

ACTS for clusterization

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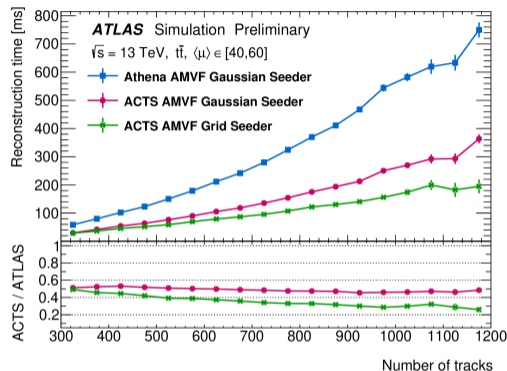
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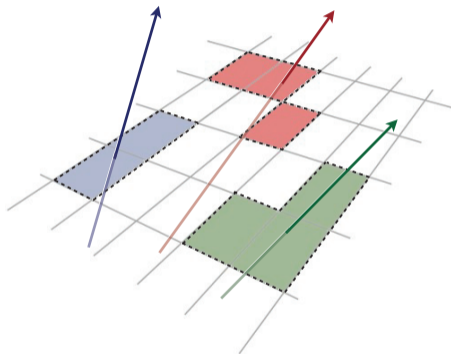
ACTS: A Common Tracking Software project

- ▶ Experiment-independent toolkit for tracking
- ▶ Free software (Mozilla Public License v2.0)
- ▶ Considered for use by Belle II, CEPC, sPHENIX, PANDA, FASER, **ATLAS**, EIC, ...
- ▶ Three overarching goals:
 1. Preserve current tracking approaches while enabling development for HL-LHC
 2. Serve as an algorithmic test bed including ML ([ONNX](#)) and accelerators ([traccc](#))
 3. Enable rapid and realistic development of new tracking detectors
- ▶ Overview paper: [\[2106.13593\]](#)
- ▶ Project webpage: [acts.readthedocs.io](#)
- ▶ Code repository: [github.com/acts-project/acts](#)



- ▶ ACTS already used in ATLAS Run 3 vertexing
- ▶ Gains in reconstruction speed when compared to Run 2 algorithm!

- ▶ Clusterization \equiv forming composite objects from discrete readout elements
- ▶ “Readout elements”: e.g. Silicon pixel (2D) or strip (1D) sensors
- ▶ Based on geometric connectivity: elements laid out on N-dimensional grid
- ▶ Connection logic can be extended e.g. with timing information
- ▶ Depending on context, clusterization can also include task of estimating crossing position



- ▶ Fast clusterization recently implemented in ACTS: [Clusterization.hpp](#)
- ▶ Based on [Hoschen-Kopelman](#) connected component labeling algorithm
- ▶ 2-pass algorithm:
 1. Assign labels to each sensor element and record equivalences (\equiv connections)
 2. Re-assign a “canonical” label to each element using recorded equivalences
- ▶ Equivalences between labels efficiently tracked by [disjoint-set](#) datastructure
- ▶ Algorithm requires sensors laid-out on 1-D or 2-D grid
 - ▶ 3-D would be possible but not implemented
 - ▶ Non-geometrical dimensions supported through extensible `Connection C++` type
- ▶ Only cluster formation is implemented; position estimation is user's responsibility
 - ▶ In general, initial position estimation requires knowledge of detector-specific readout geometry
 - ▶ Not available in ACTS detector geometry implementation
 - ▶ Precise position calibration requires knowledge of trajectory: Interface being designed

The Hoschen-Kopelman algorithm

- ▶ Input: A one-dimensional collection of “cells” (readout elements)
 - ▶ Output: None, the algorithm simply assigns labels to the cells
-
1. Sort cells (Column-wise order used for 2-D cells)
 2. Loop over cells in forward direction. For each cell:
 - 2.1 Assign a new label to the cell
 - 2.2 Loop backwards starting at nearest cell. For each cell pair:
 - 2.2.1 If they are more than one column away: break out of inner loop
 - 2.2.2 Else, if they are more than one row away: continue inner loop
 - 2.2.3 Else, they are connected: Mark their labels as equivalent and continue
 3. Loop over cells in forward direction. For each cell:
 - 3.1 Replace its label by “canonical” label based on recorded equivalences
-
- ▶ After second pass, all cells in a cluster have the same label

Clusterization algorithm details

► First pass

```
// Sort cells by position to enable in-order scan
std::sort(cells.begin(), cells.end(), internal::Compare<Cell, GridDim>());

// First pass: Allocate labels and record equivalences
for (auto it = cells.begin(); it != cells.end(); ++it) {
    const internal::Connections<GridDim> seen =
        internal::getConnections<Cell, Connect, GridDim>(it, cells, connect);
    if (seen.nconn == 0) {
        // Allocate new label
        getCellLabel(*it) = ds.makeSet();
    } else {
        // Sanity check: first element should always have
        // label if nconn > 0
        if (seen.buf[0] == NO_LABEL) {
            throw std::logic_error("seen.nconn > 0 but seen.buf[0] == NO_LABEL");
        }

        // Record equivalences
        for (size_t i = 1; i < seen.nconn; i++) {
            // Sanity check: since connection lookup is always backward
            // while iteration is forward, all connected cells found here
            // should have a label
            if (seen.buf[i] == NO_LABEL) {
                throw std::logic_error("i < seen.nconn but see.buf[i] == NO_LABEL");
            }

            // Only record equivalence if needed
            if (seen.buf[0] != seen.buf[i]) {
                ds.unionSet(seen.buf[0], seen.buf[i]);
            }
        }

        // Set label for current cell
        getCellLabel(*it) = seen.buf[0];
    }
}
```

► Default connection logic

```
// Cell collection logic
template <typename Cell, typename Connect, size_t GridDim>
Connections<GridDim> getConnections(typename std::vector<Cell>::iterator it,
                                     std::vector<Cell>& set, Connect connect) {

    Connections<GridDim> seen;
    typename std::vector<Cell>::iterator it_2(it);

    while (it_2 != set.begin()) {
        it_2 = std::prev(it_2);

        ConnectResult cr = connect(*it, *it_2);
        if (cr == eNoConnStop) {
            break;
        }
        if (cr == eNoConn) {
            continue;
        }
        if (cr == eConn) {
            seen.buf[seen.nconn++] = getCellLabel(*it_2);
            if (seen.nconn == seen.buf.size()) {
                break;
            }
        }
    }
    return seen;
}
```

► Second pass

```
// Second pass: Merge labels based on recorded equivalences
for (auto& cell : cells) {
    Label& lbl = getCellLabel(cell);
    lbl = ds.findSet(lbl);
}
```

- ▶ Algorithm does not require a specific “Cell” type
- ▶ Instead, algorithm is templated on cell type
- ▶ User needs to define three functions:
 - ▶ `int getColumn(const Cell&)`, → The primary sorting key
 - ▶ `int getRow(const Cell&)`, → The secondary sorting key (only if 2-D grid)
 - ▶ `int& getCellLabel(Cell&)` → Where the label gets stored
- ▶ This enables use of `labelClusters` function:

```
template <typename CellCollection, size_t GridDim = 2,  
         typename Connect =  
             DefaultConnect<typename CellCollection::value_type, GridDim>>  
void labelClusters(CellCollection& cells, Connect connect = Connect());
```

- ▶ For 2-D grid with default connection (geometric-only), use is as simple as:

```
std::vector<Cell> cells = ...;  
labelClusters(cells)
```

How to use? Cell connection

- ▶ Default connection is geometric only, based on adjacency
- ▶ Cell connection is extensible
- ▶ Example if want to use a pixel detector with timing information:

```
class MyCell;
struct MyConnect : Acts::Ccl::DefaultConnect<MyCell, 2> {
    ConnectResult operator()(const Cell& ref, const Cell& iter) {
        switch(Acts::Ccl::DefaultConnect<MyCell, 2>(ref, iter)) {
            case eNoConn:      return eNoConn;
            case eNoConnStop: return eNoConnStop;
            case eConn:        return timing_match(ref, iter)? eConn : eNoConn;
        }
    }
}
```

- ▶ Then:

```
std::vector<MyCell> cells = ...;
labelClusters<std::vector<MyCell>, 2, MyConnect>(cells, MyConnect());
```

- ▶ After labelling, creating clusters based on labels is easy
- ▶ For convenience, ACTS provides `mergeClusters` function
- ▶ Again, not imposing a specific cluster or cell type
- ▶ User needs to provide one function:
 - ▶ `void clusterAddCell(Cluster&, const Cell&)`
 - ▶ the `Cluster` type needs a default constructor
- ▶ Usage is straightforward:

```
std::vector<MyCell> cells = ...;
labelClusters(cells);
std::vector<MyCluster> clusters
    = mergeClusters<std::vector<MyCell>, std::vector<MyCluster>>(cells)
```

- ▶ In most use-cases, always chaining call to `labelClusters` and `mergeClusters`
- ▶ For convenience, ACTS provide a wrapper that does exactly this: `createClusters`
- ▶ Usage is straightforward. E.g. In default scenario (2-D grid, default connectivity)

```
std::vector<MyCell> cells = ...;  
std::vector<MyCluster> clusters= createClusters(cells)
```

- ▶ I gave some simple examples in previous slides
- ▶ Call signatures can get messier when deviate from default use case
- ▶ Can find real examples of the 1-D and 2-D cases in ATLAS Run 4 ACTS-based tracking:
 - ▶ 1-D: [ActsStripClusteringTool](#)
 - ▶ 2-D: [ActsPixelClusteringTool](#)
- ▶ Examples in ACTS itself:
 - ▶ In Digitization example: [ModuleClusters.cpp](#)
 - ▶ In test suite: [Clusterization unit tests](#)
- ▶ Otherwise, documentation is sorely lacking at this point
- ▶ Will be fixed Soon™
- ▶ I'm always happy to answer any questions

- ▶ Unfortunately no plots to show: my benchmarking uses ATLAS MC, so has to stay internal. . .
- ▶ Implementation performs best when:
 - ▶ Detector modules are \approx sparsely activated; or
 - ▶ if occupancy high, dominated by many small clusters
- ▶ In practice, these assumptions hold for overwhelming majority of ATLAS data
- ▶ Other experiments *may* have other requirements: let us know!
 - ▶ Optimal solution may be different in this case, and we want to support everyone

- ▶ ACTS has fast clusterization implementation based on Hoschen-Kopelman algorithm
- ▶ Only cluster formation – position estimation for now left to users
 - ▶ In future, will have support for at least position calibration
- ▶ Gave a few examples of usage and references to real-world code
- ▶ Implementation performs best for sparse module and/or small-cluster regime
 - ▶ Let us know if that's not your use case!
- ▶ More documentation will come at some point!