIO DST.
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Study README, and README in the [samples](#) sub-directory, to eg.:

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- Update **documentation** and in-line comments, to reflect new structure.
- Consolidate use of **Fortran 95/203/2008** features. Possibly - when gcc/gfortran 4.4 (ie. Fortran 2003) is common-place - **Object Orientation**, **if there is no performance penalty**.
  - Use of user-defined types.
  - Use of PURE and ELEMENTAL routines,
  - Optimal choice between pointer, allocatable and automatic and/or assumed-size, assumed-shape, and explicit arrays.
- I/O over **FIFO**:s to avoid storage and I/O rate limitations.
- The **Grid**.
- Investigate running on **GPU**:s.

- Further reduce CERNLIB dependence - at a the cost of backward compatibility on steering files ? HBOOK dependence will remain in the foreseeable future - but only for user convenience : SGV itself doesn't need it.