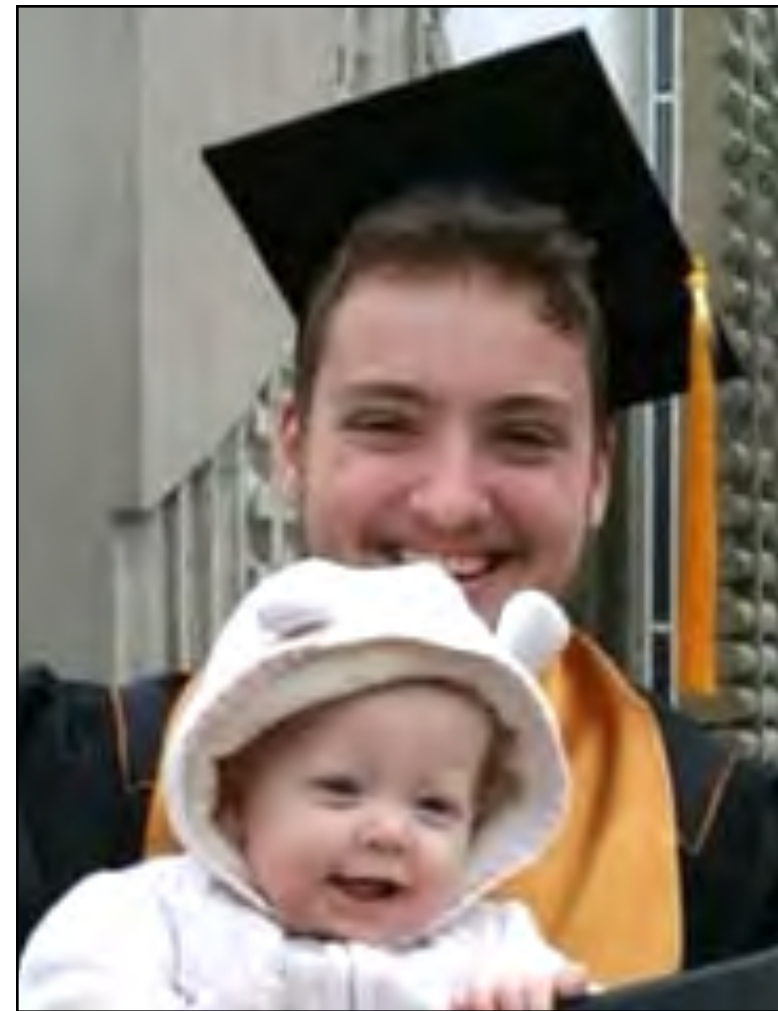


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# Reconstructing Proton-Proton Collision Positions at the Large Hadron Collider with D-Wave

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**Andrew J. Wildridge**



**Sachin B. Vaidya**



**Andreas Jung**



**Souvik Das**

- **Introduction to the Large Hadron Collider and the Compact Muon Solenoid detector**
- **QUBO formulation of the problem for D-Wave**
- **D-Wave performance, benchmarked against CPU**
- **Conclusions and outlook**

# The Large Hadron Collider



CMS Experiment

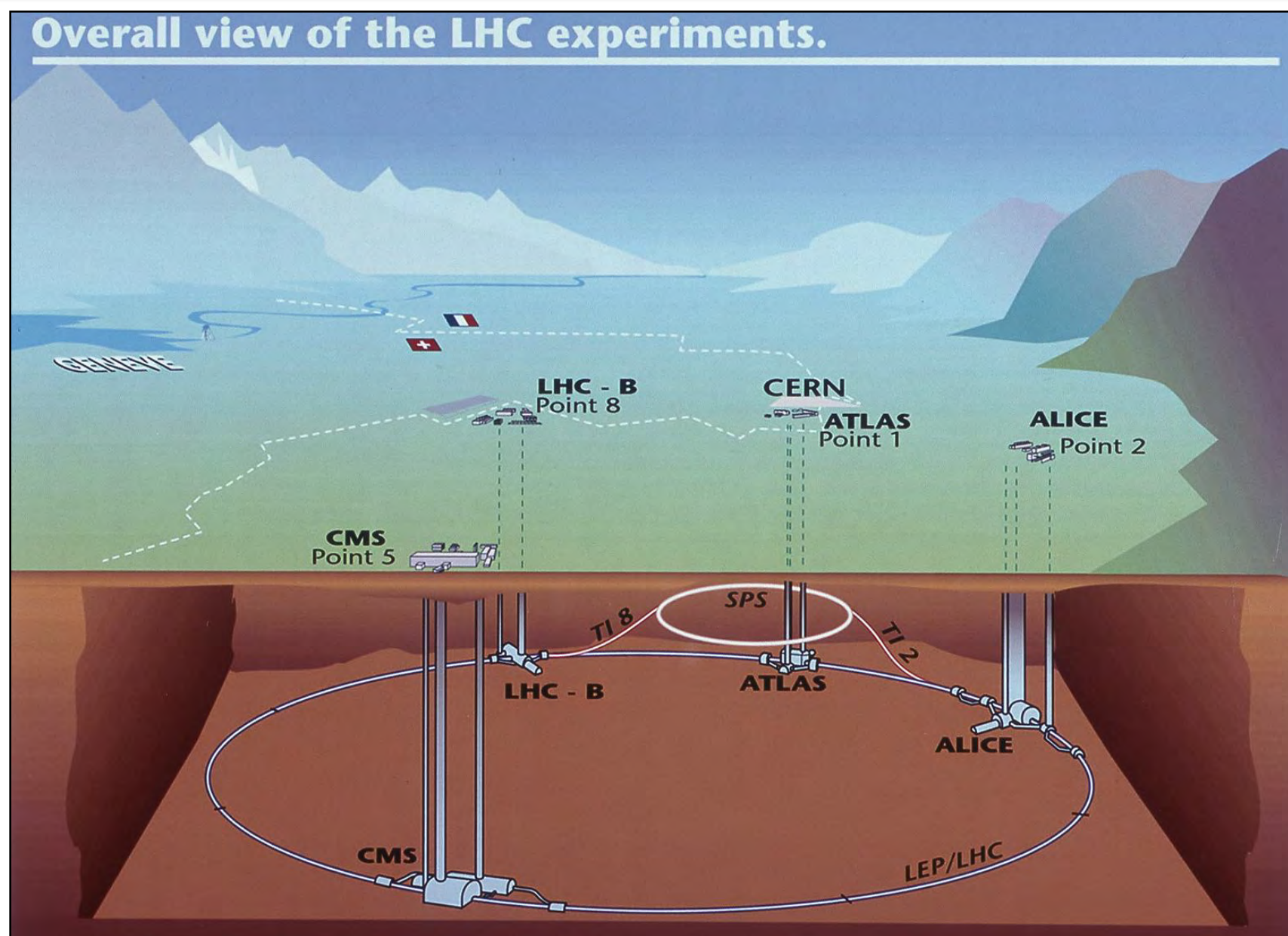
The Large Hadron Collider

# The Large Hadron Collider



- ◎ The Higgs boson
- ◎ Large extra-dimensions
- ◎ Supersymmetry
- ◎ Dark matter
- ◎ Baryogenesis

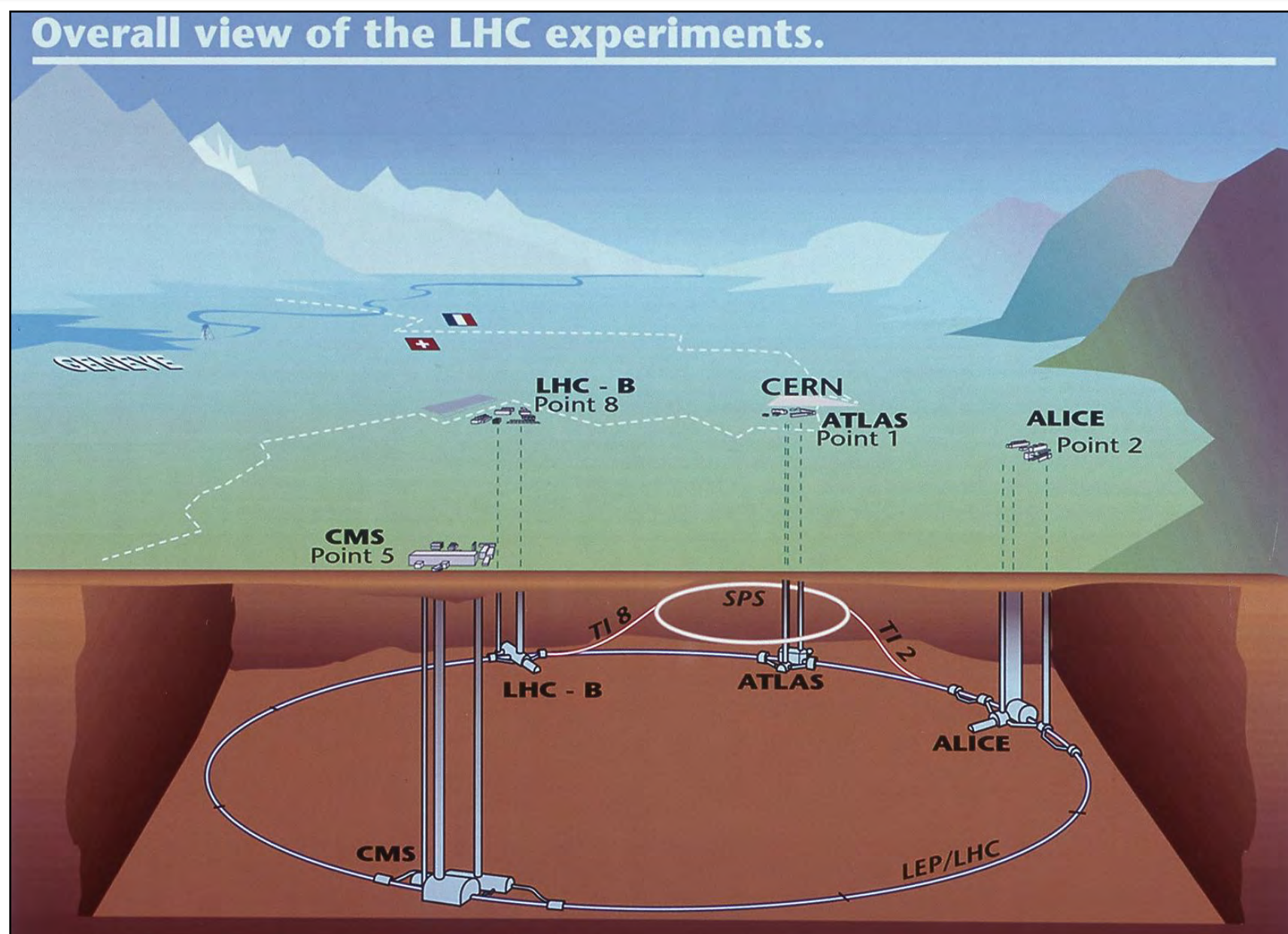
# Proton-proton collisions at the Large Hadron Collider



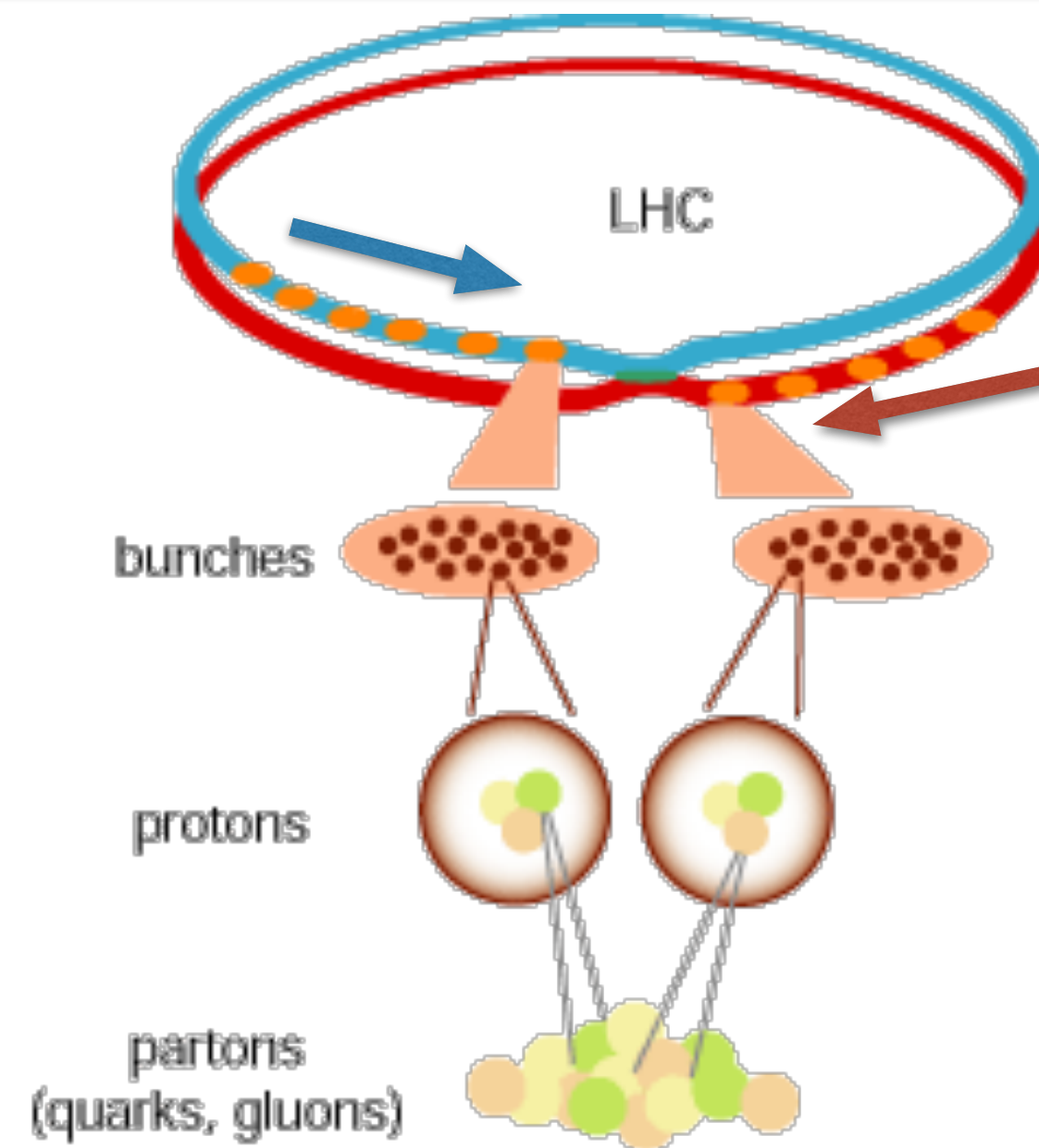
The LHC is 100 m below the surface, 27 km in circumference

**Looking for:** A high-energy physics problem that has a natural formulation for quantum annealing, and is simple

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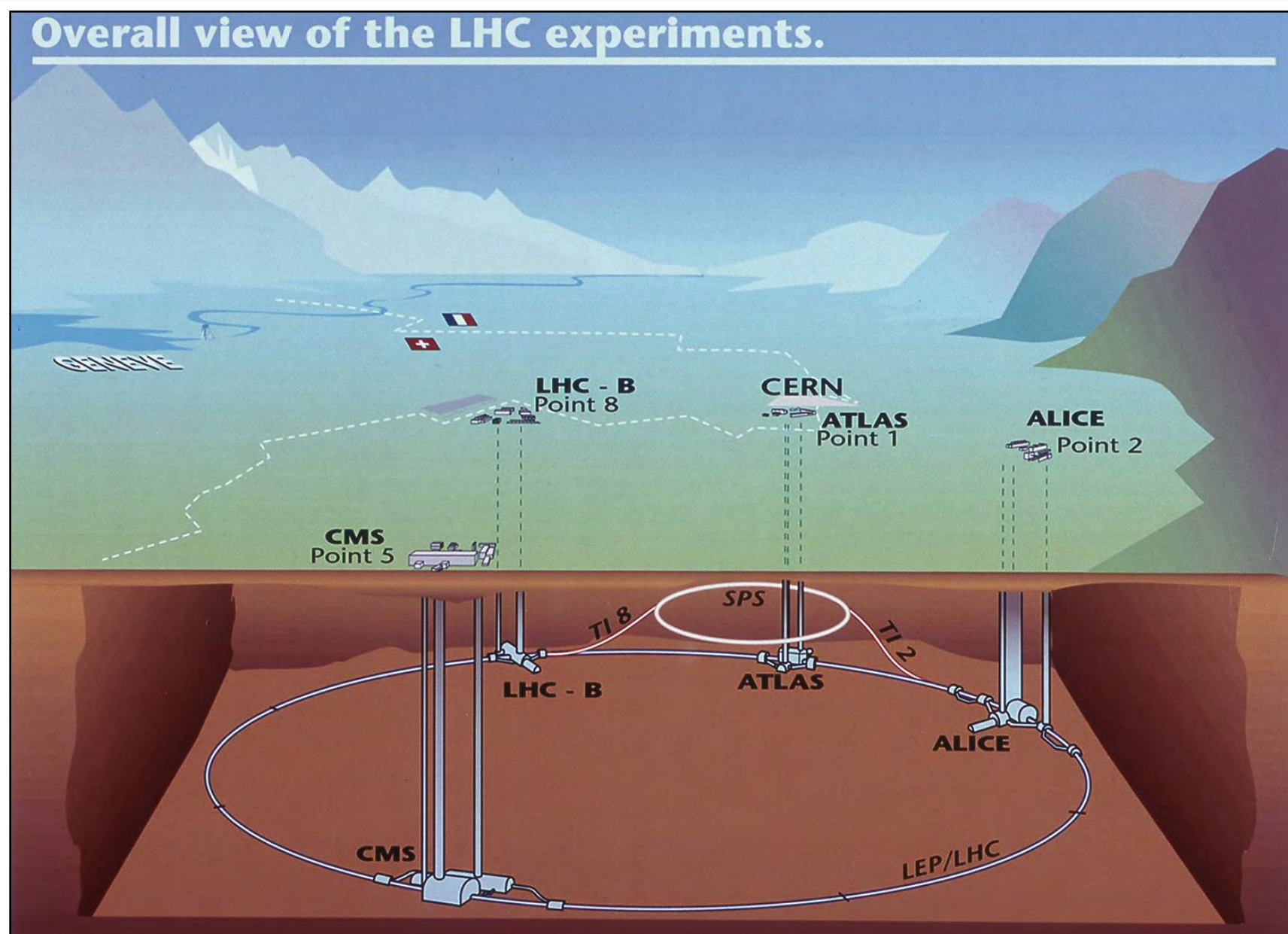
Structure of colliding proton bunches

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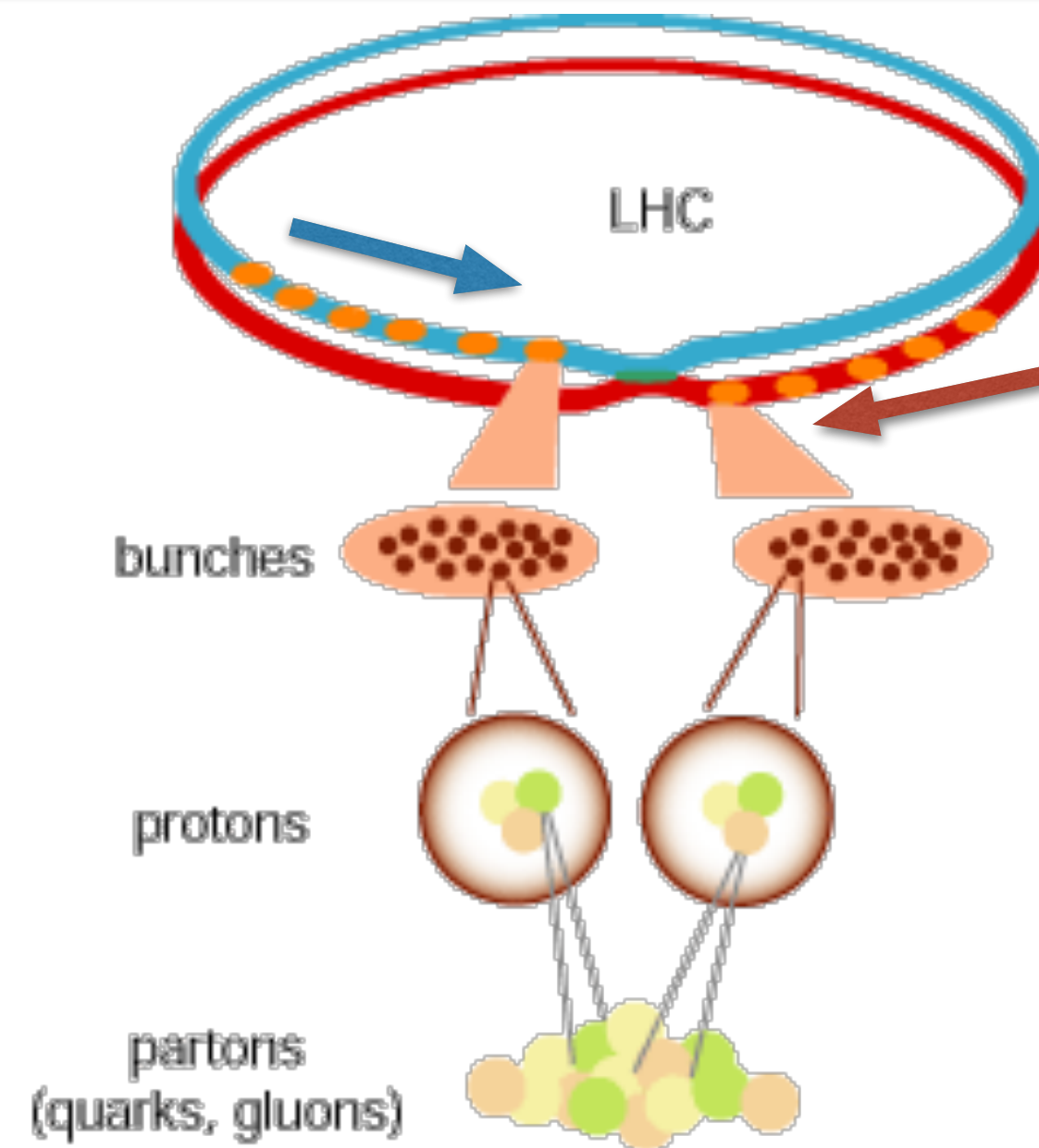
**Chosen problem: Reconstructing proton-proton collision positions at the Large Hadron Collider (LHC)**

- The LHC circulates protons inside its beam-pipes not in a continuous stream but in several closely packed **bunches**.
- Each bunch contains ~ 100 billion protons

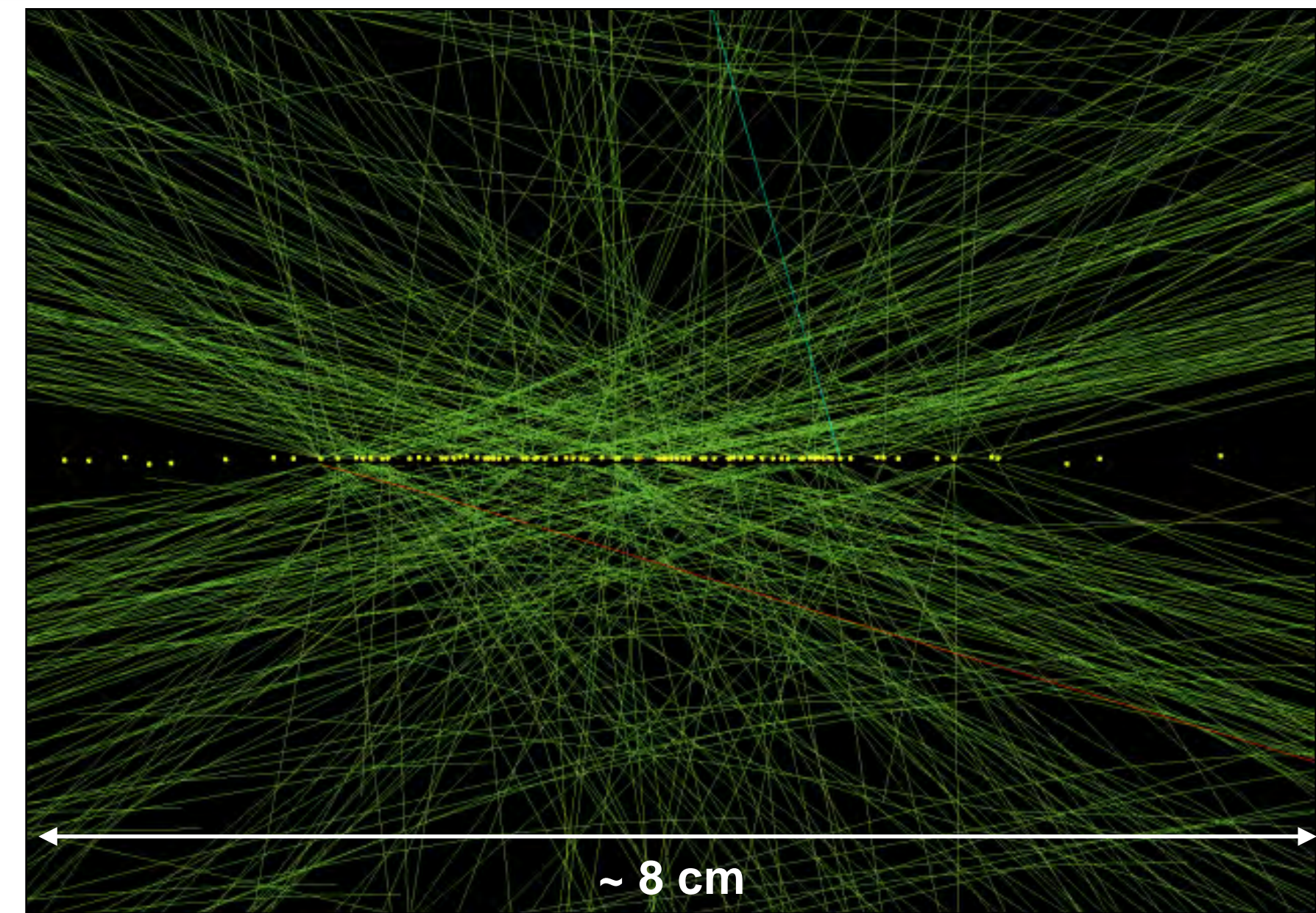
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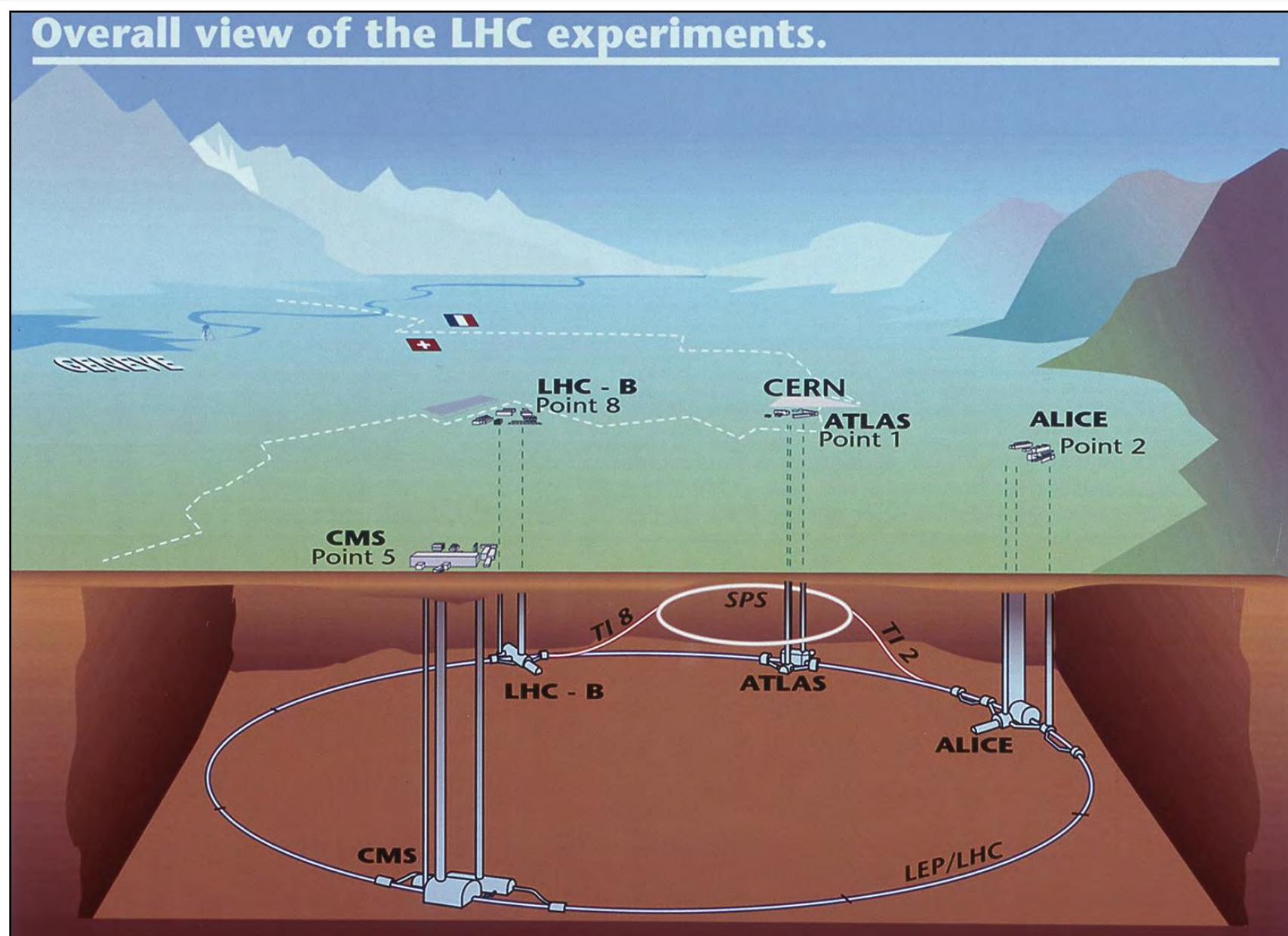
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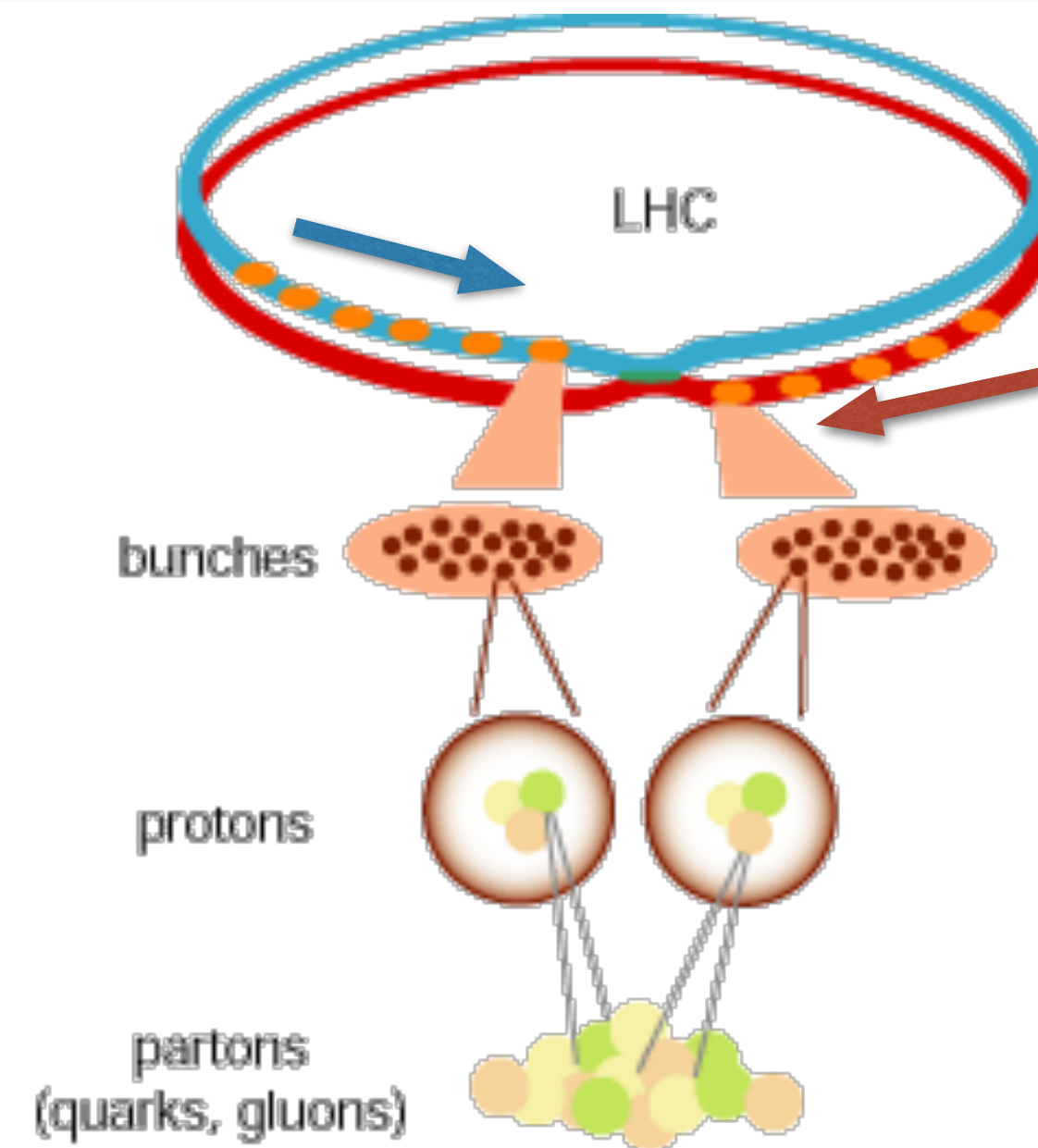
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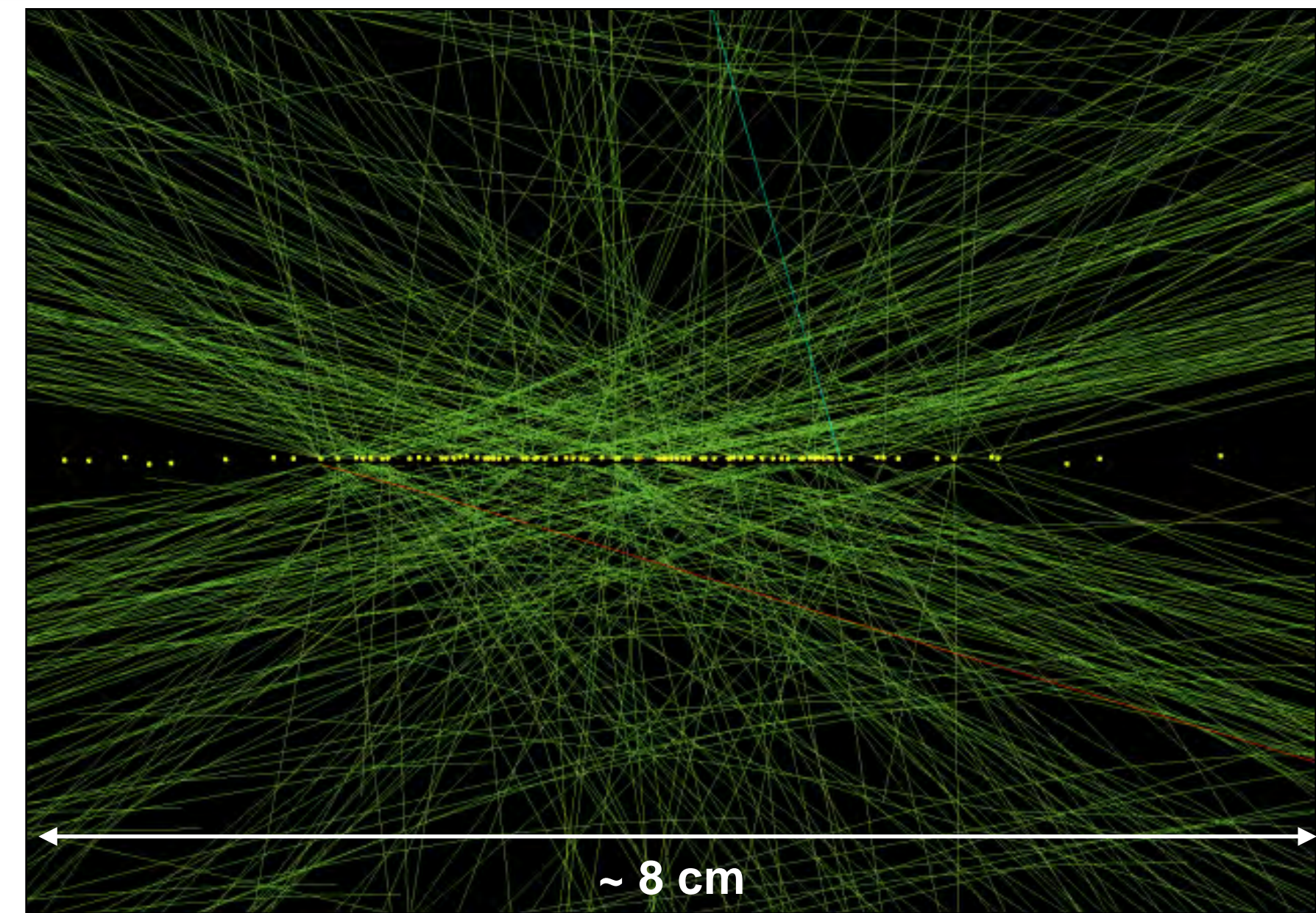
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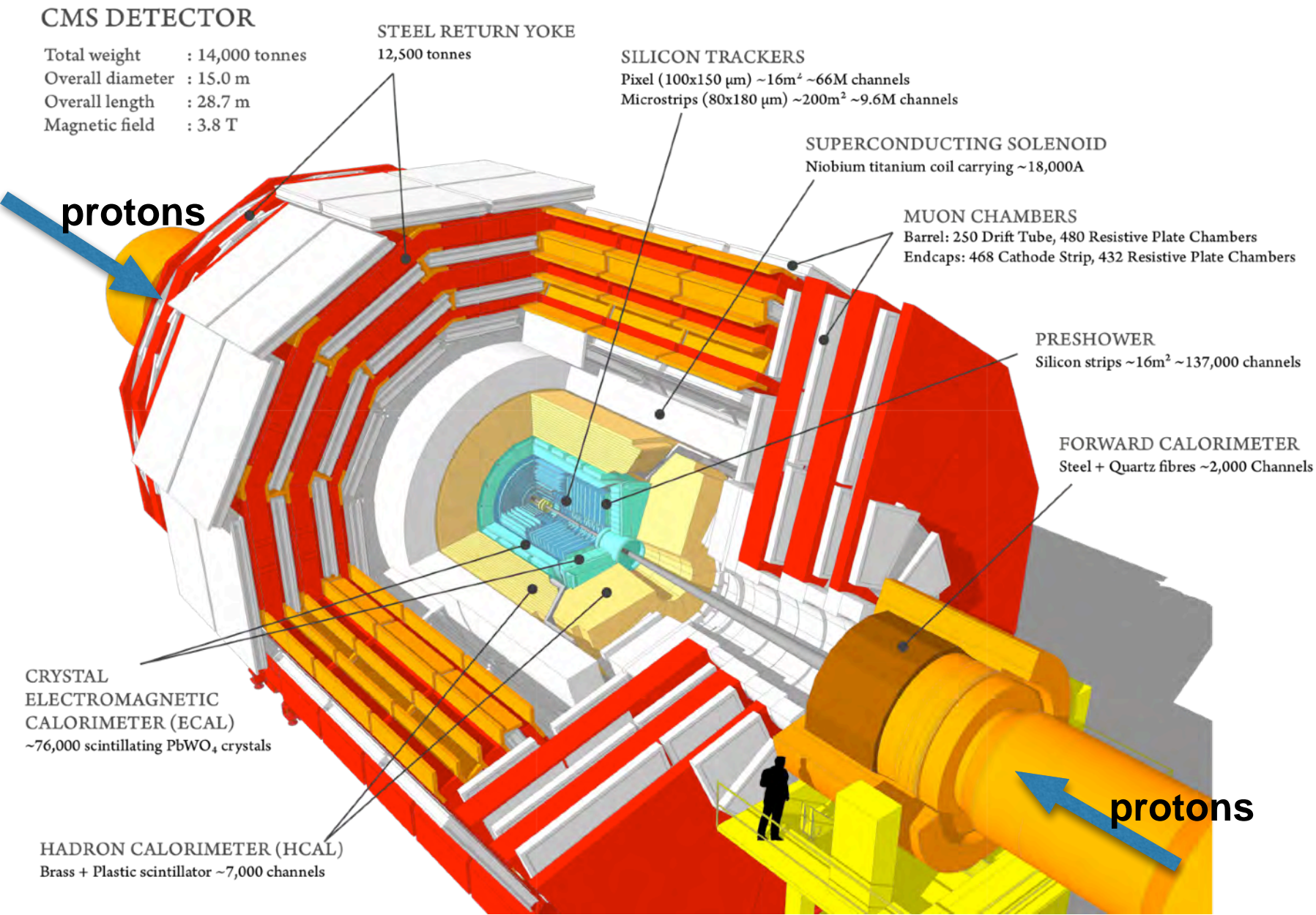
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*Which tracks come from which p-p collision?*

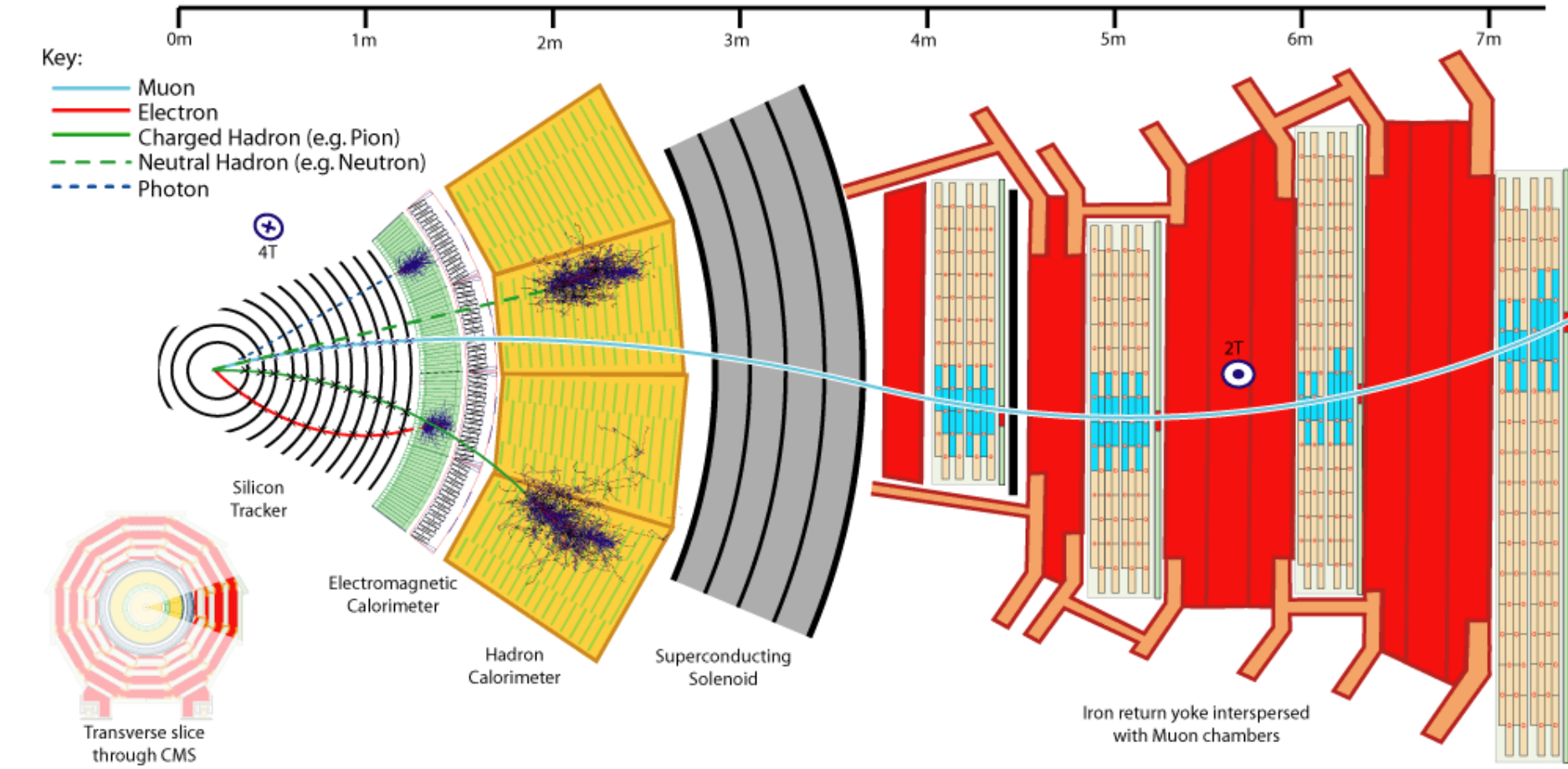
*Where are the p-p collision points in a bunch?*



# p-p collision position reconstruction at the Compact Muon Solenoid <sup>5</sup>



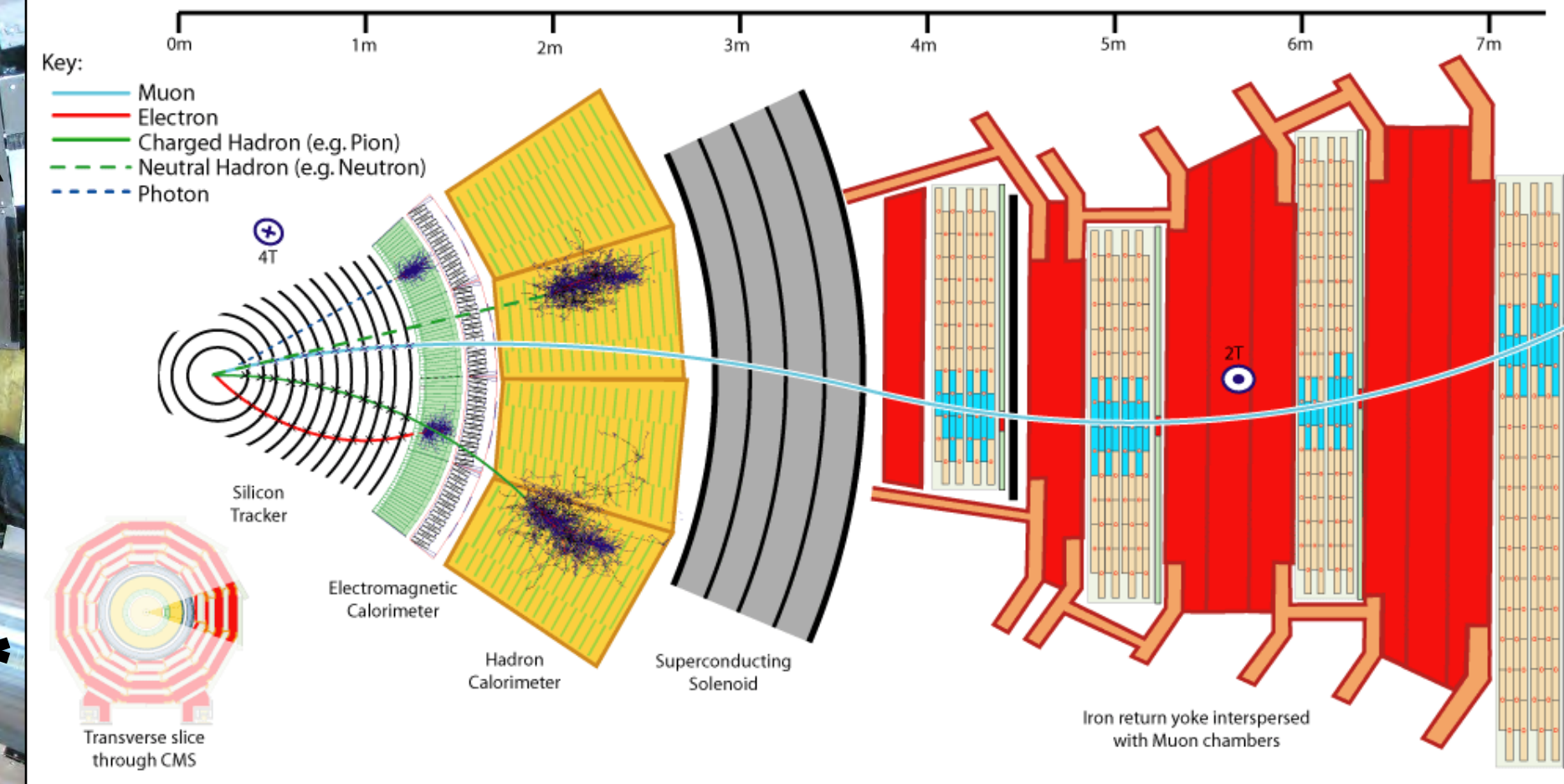
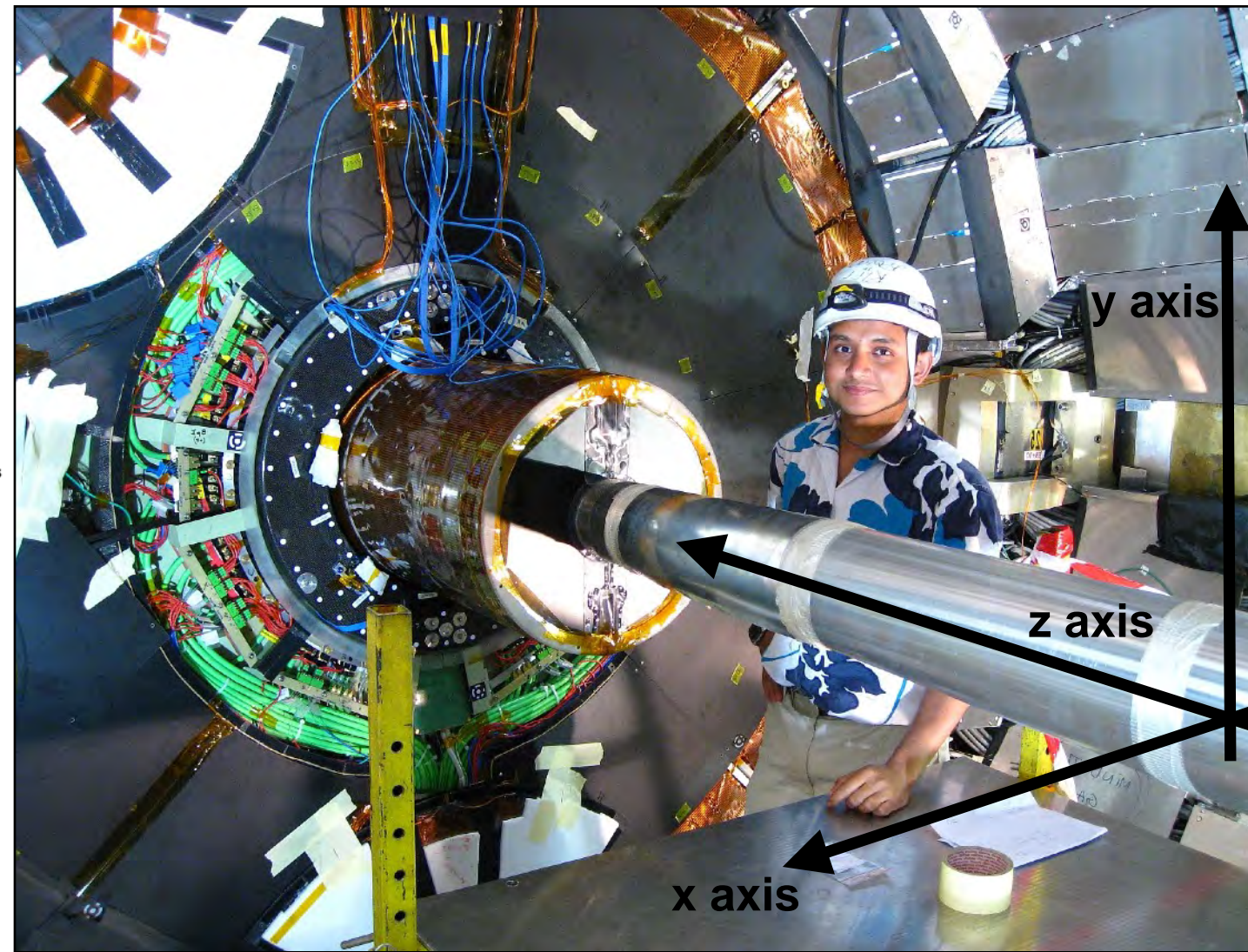
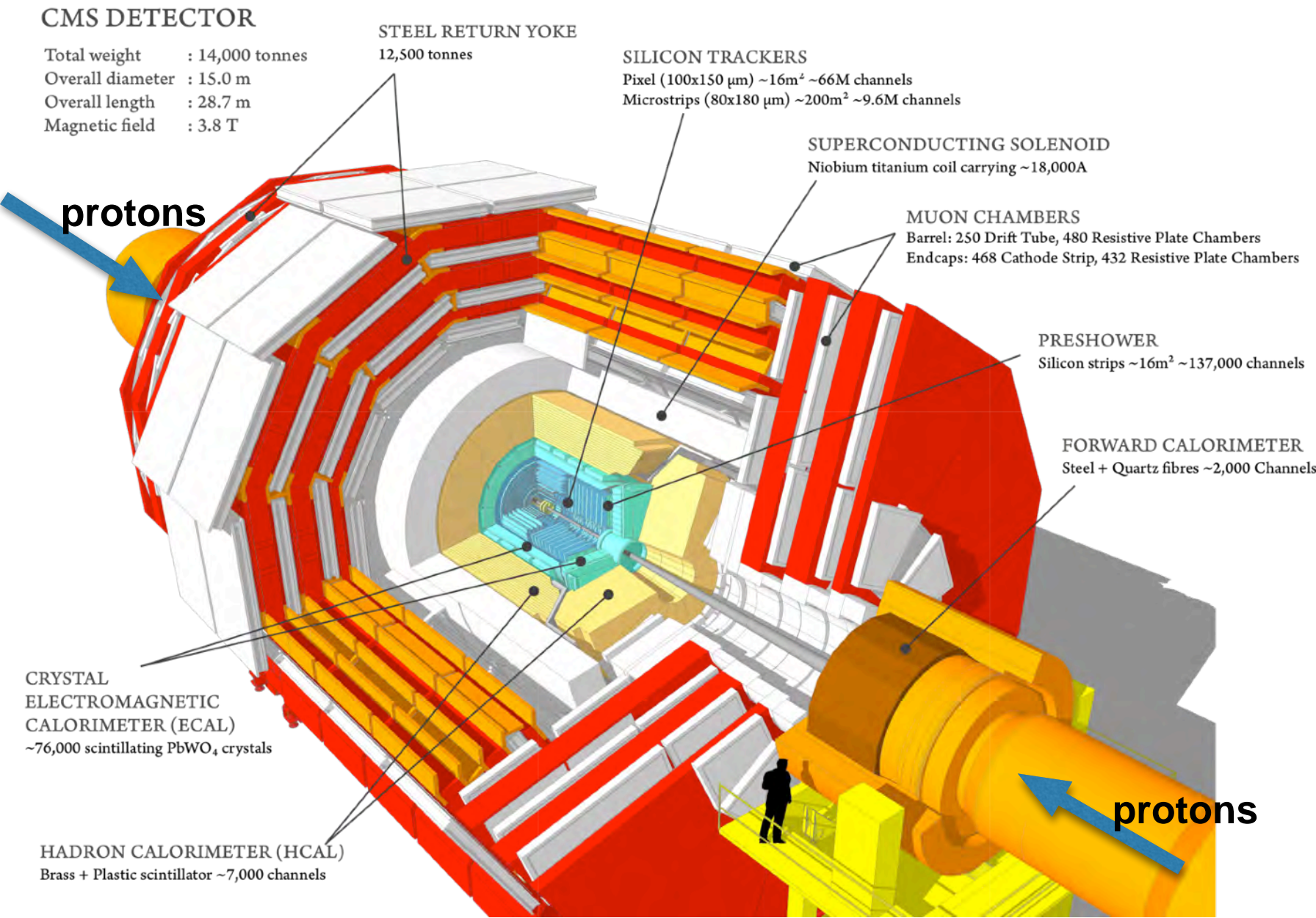
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In the x-y plane. Particle trajectories reconstructed as tracks. Reconstructions come with uncertainties

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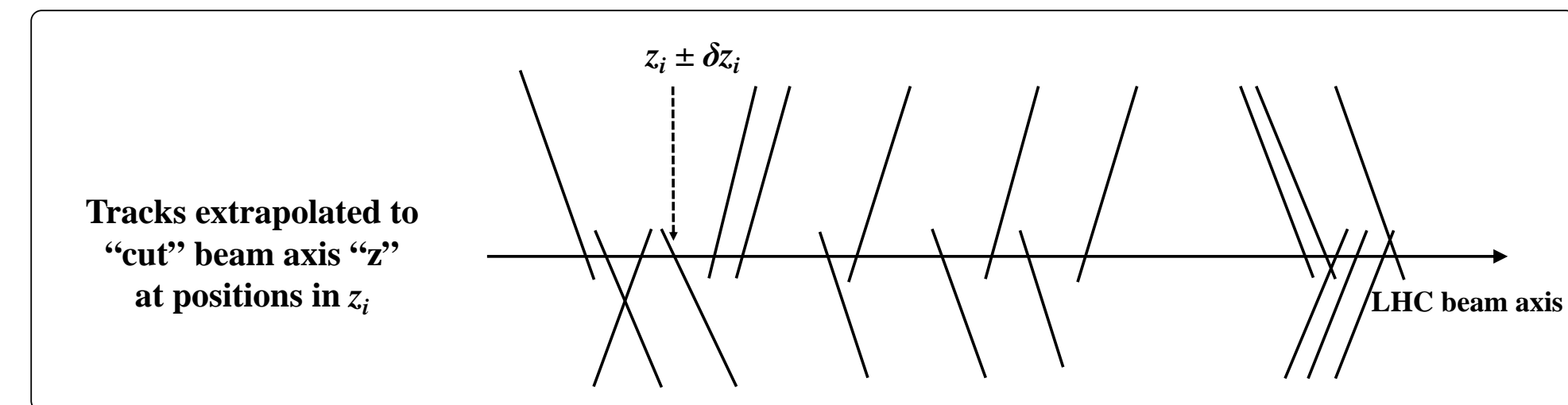
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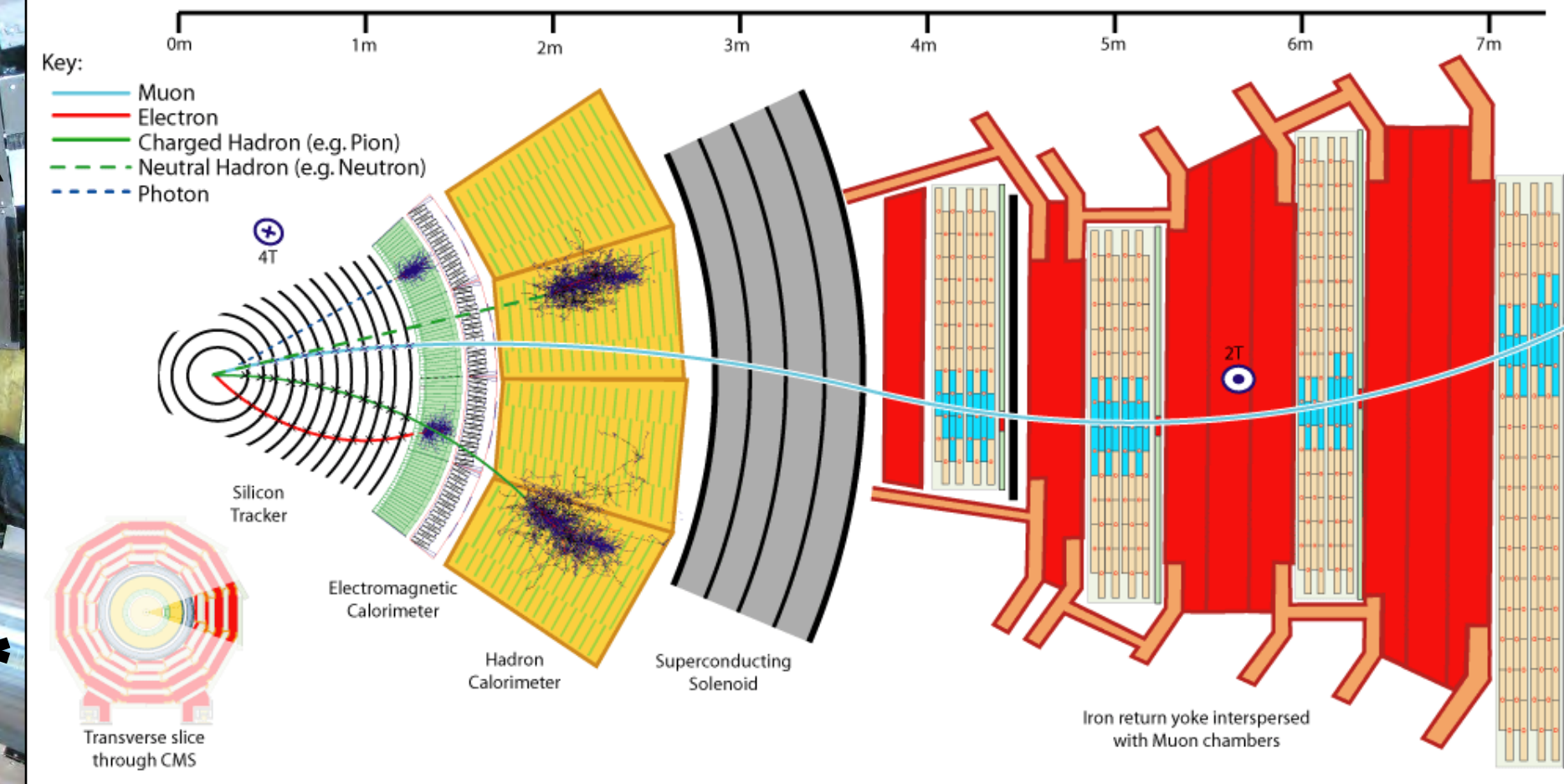
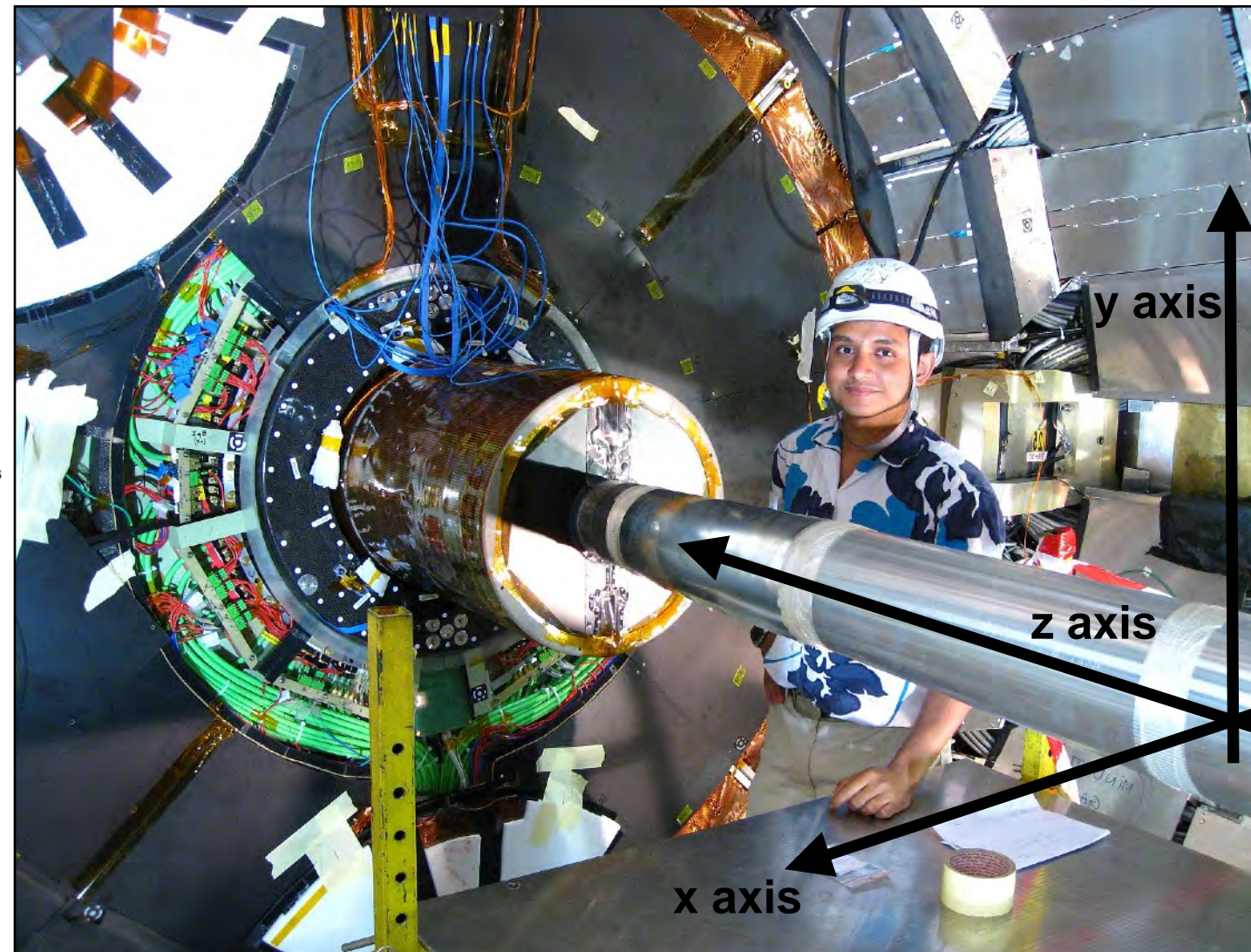
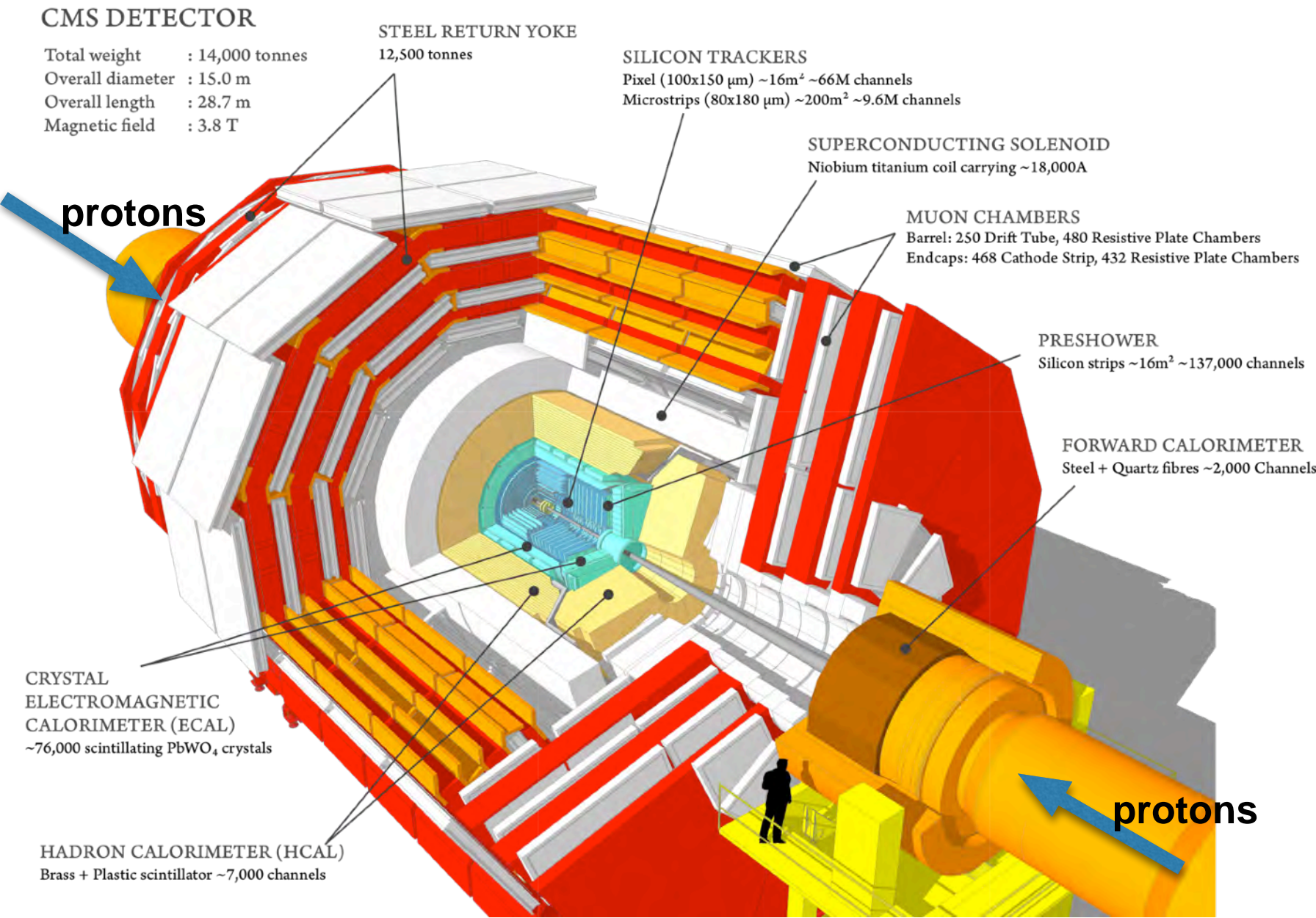
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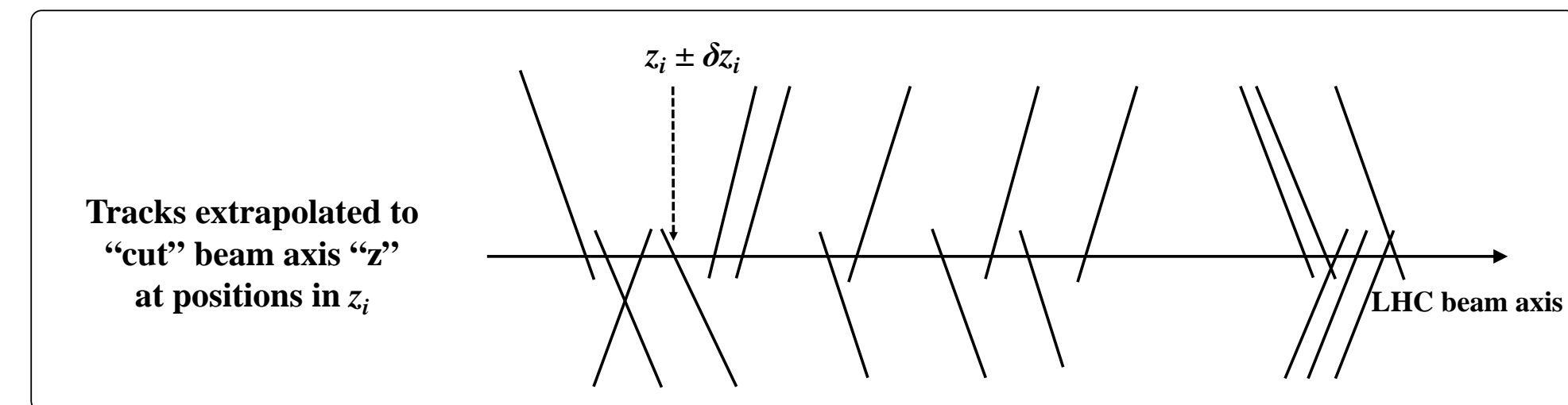
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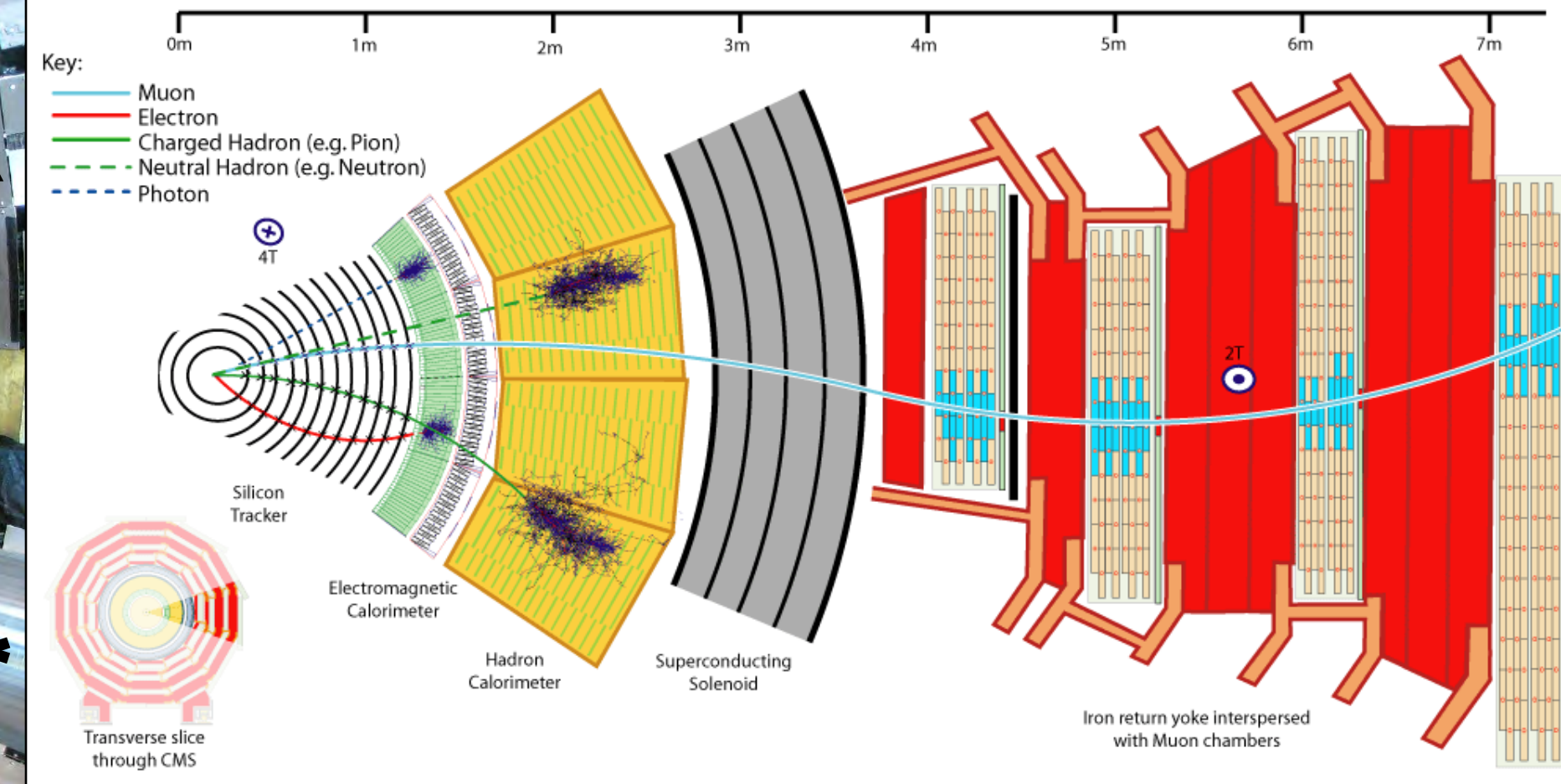
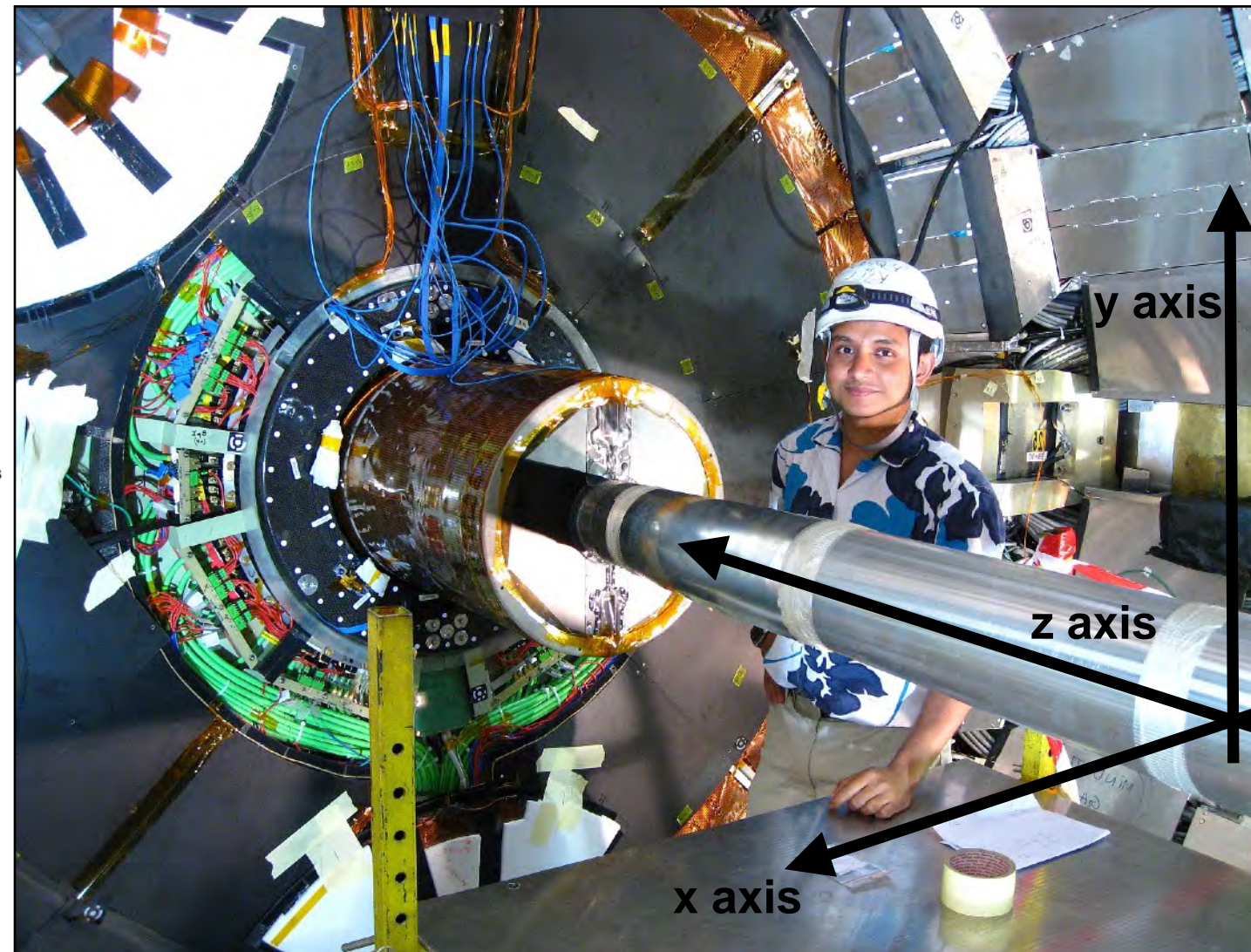
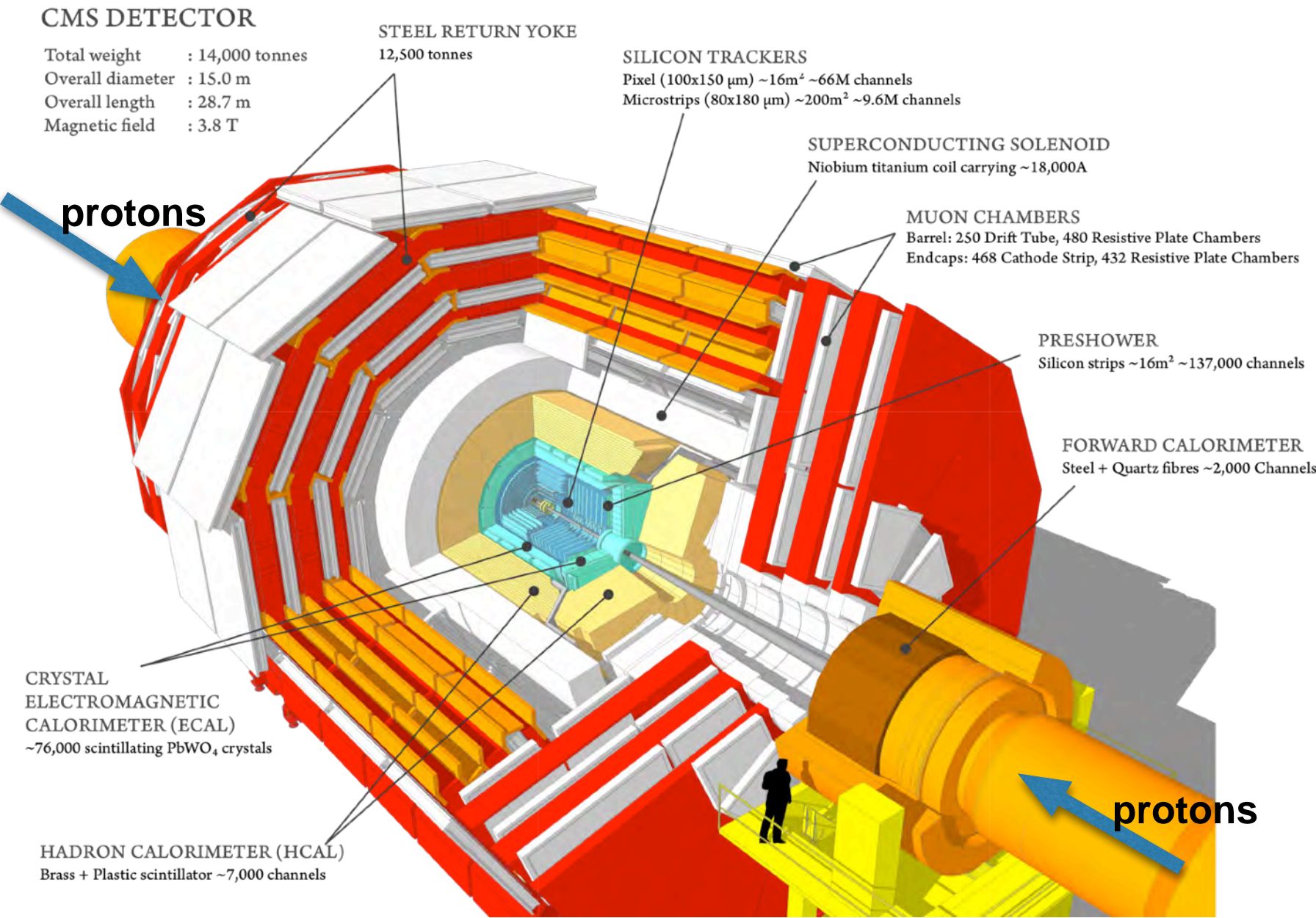
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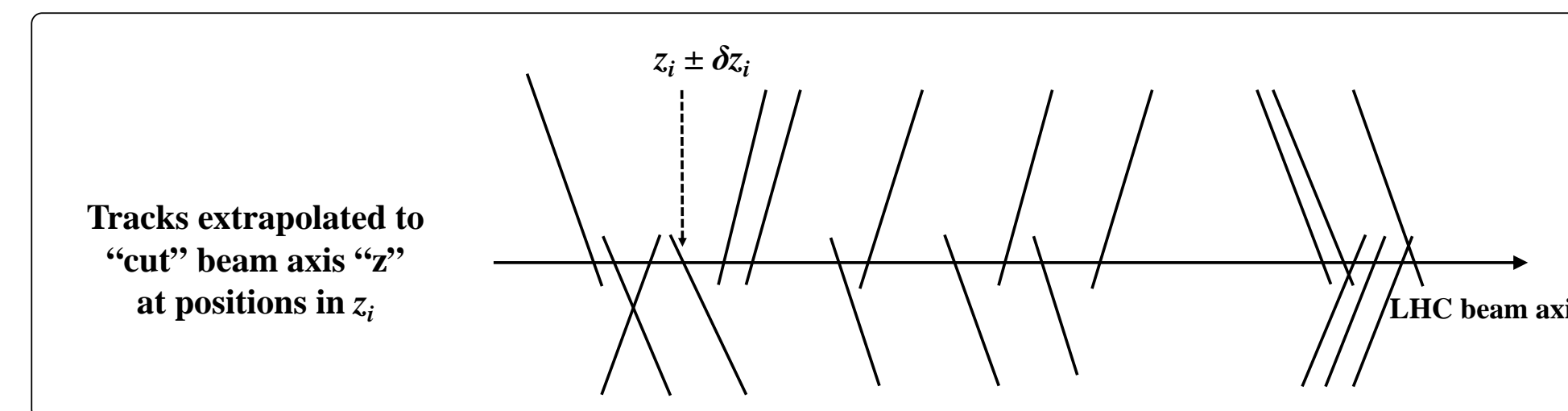
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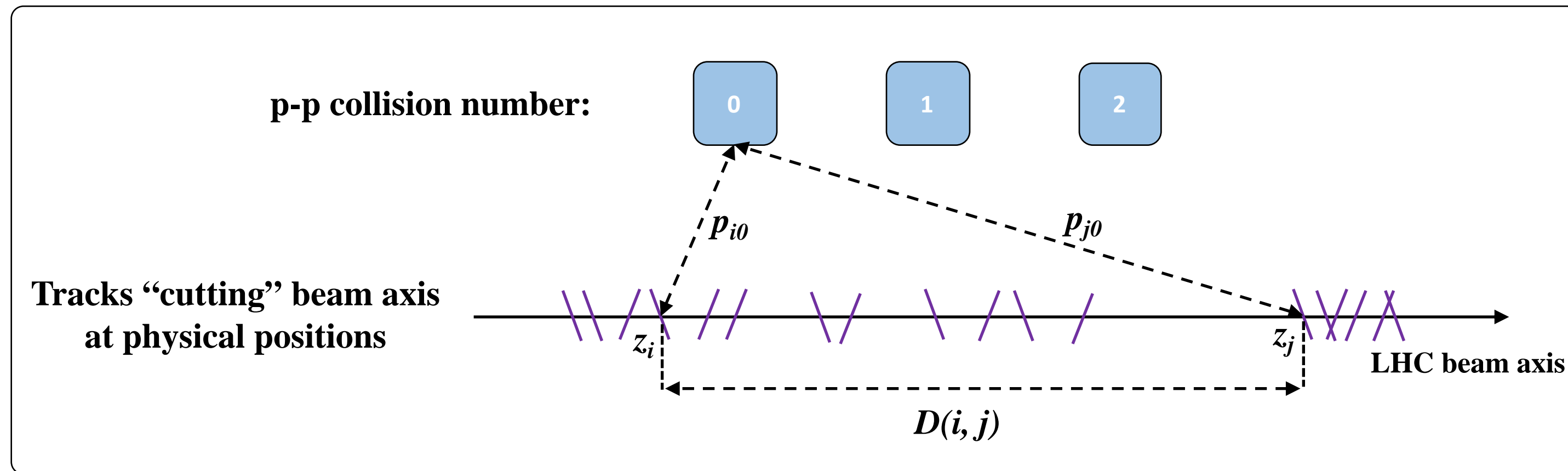
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Can D-Wave solve it using quantum annealing?

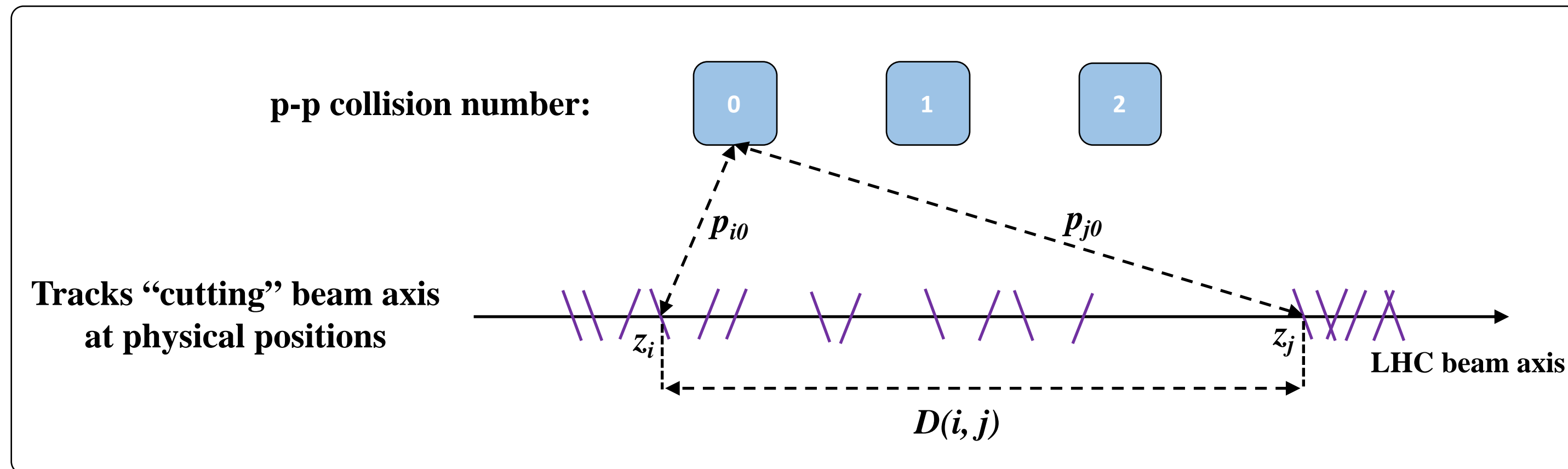
# Track clustering QUBO formulation for D-Wave



$$H_p = \sum_k^{n_V} \sum_i^{n_T} \sum_{j>i}^{n_T} p_{ik} p_{jk} g(D(i, j); m) + \lambda \sum_i^{n_T} \left( 1 - \sum_k^{n_V} p_{ik} \right)^2,$$

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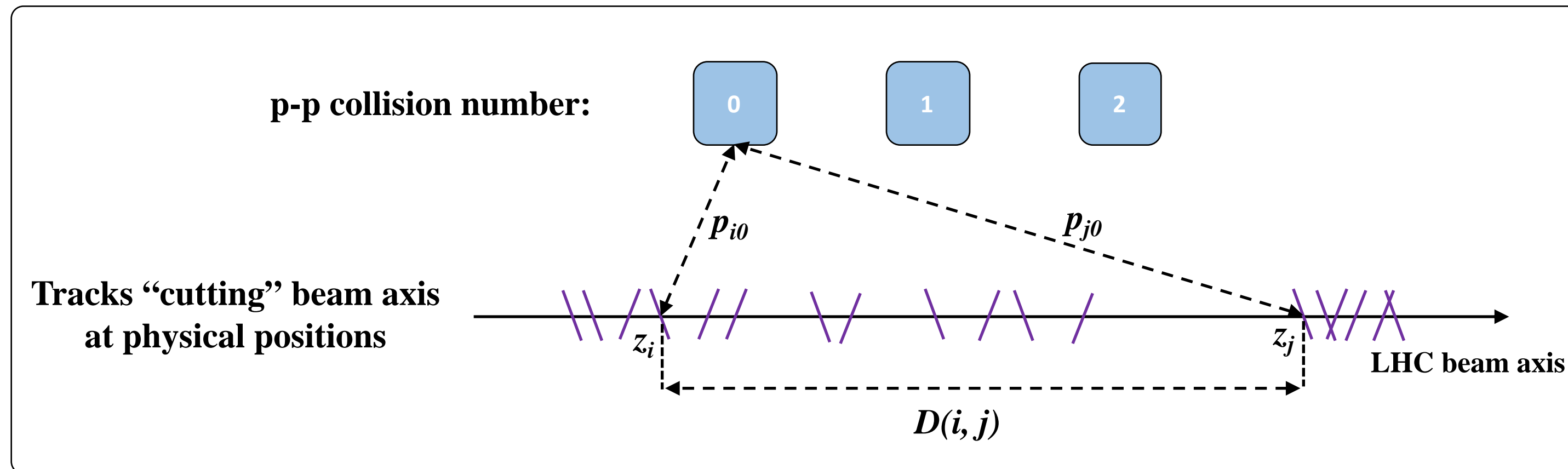
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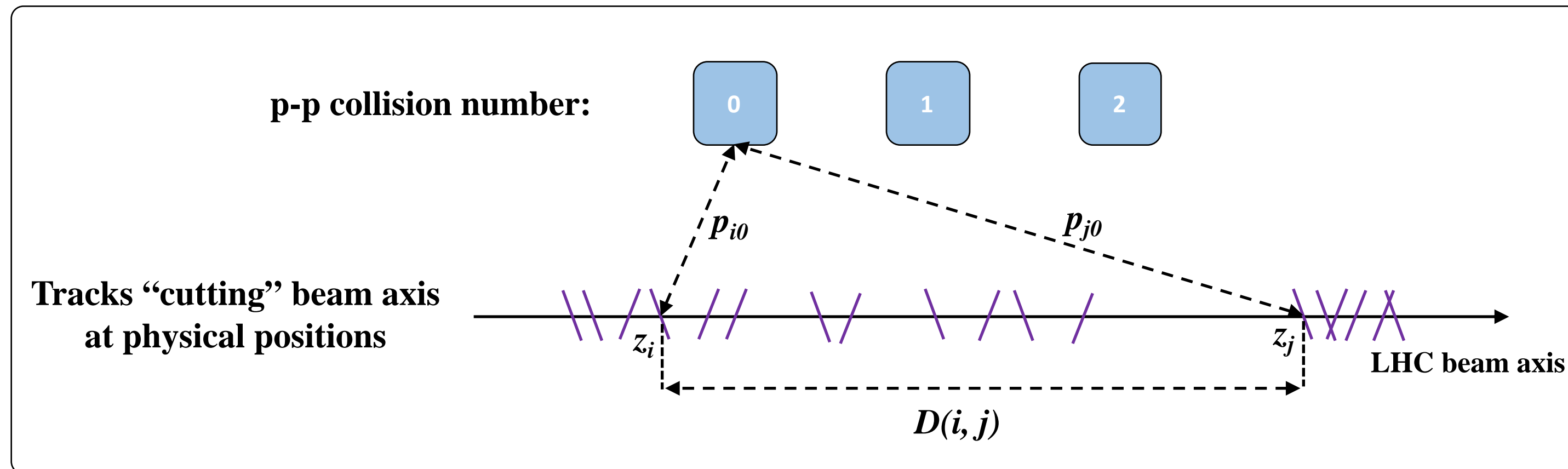
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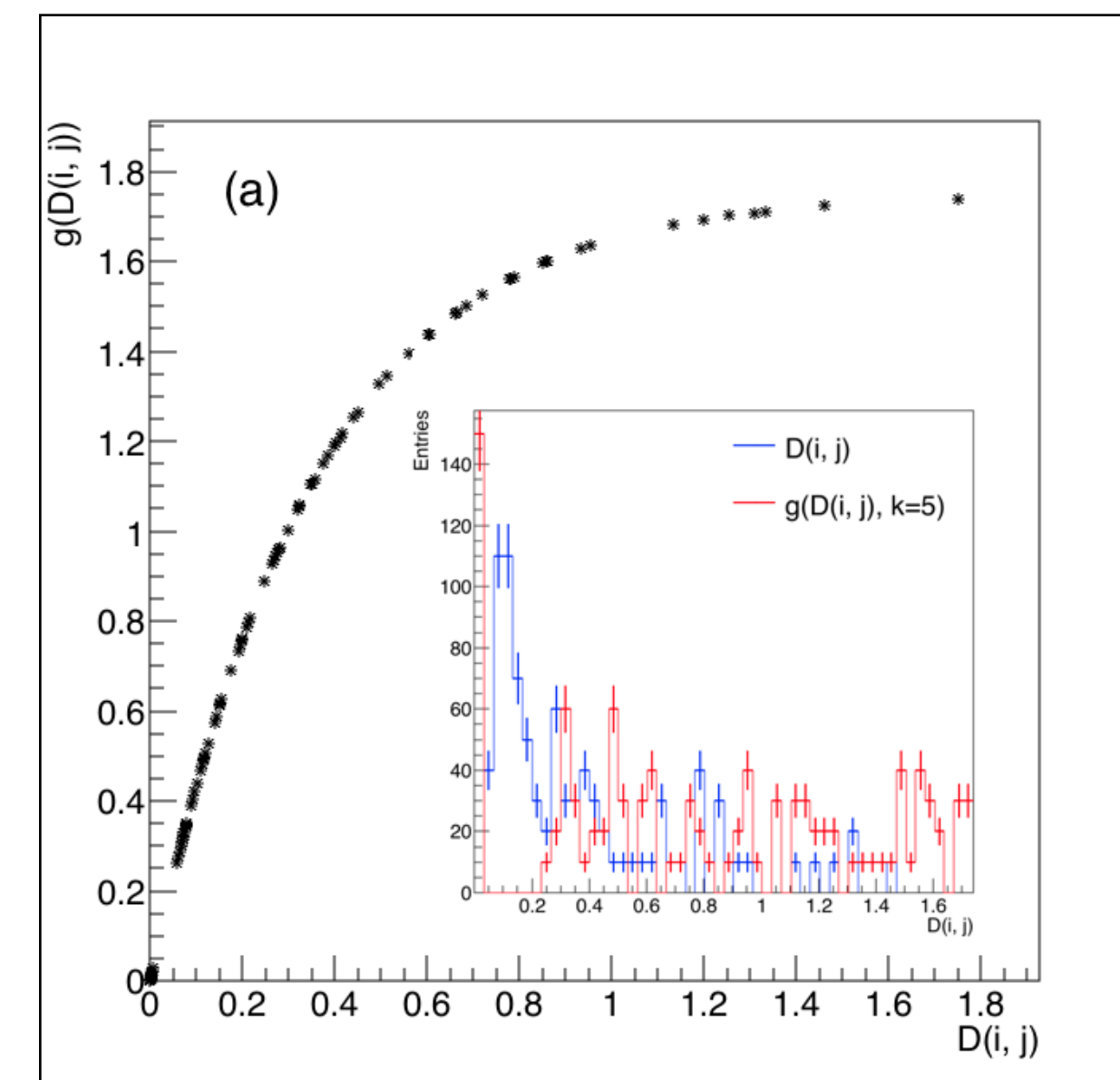
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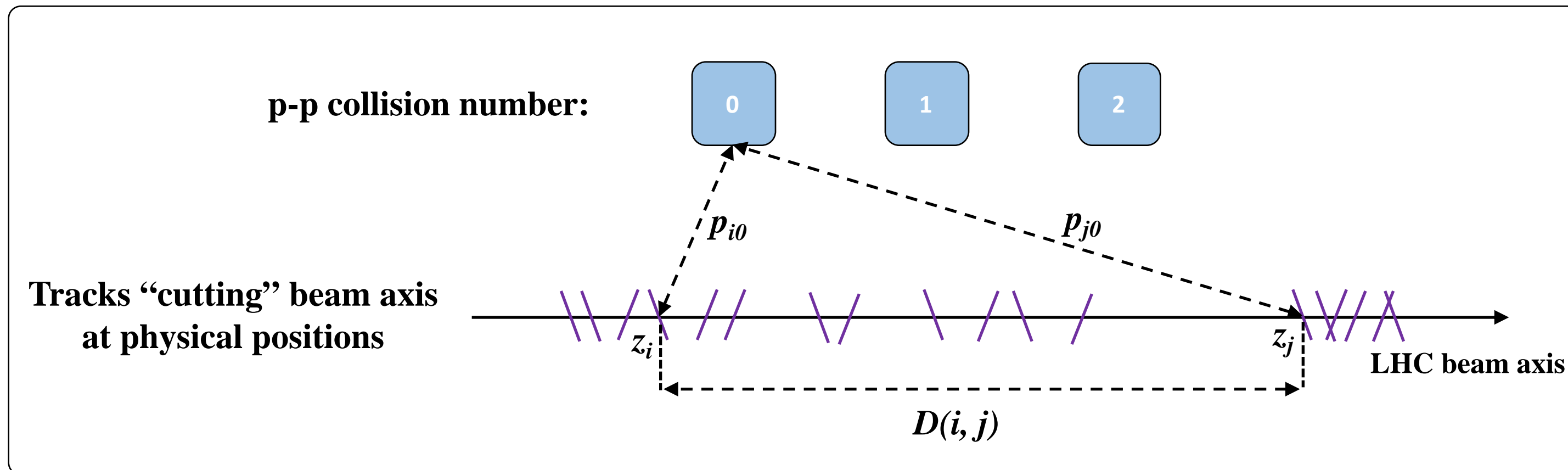
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$$g(x; m) = 1 - e^{-mx},$$





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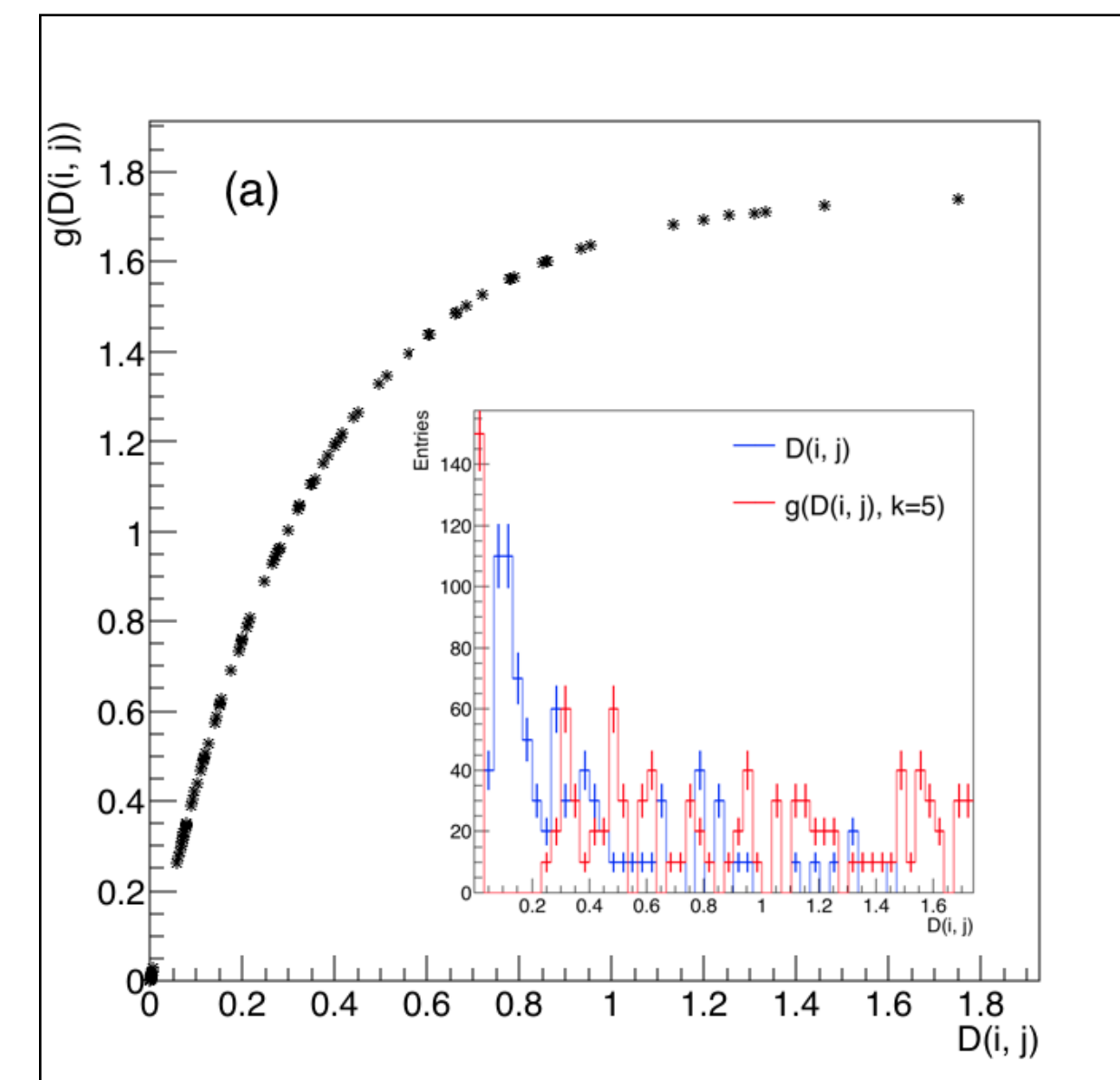
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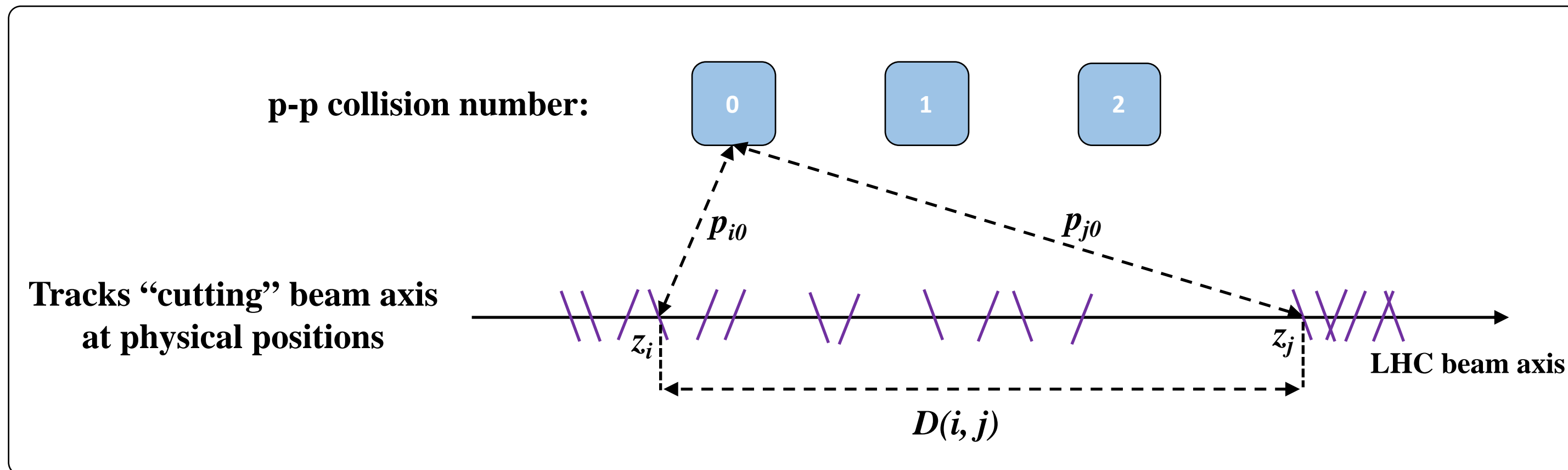
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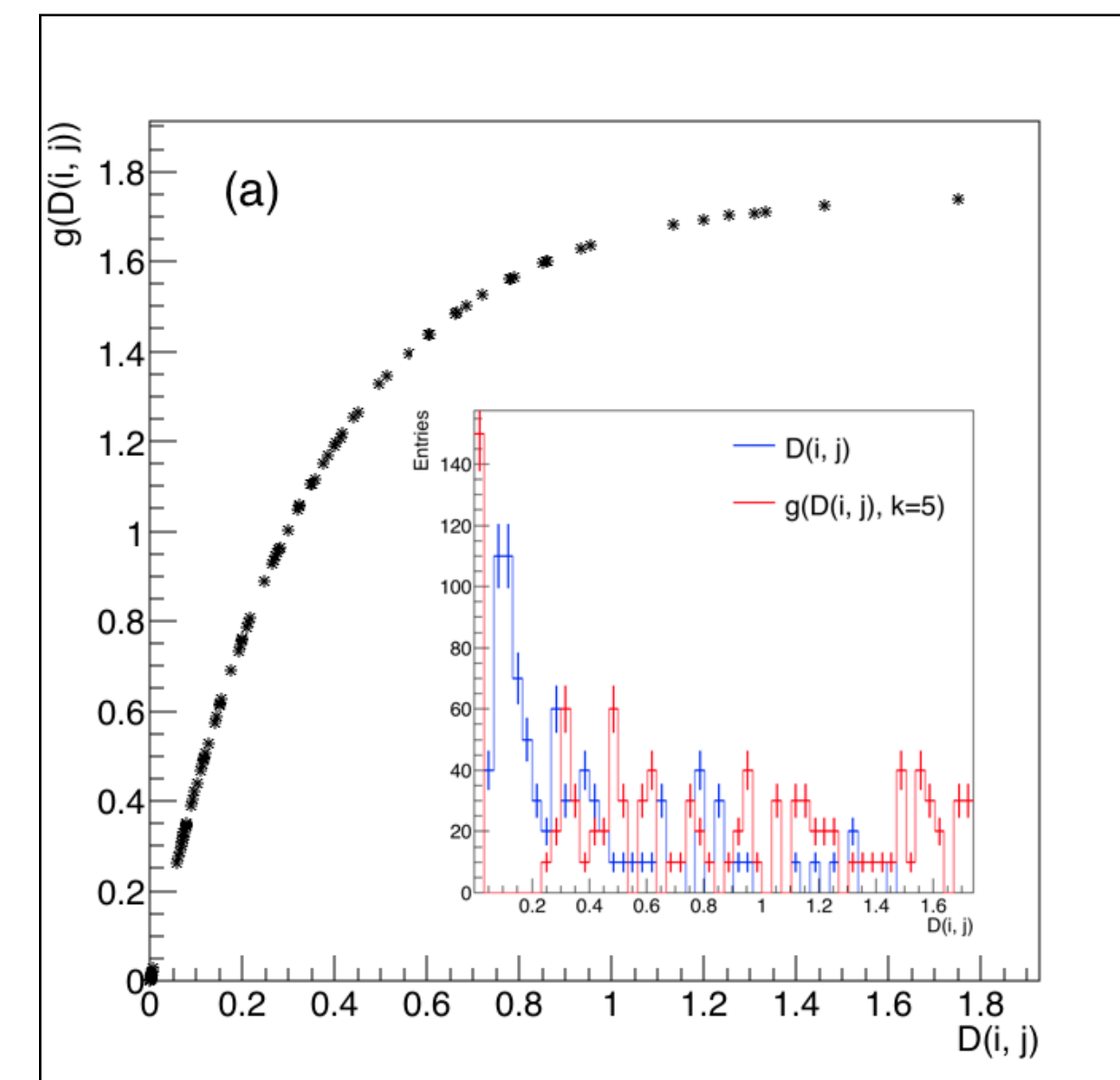
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• We use D-Wave’s default embedding algorithms

# Performance on one event with 3 p-p collision, 15 tracks

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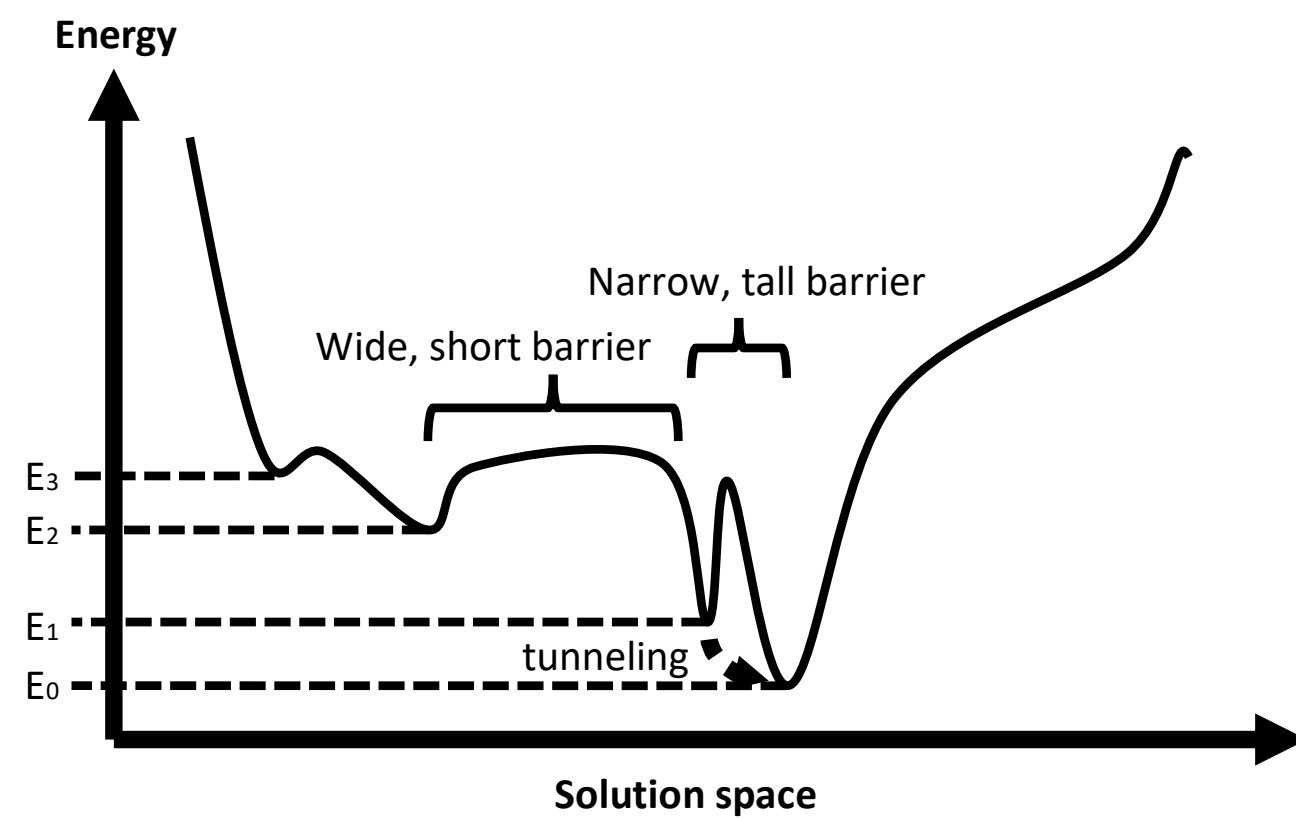
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- Algorithm tested on artificial events drawn from **measured LHC distributions of collision points** and **measured CMS distributions of tracks**
- **Realistic track reconstruction uncertainties** used. *CMS Collaboration, JINST 9 (2014) P10009*

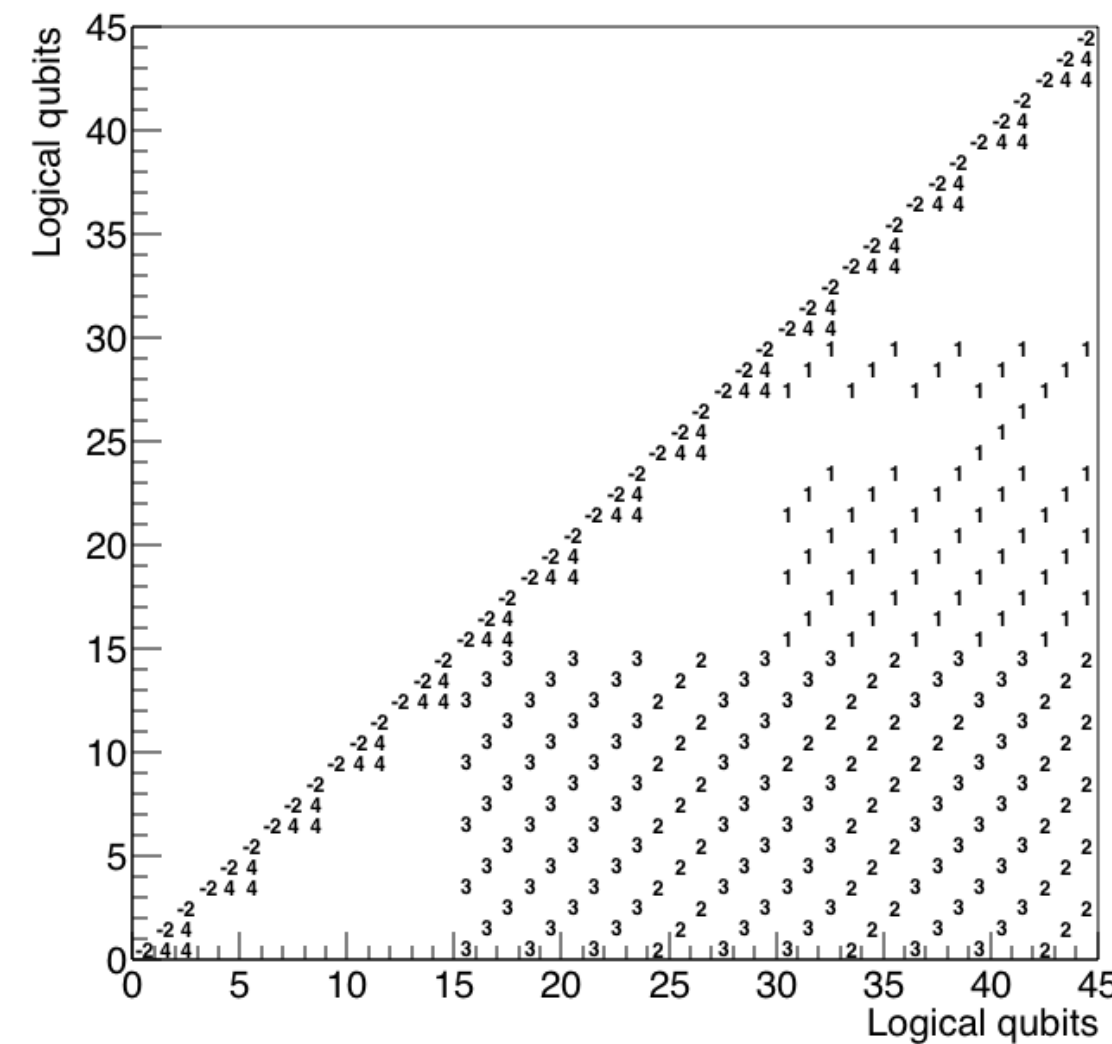
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Cartoon of energy profile in solution space.  
Tunneling is easy through tall narrow barriers.  
Classically difficult

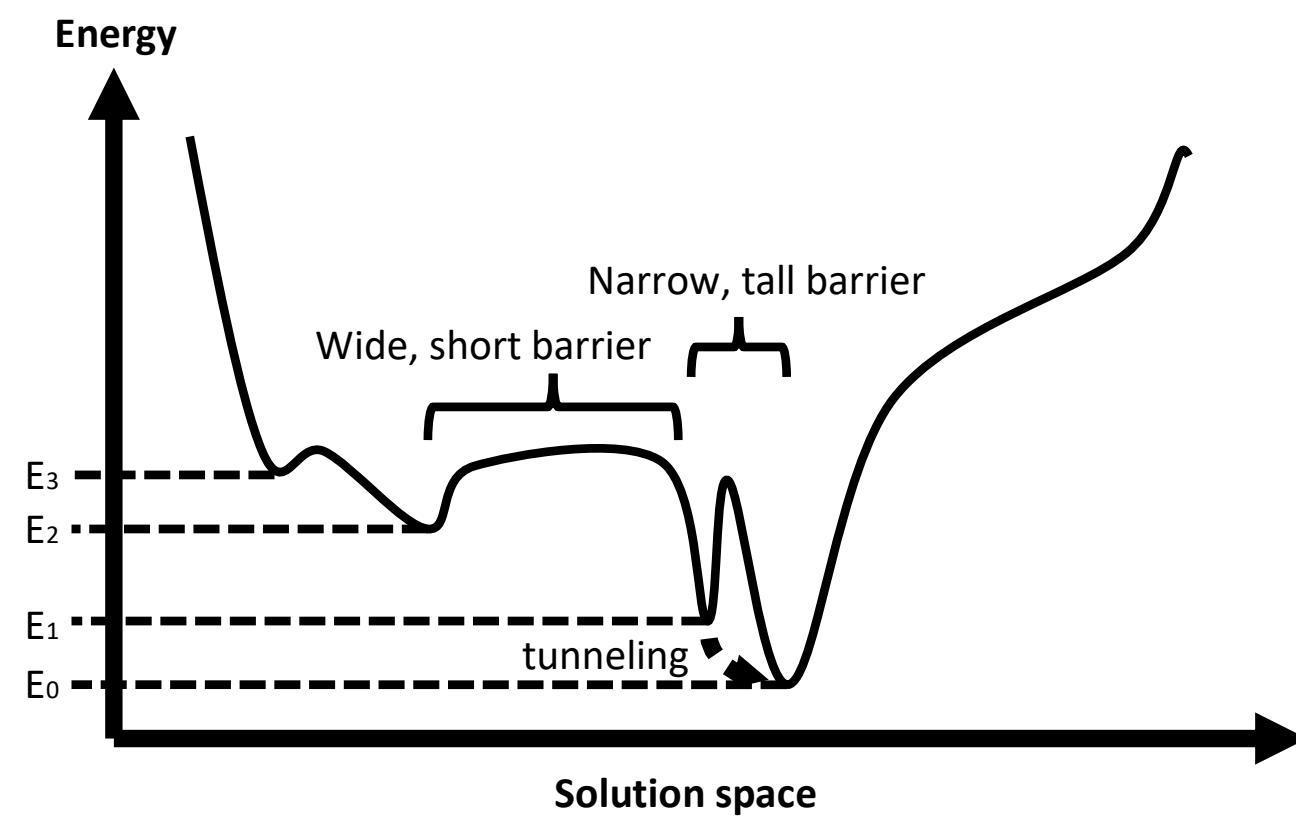


Coefficients of QUBO form used to solve  
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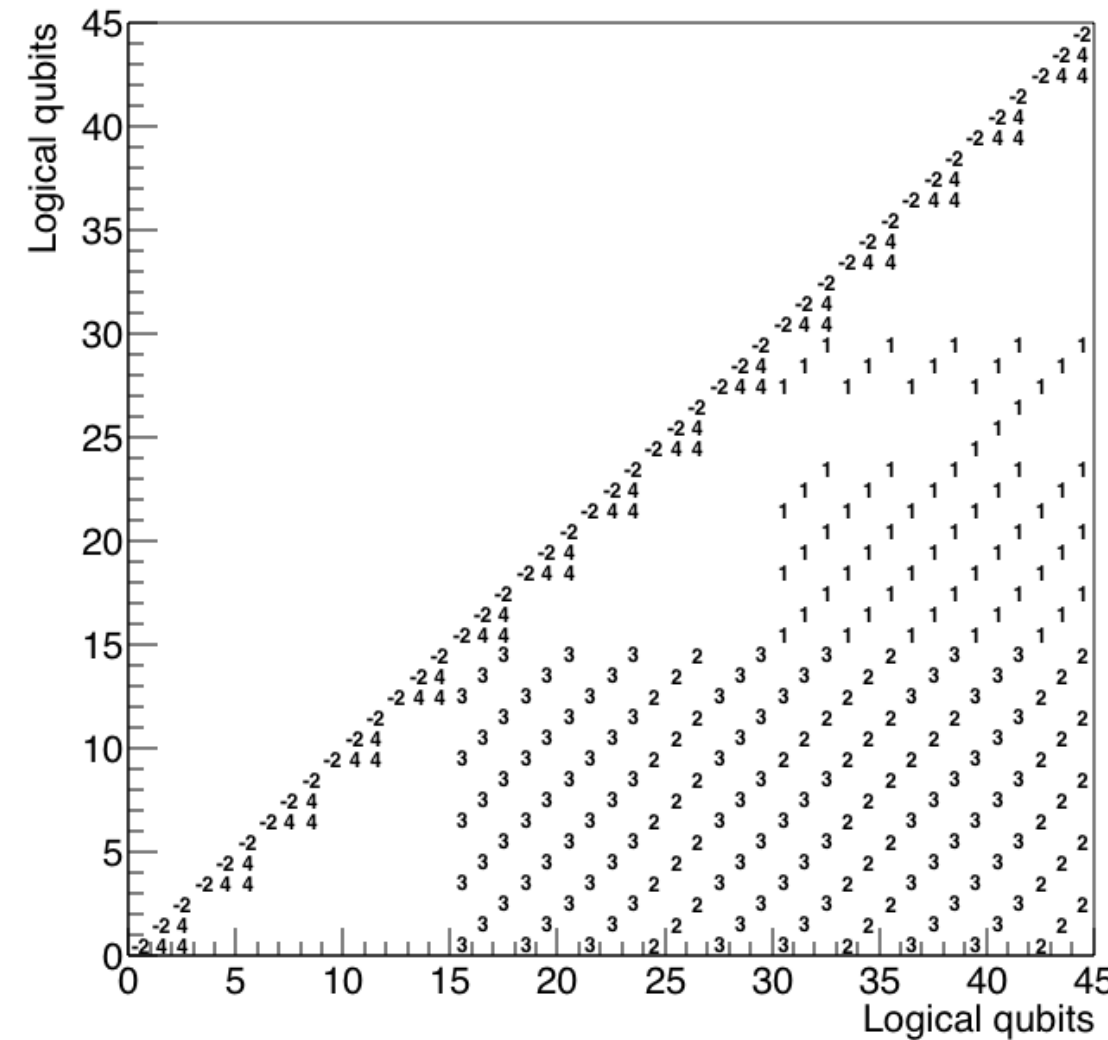
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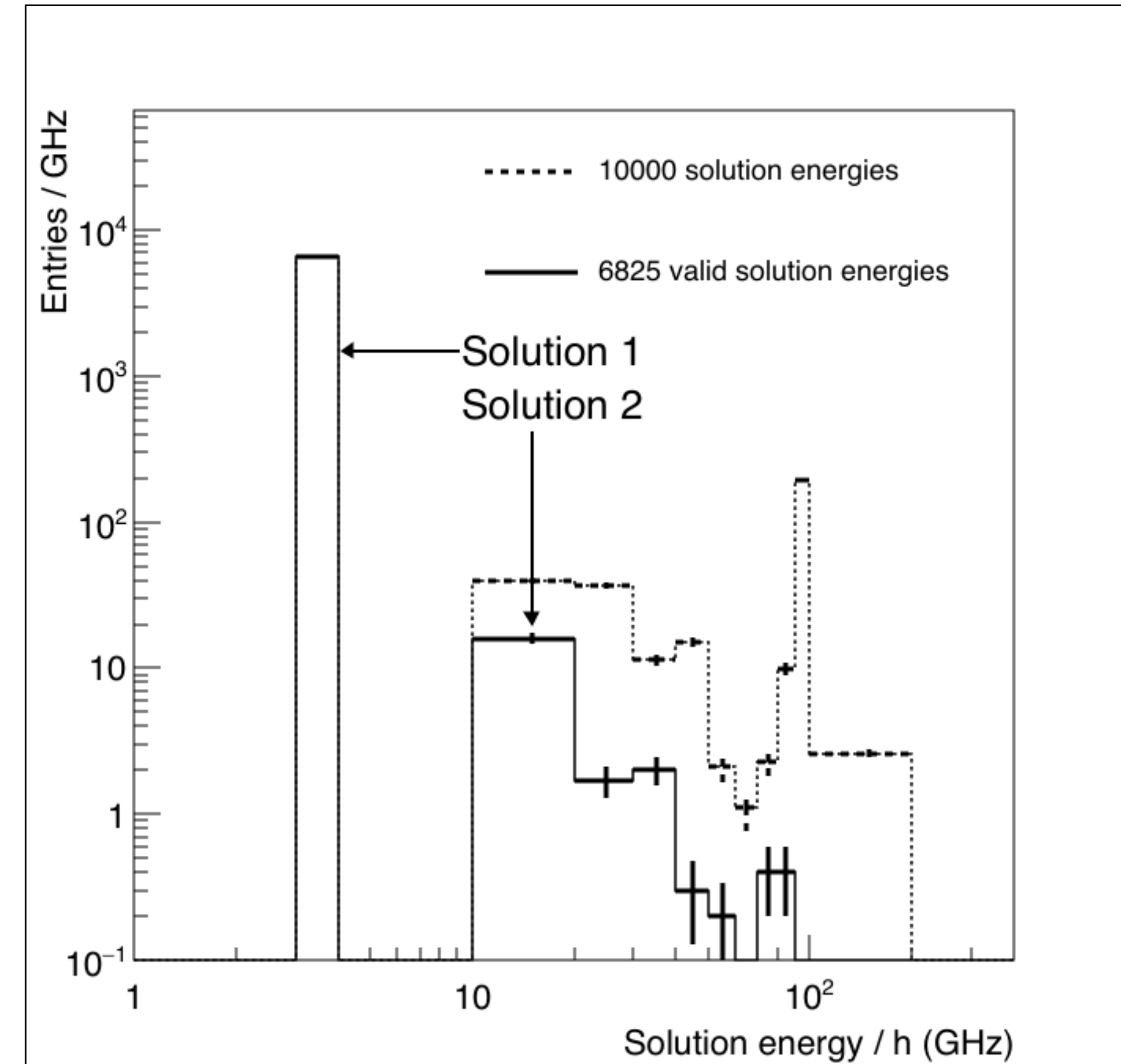
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Energy spectrum of solutions for one event with 3 p-p collisions and 15 tracks explored by the D-Wave 2000Q\_2\_1 with 10,000 samples. Energies corresponding to valid solutions, where  $p_{ik}$  add up to 1 for every track, are plotted with solid lines while invalid solutions are plotted with dashed lines. Error bars correspond to statistical uncertainties. The best and next-to-best valid solutions are indicated as Solutions 1 and 2, respectively. For clarity, the histogram is binned by 1 GHz below 10 GHz, by 10 GHz for 10 — 100 GHz, and by 100 GHz above 100 GHz

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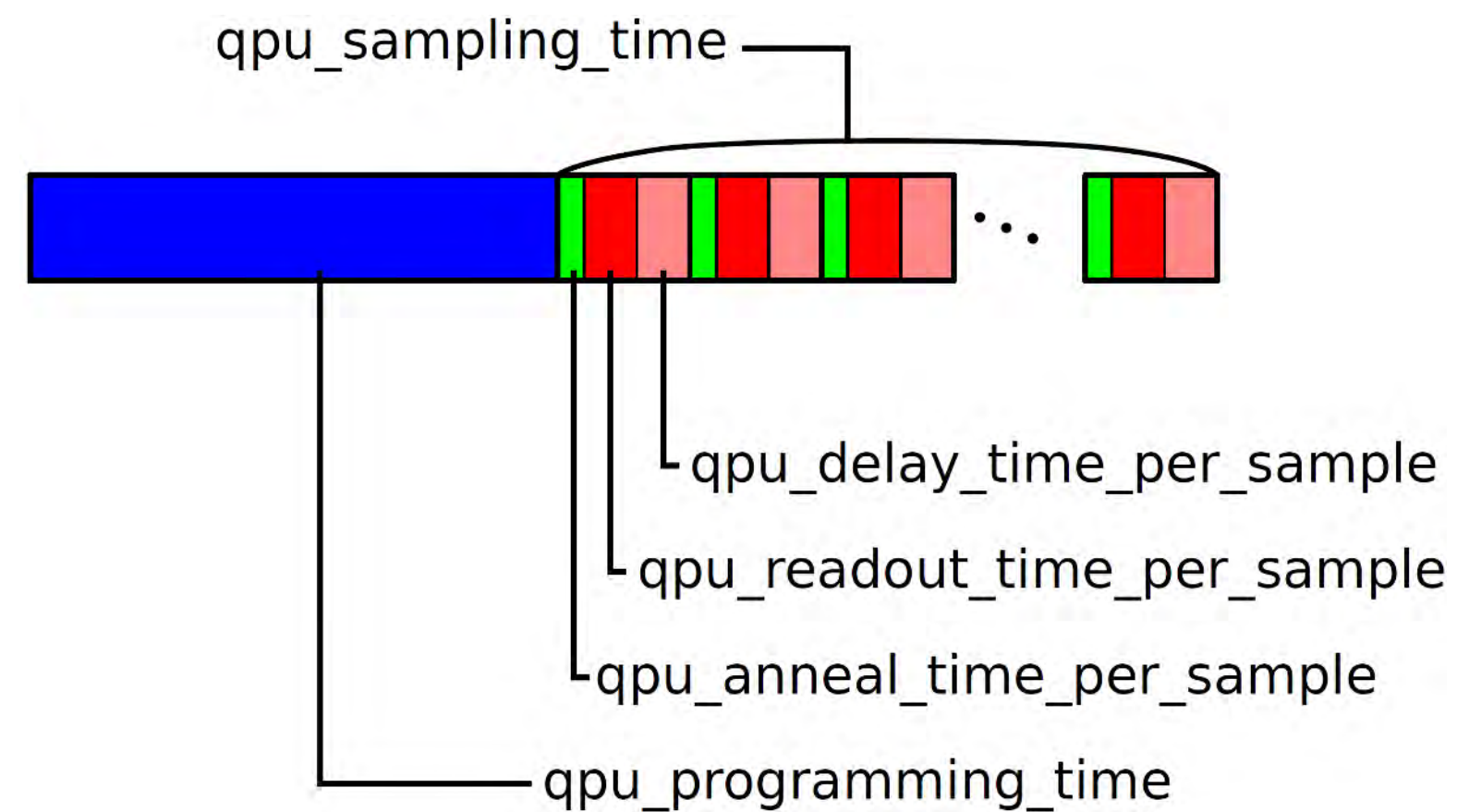
- QUBO bias terms are equal and come from the  $\lambda$  constraint
- Quantum state prepared and annealed 10,000 times. **DW\_2000Q\_2\_1** used
  - 6,825 solutions are valid, i.e.  $\lambda$  constraint is strictly met ( $\sum_k p_{ik} = 1$  for all tracks)
  - 6,615 solutions are correct (Solution 1). **Convergence efficiency = 66%**
  - Small number of valid secondary solutions where one track has been misassociated

# Benchmarking against simulated annealing in equal time

**Principle: Equalize working time between CPU and QPU, and compare convergence efficiency**

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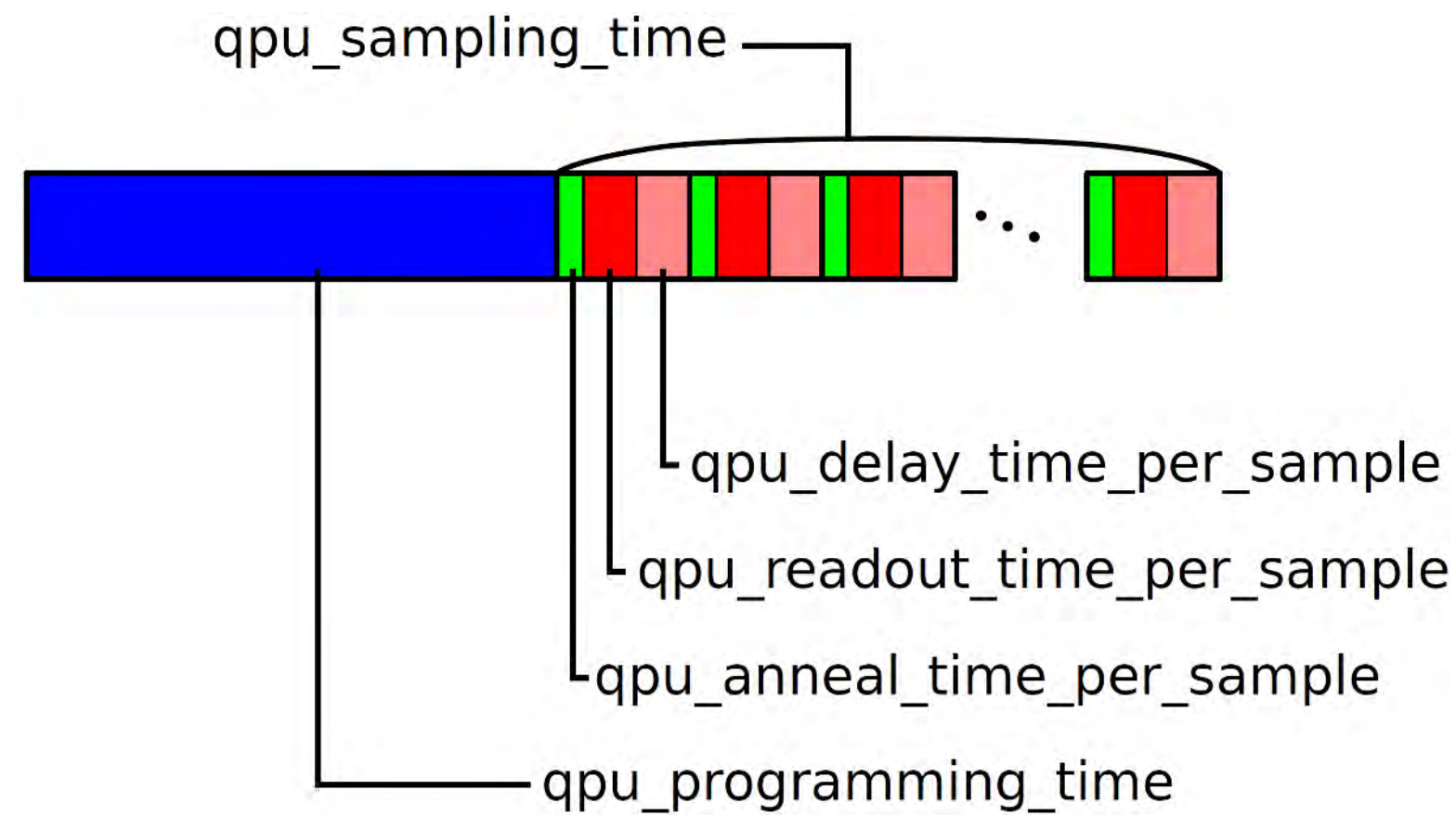
## Sampling time of D-Wave 2000Q\_2\_1

- Total sample time = **164  $\mu$ s**
  - Anneal time = 20  $\mu$ s
  - Readout time = 123  $\mu$ s
  - Delay time = 21  $\mu$ s
- How many Simulated Annealing sweeps can we fit in this?
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**CPU:** 3.1 GHz Intel Core i7-5557U (MacBook pro 2017)

**Algorithm:** Simulated annealing. Time optimized

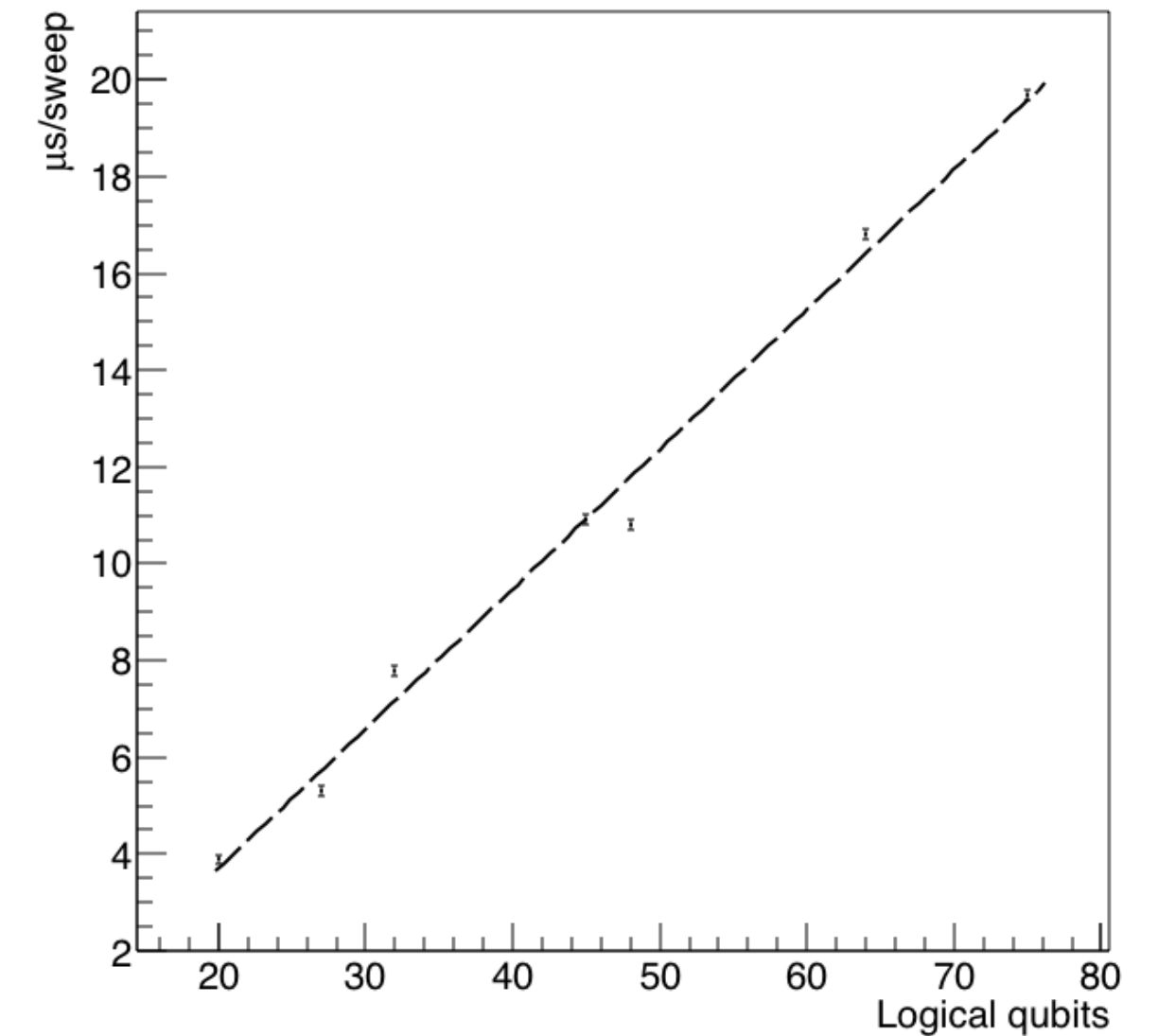
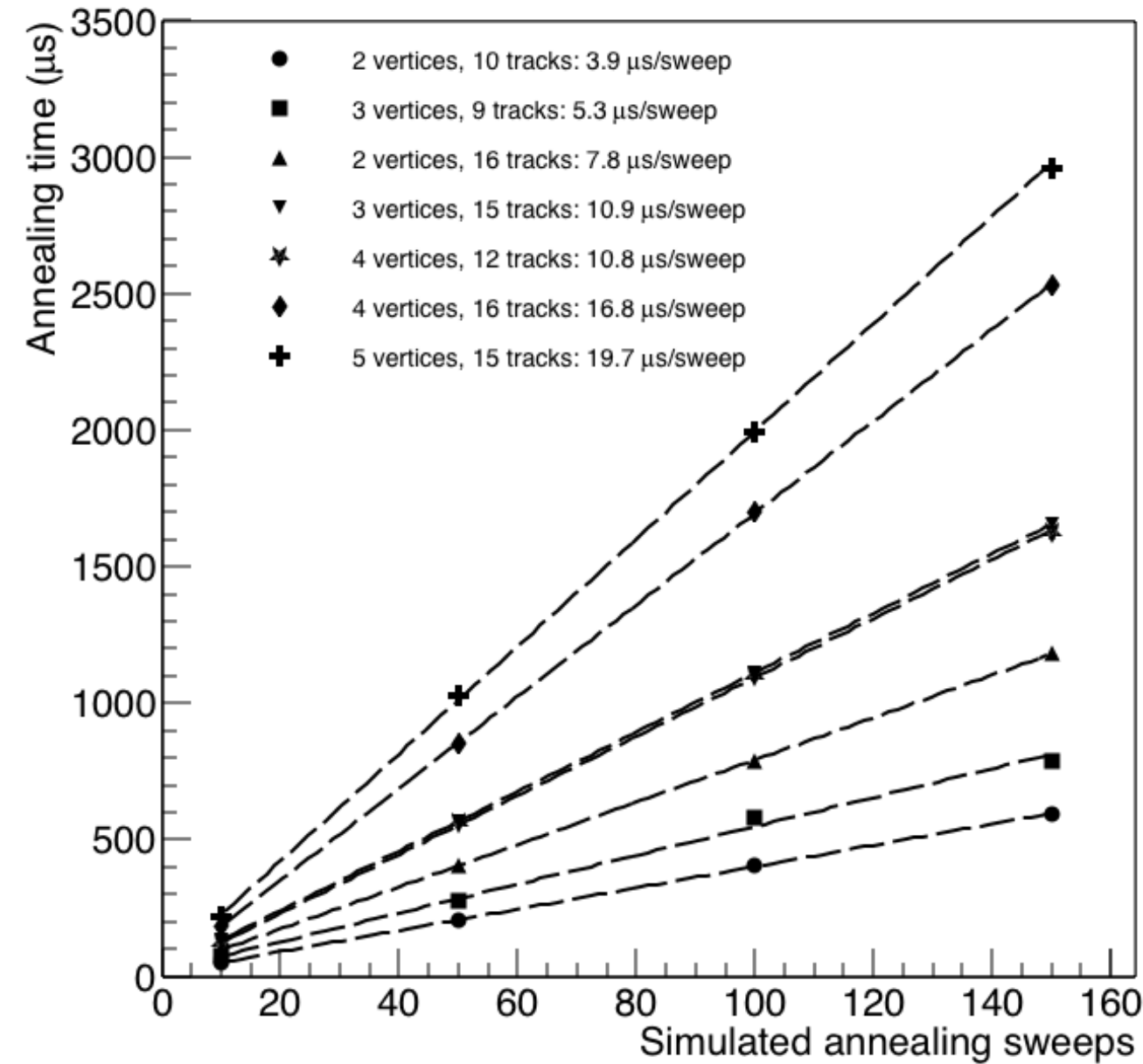
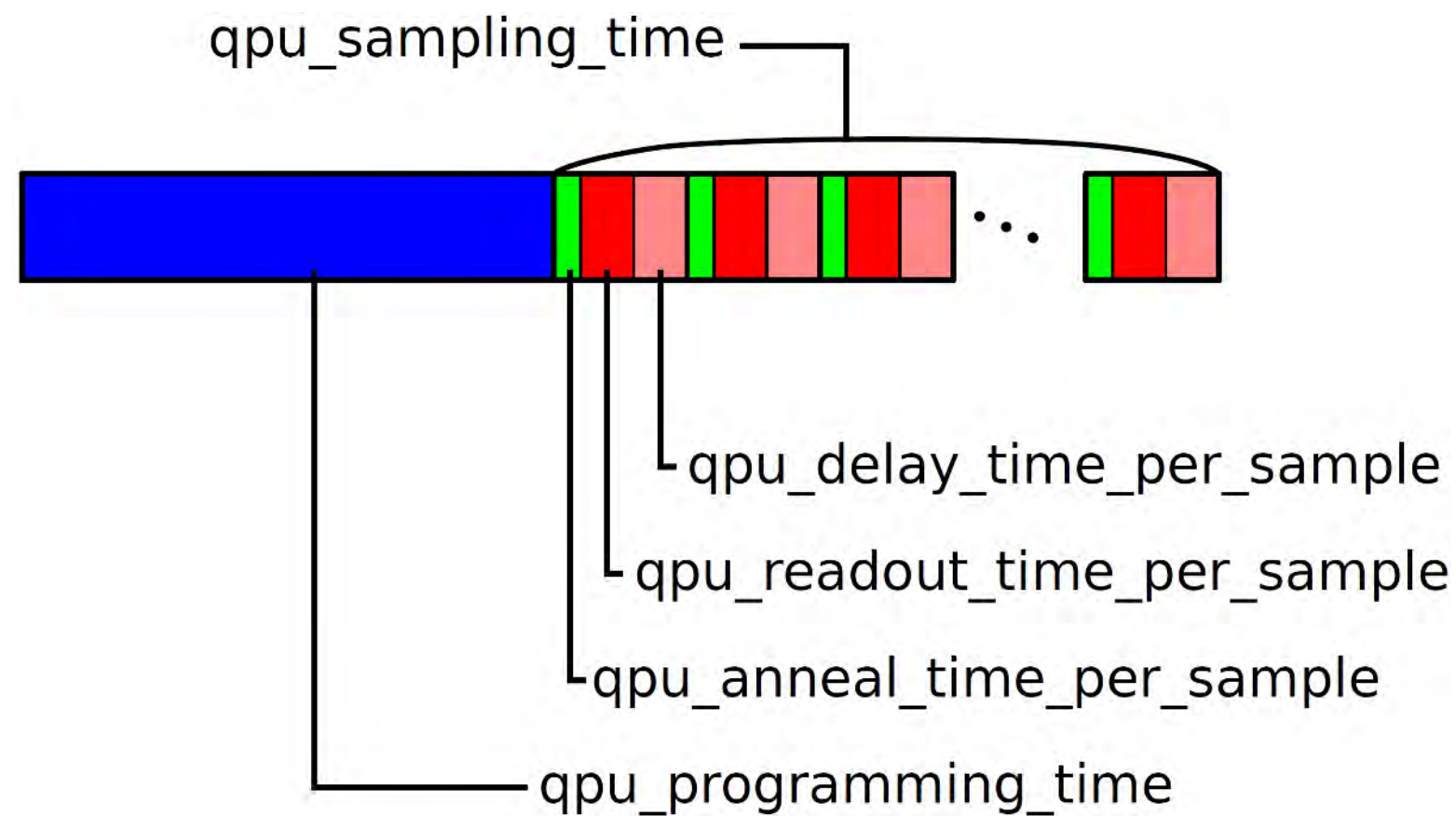
- Use a sorted `std::map` with keys = bit index, value = list of other bits it couples to and the coupling
- Bit flip only requires to compute energy difference

**Compiler:** C++, -O2 optimization



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Time per sweep scales linearly with number of bits involved

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  - Anneal time = 20 μs
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- How many Simulated Annealing sweeps can we fit in this?
  - Depends on problem complexity

CPU process time against number of sweeps for various event topologies under consideration

**CPU:** 3.1 GHz Intel Core i7-5557U (MacBook pro 2017)

**Algorithm:** Simulated annealing. Time optimized

- Use a sorted `std::map` with keys = bit index, value = list of other bits it couples to and the coupling
- Bit flip only requires to compute energy difference

**Compiler:** C++, -O2 optimization

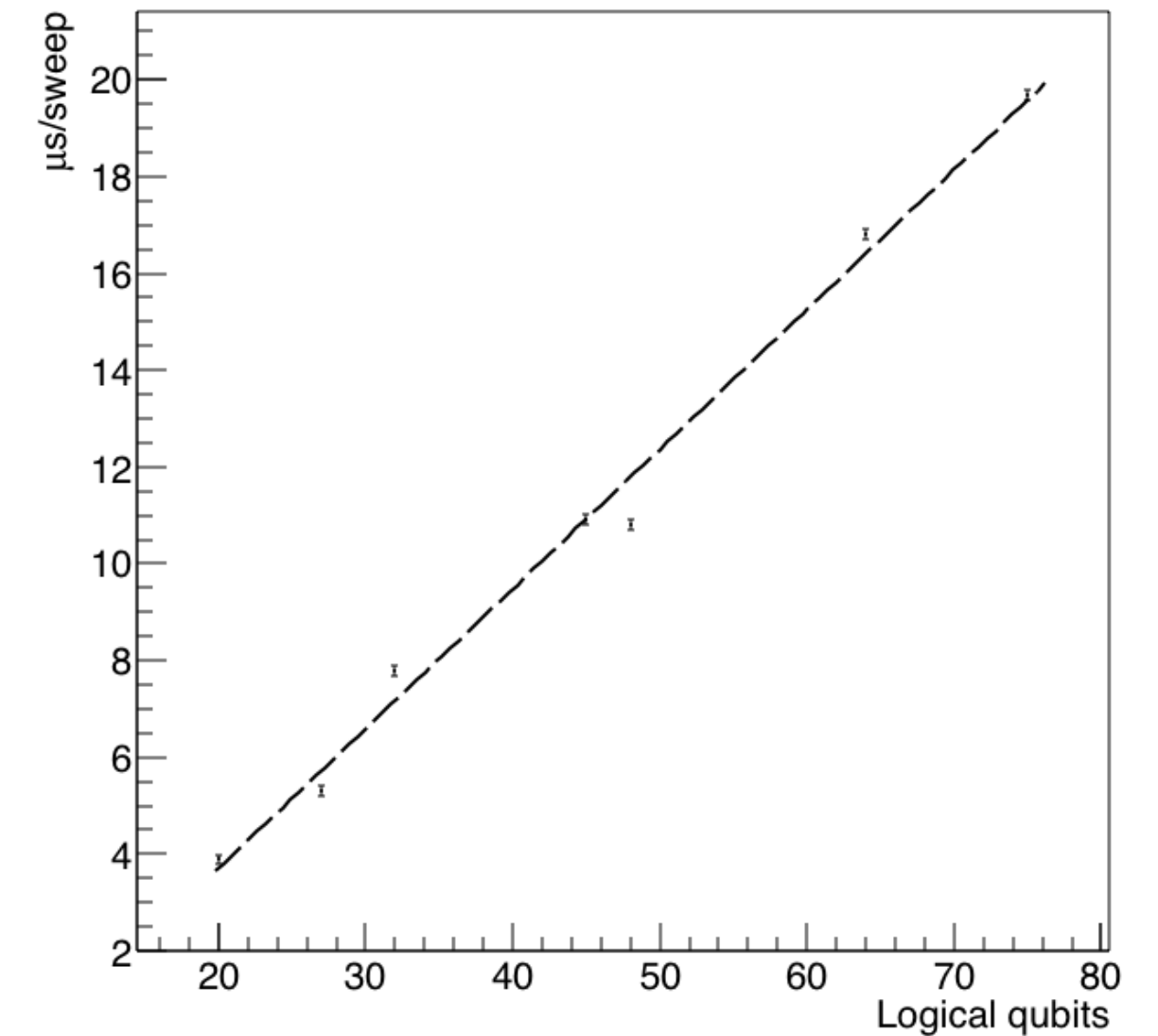
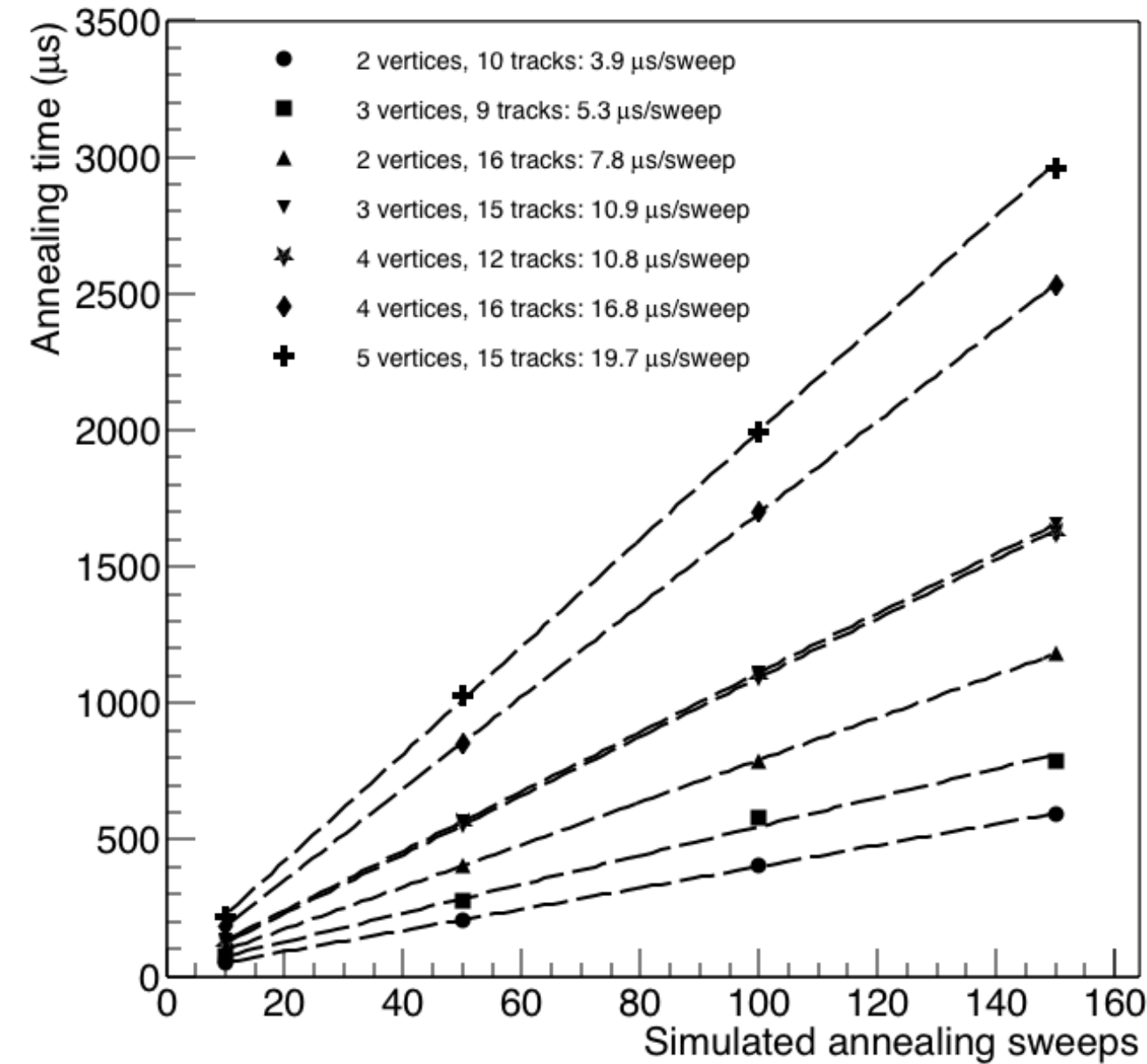
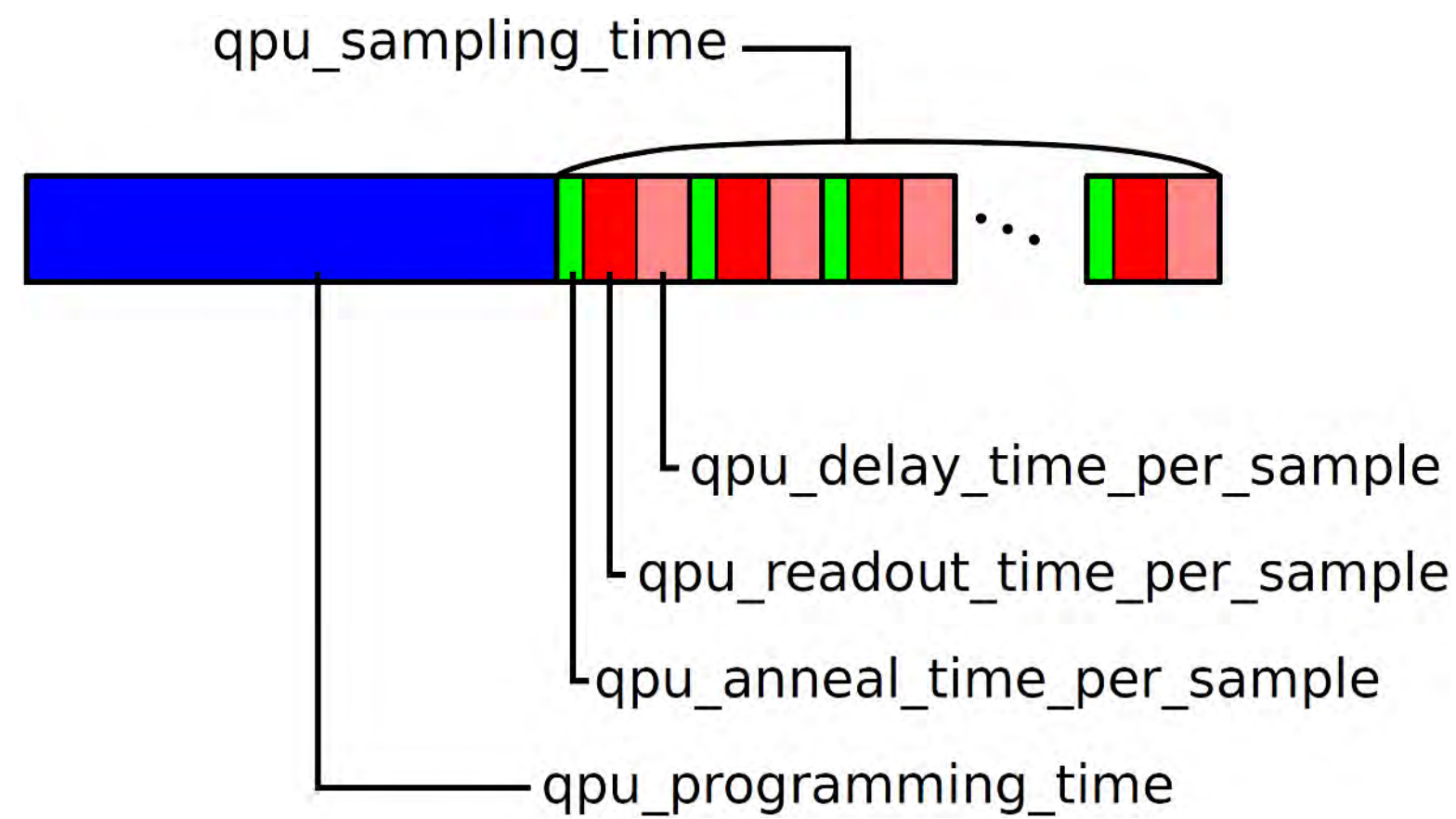


## Estimating CPU time per sweep

- Measure **process time**, not wall time
- Plot **process time against nSweeps** for different event topologies
- Discard overhead. **Slope** is time per sweep.
  - For 3 collision 15 tracks, 10.9 μs/sweep. Thus, 15 sweeps would fit in D-Wave's sampling time

# Benchmarking against simulated annealing in equal time

Principle: Equalize working time between CPU and QPU, and compare convergence efficiency



Time per sweep scales linearly with number of bits involved

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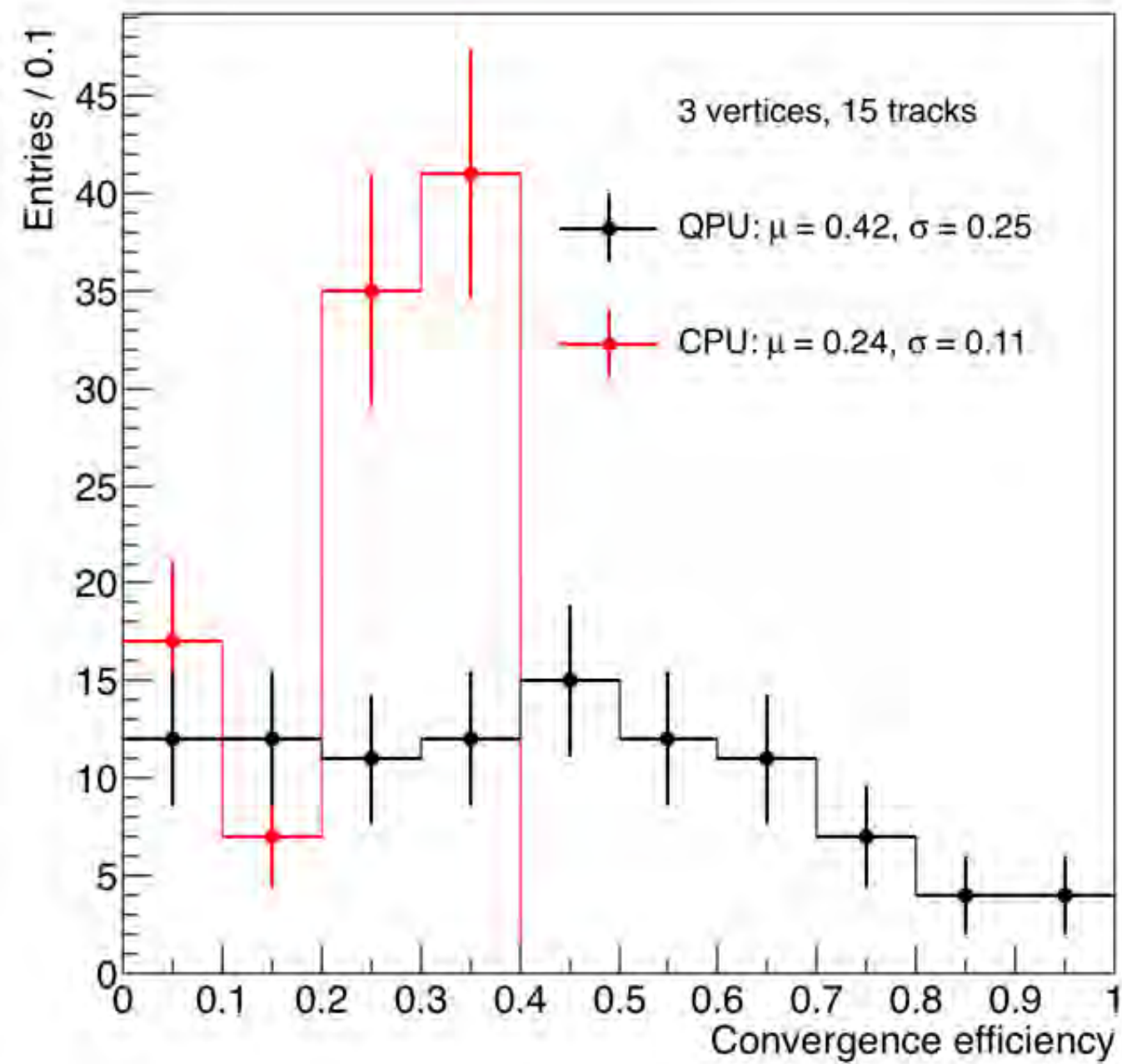


Simulated annealing on CPU is only allowed as many iterations between  $\beta_{init} = 0.1$  and  $\beta_{final} = 10$  as would fit 164 μs

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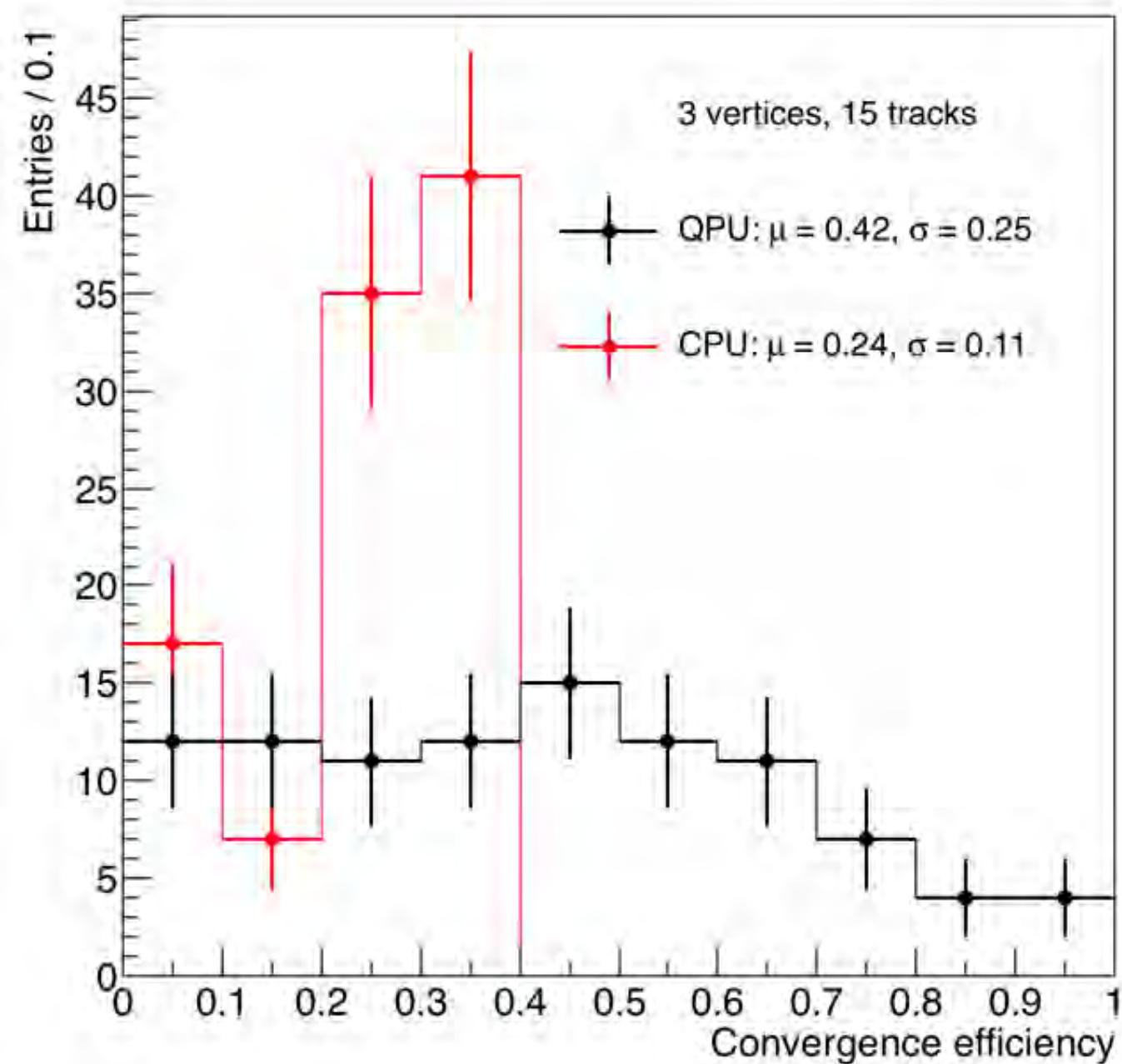
# Performance on 100 events with 3 p-p collisions and 15 tracks



## Performance on an ensemble of events

- 100 events with 3 p-p collisions and 15 tracks are thrown from measured CMS distributions
  - Each event is sampled 10,000 times by both the QPU and the CPU (in equivalent time)
- Events with *collisions spread closely compared to the spread of their tracks are hard* for both QPU and CPU to solve
- A distribution of convergence efficiencies is observed
  - QPU: mean = 42%, std. dev. = 25%
  - CPU: mean = 24%, std. dev. = 11%

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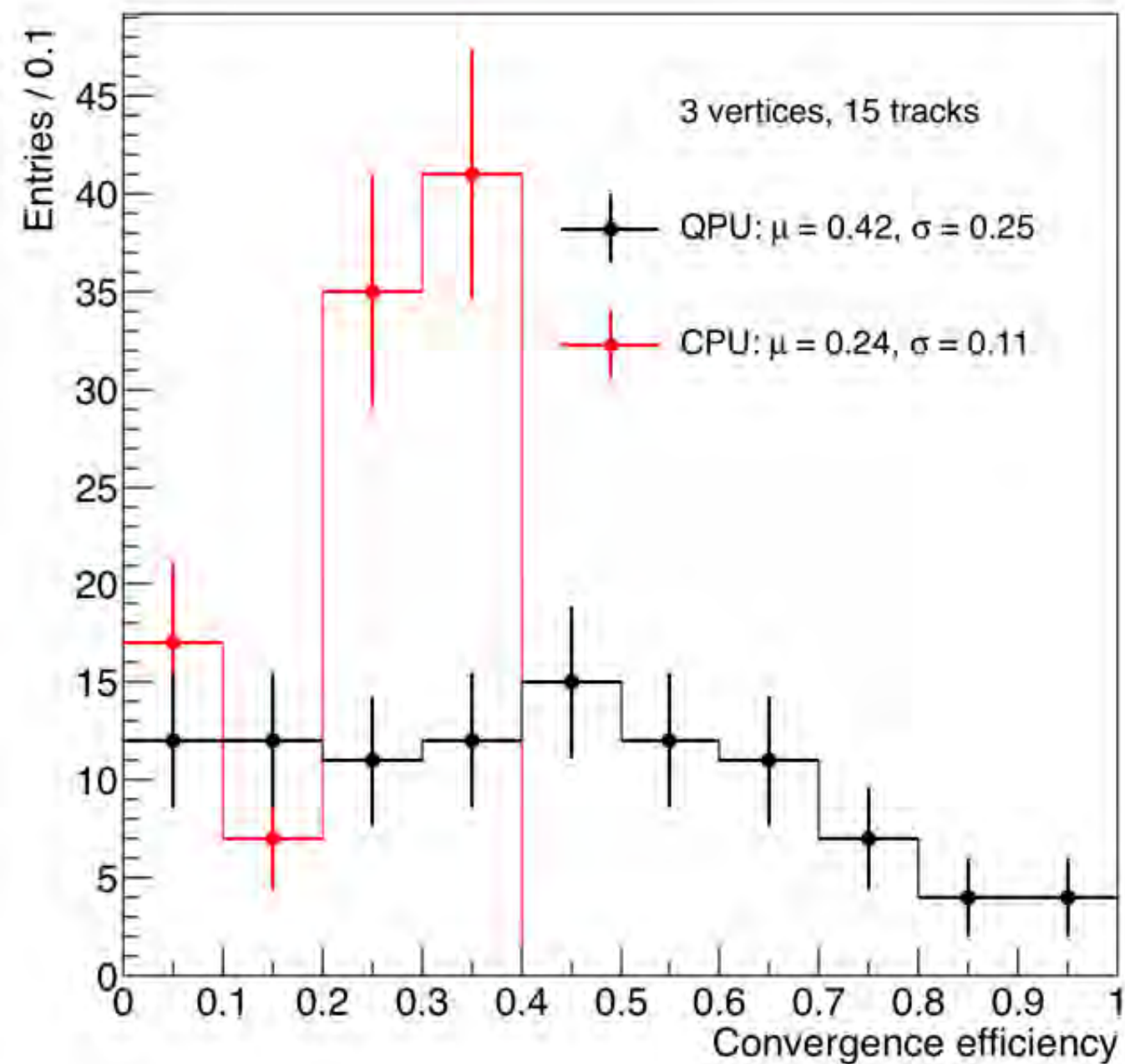


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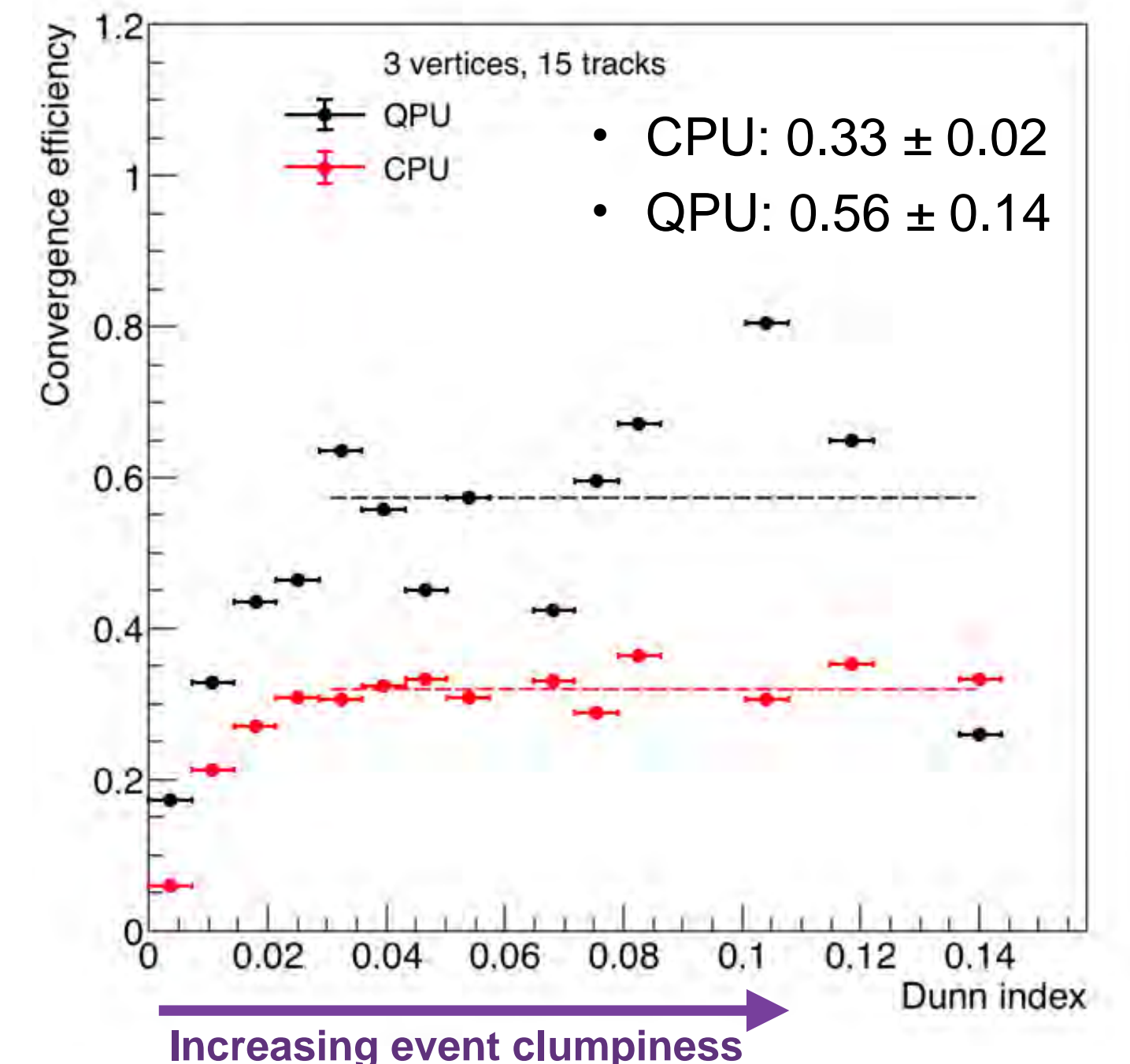
## Performance against event “clumpiness”

- A measure of event clumpiness is the Dunn index. Low Dunn index = diffuse event, high = clumpy event

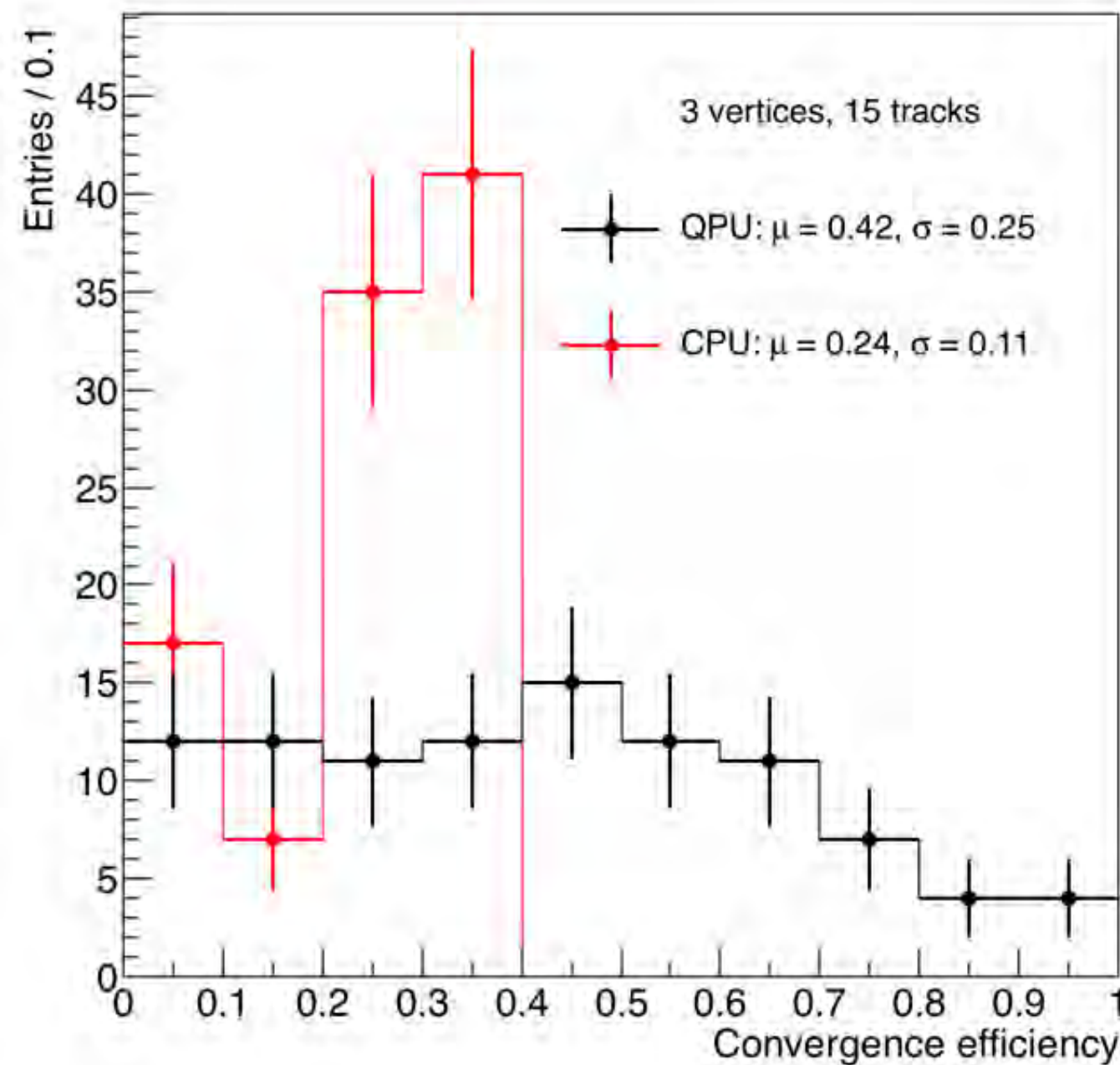
$$D = \frac{\min_{1 \leq i \leq j \leq n} d(i, j)}{\max_{1 \leq k \leq n} d'(k)}$$

numerator = minimum inter-cluster distance  
denominator = maximum intra-cluster distance between tracks

- Convergence efficiency is plotted as a function of Dunn index. Shows expected structure



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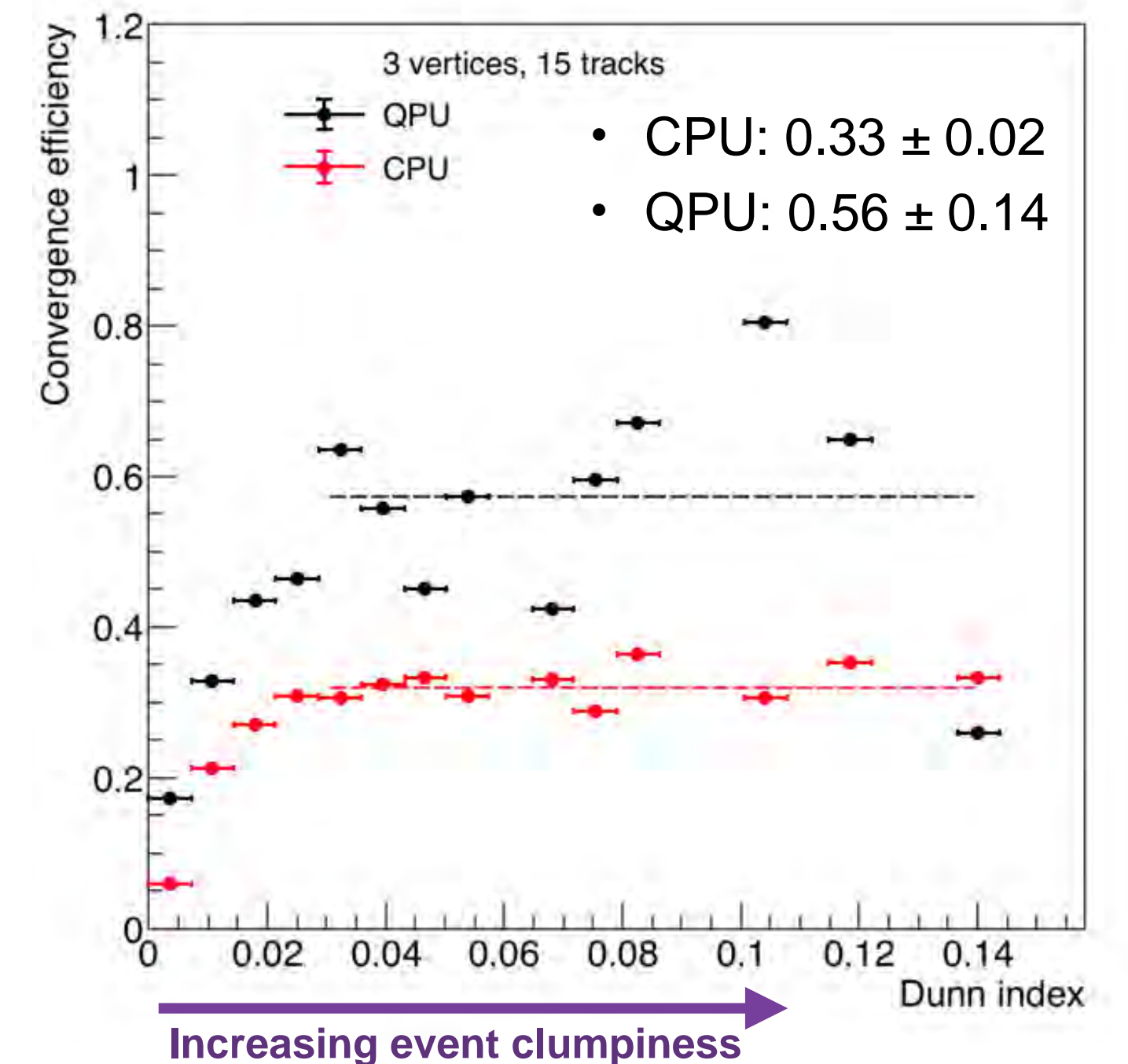
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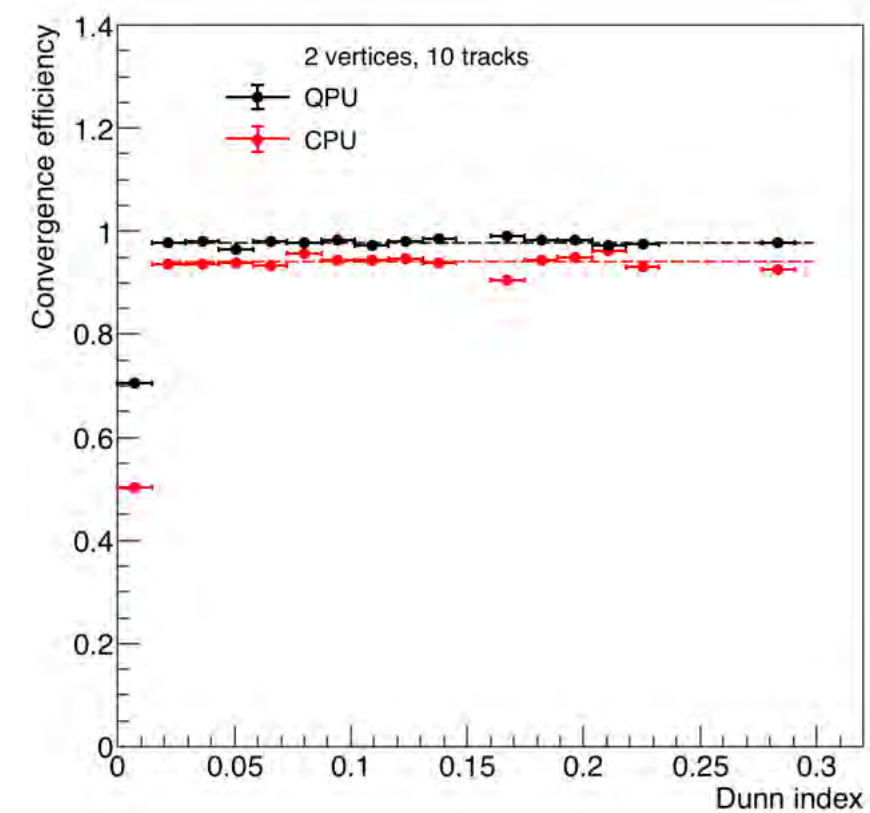
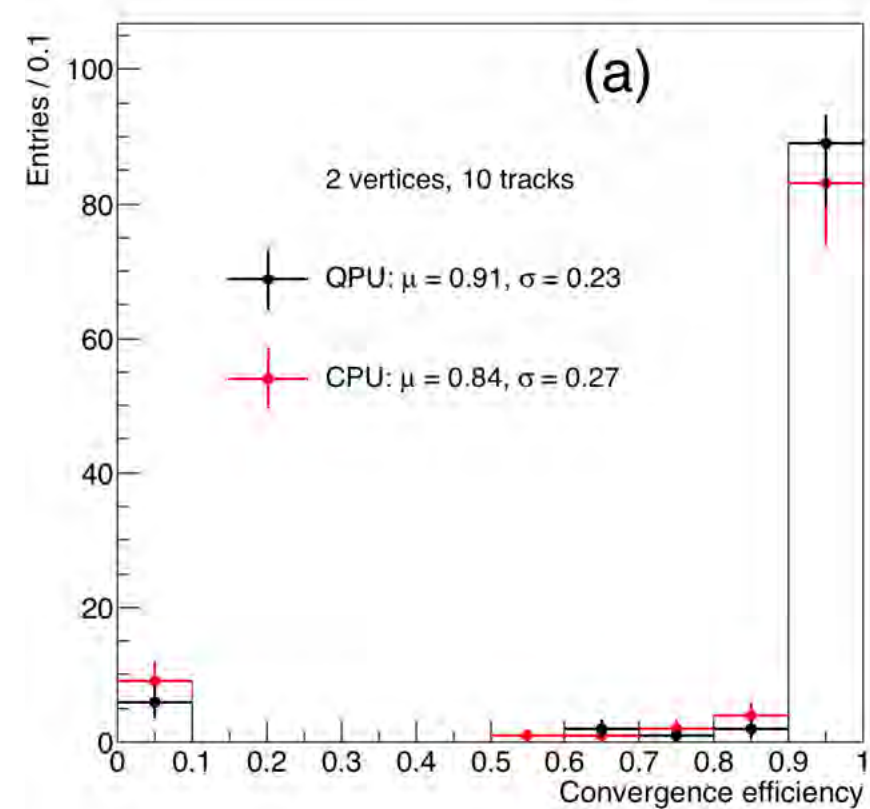


**Convergence efficiency increases with event clumpiness. QPU beats CPU in efficiency for same running time**

# QPU vs CPU scaling with event complexity

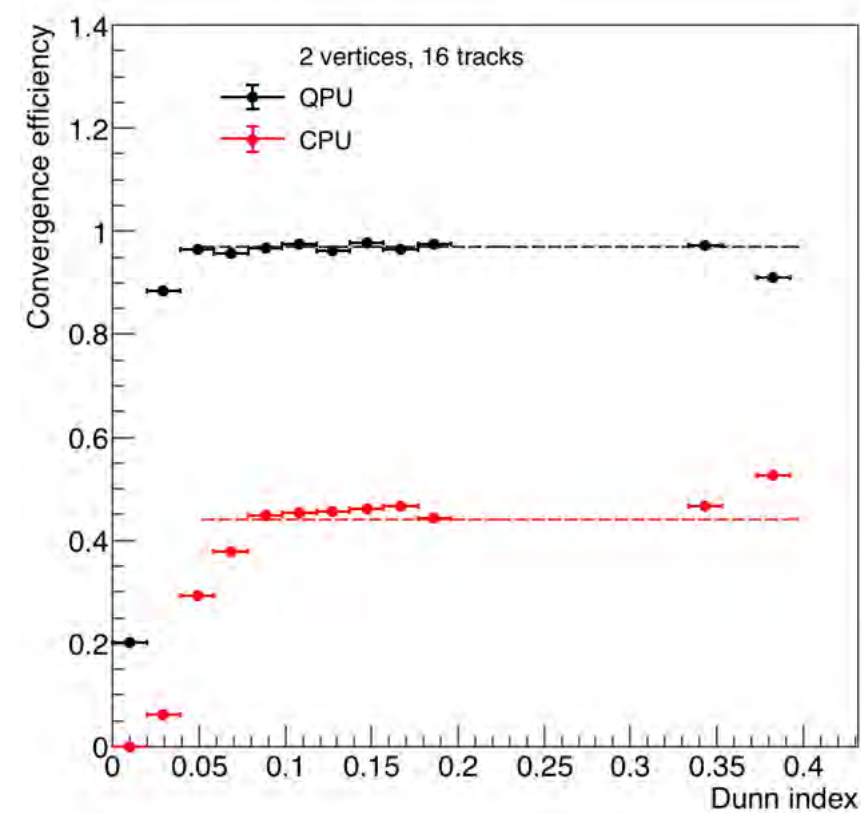
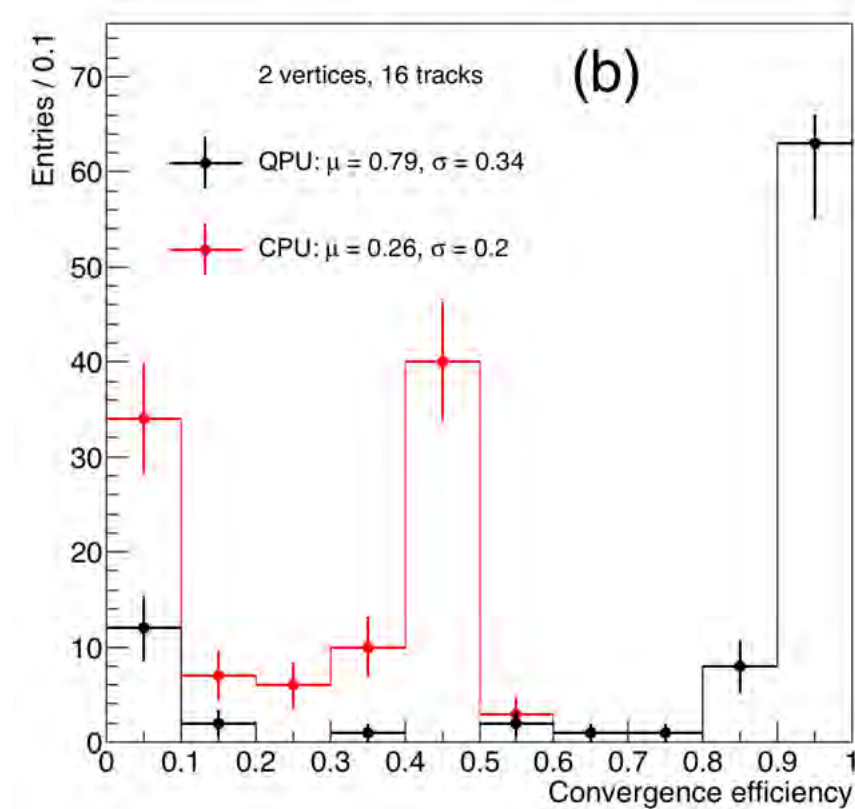
We scan over event topologies with increasing complexity

2 collisions, 10 tracks



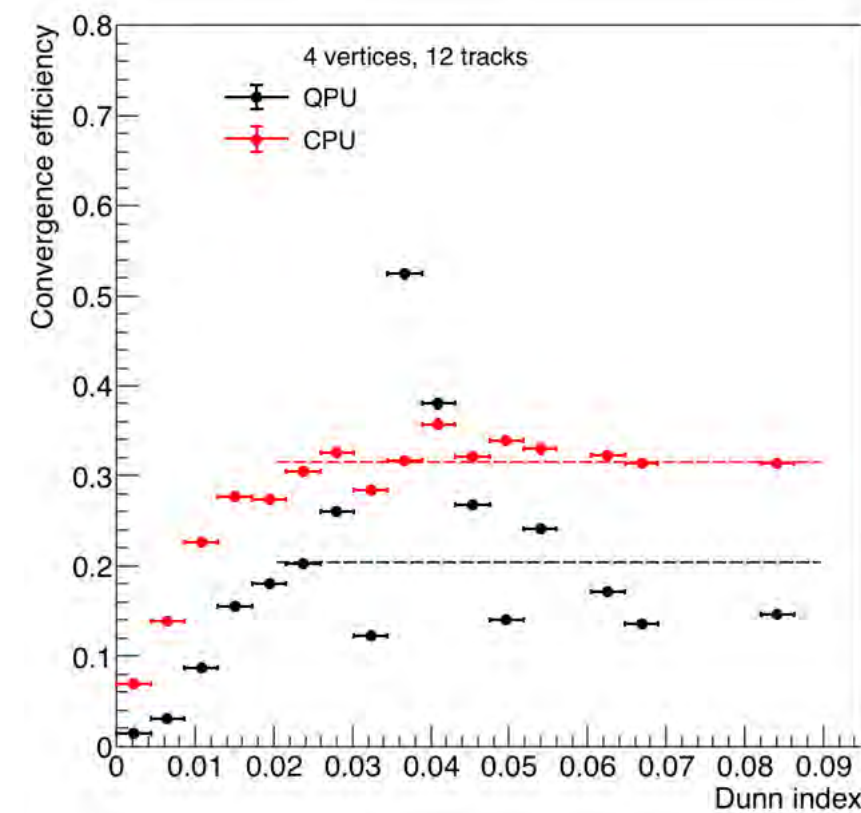
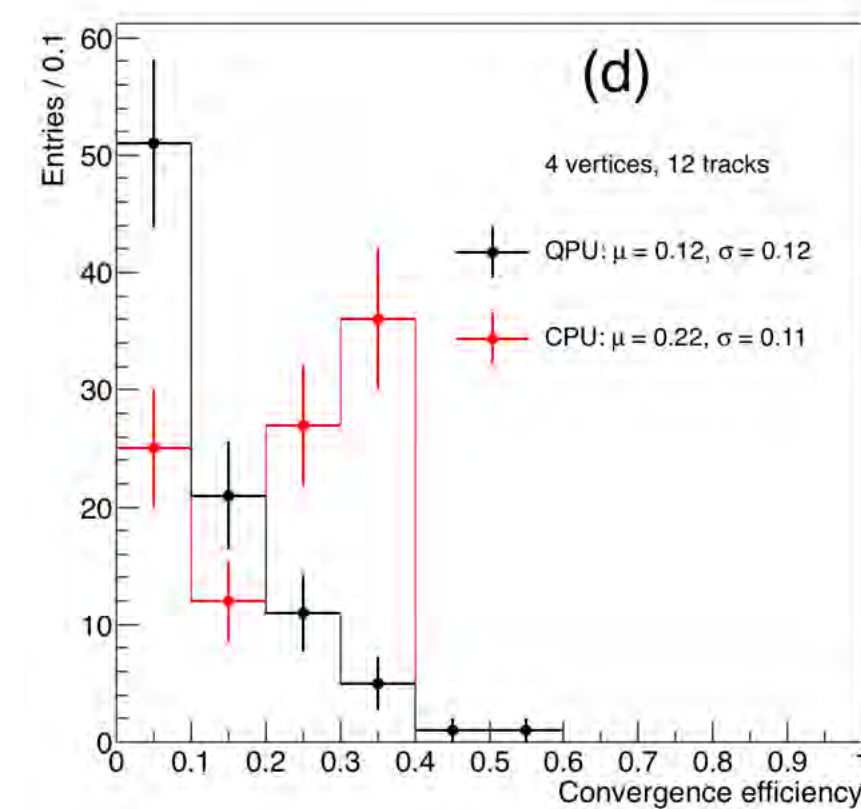
- CPU:  $0.94 \pm 0.01$
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2 collisions, 16 tracks



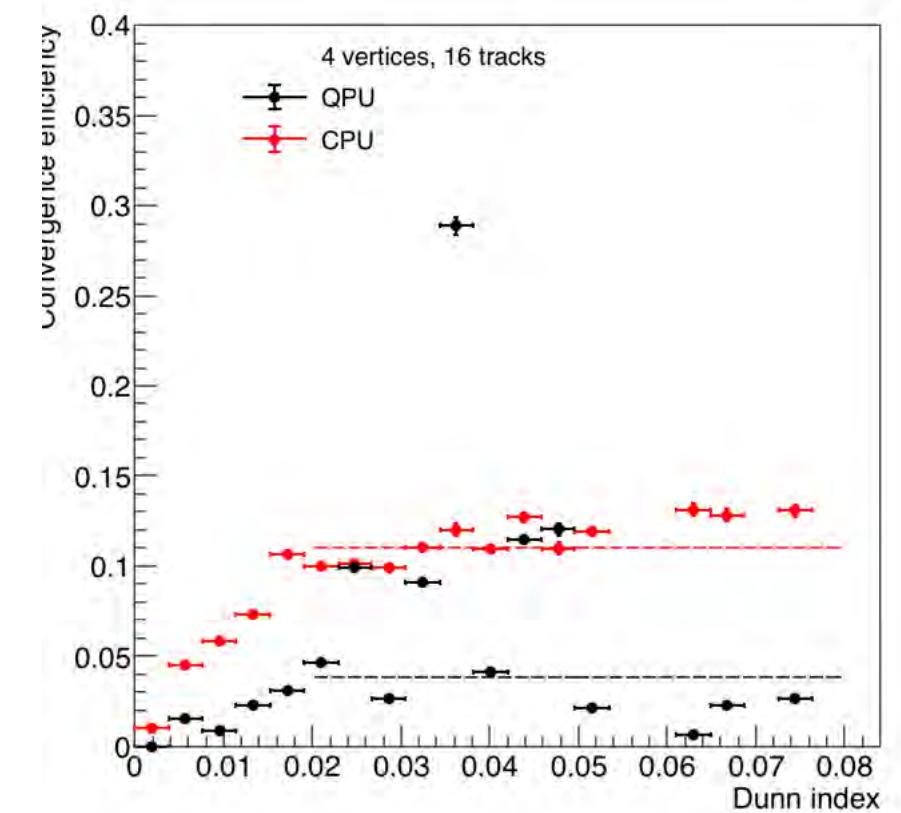
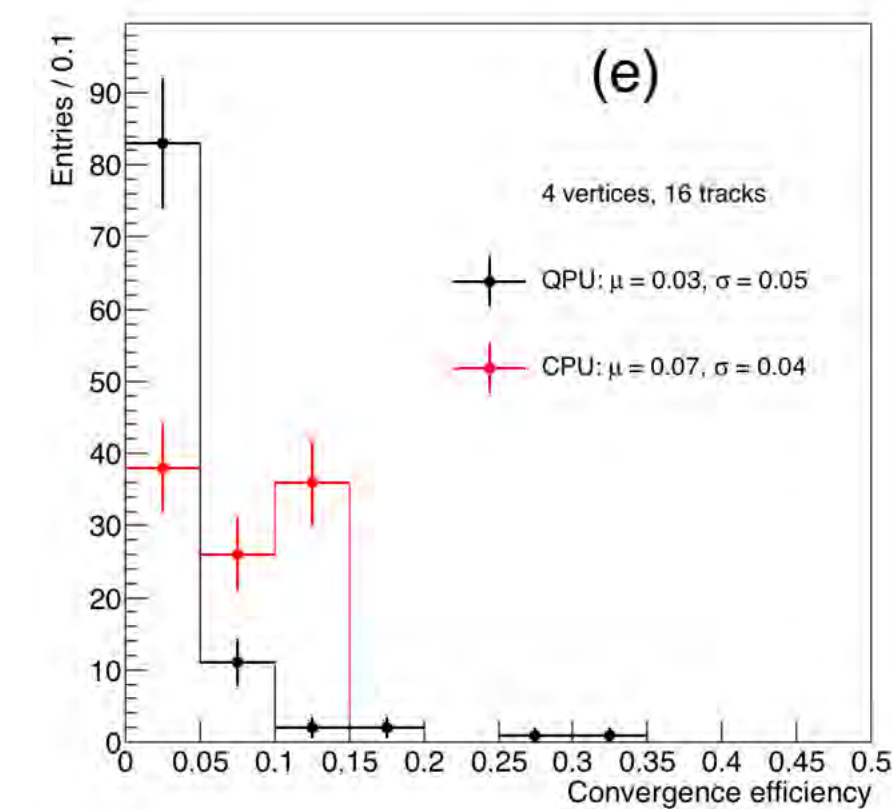
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4 collisions, 12 tracks



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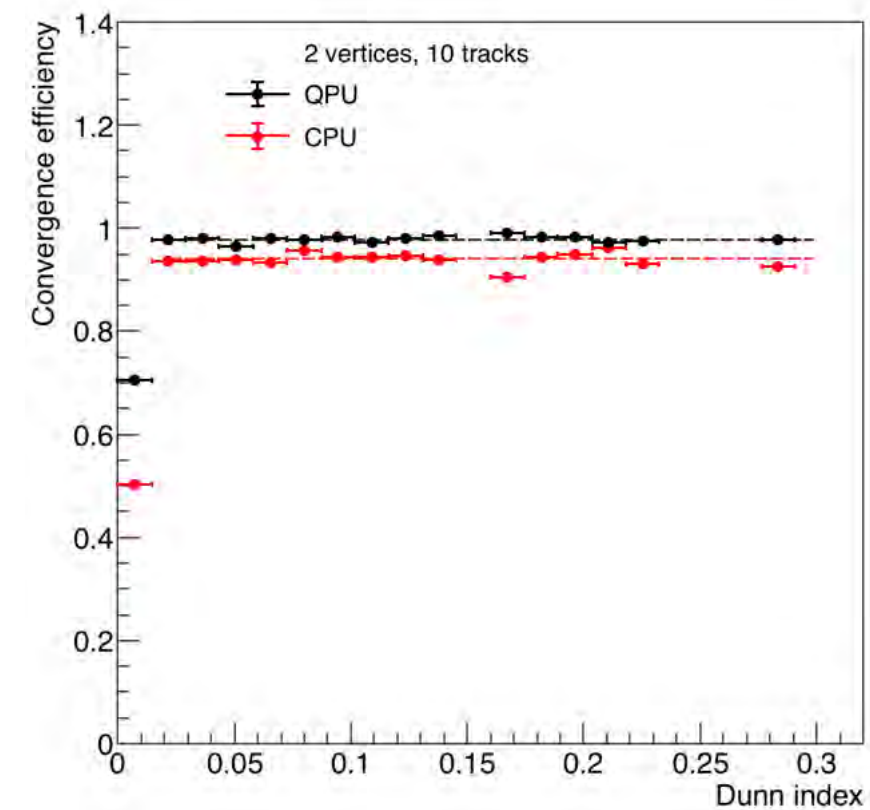
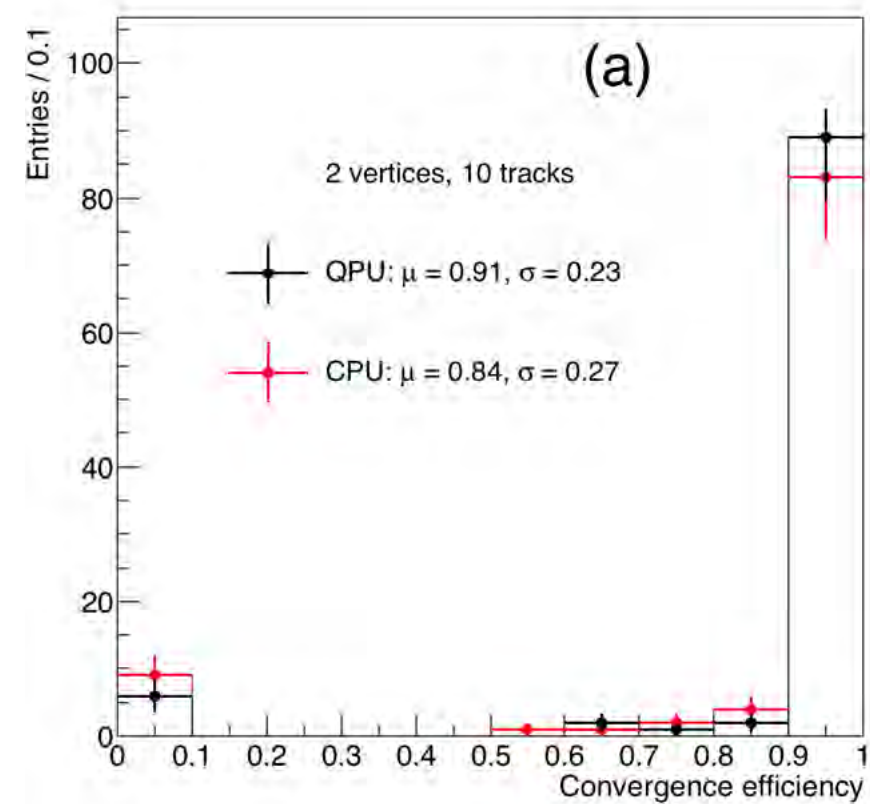
Maximum  
convergence  
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# QPU vs CPU scaling with event complexity

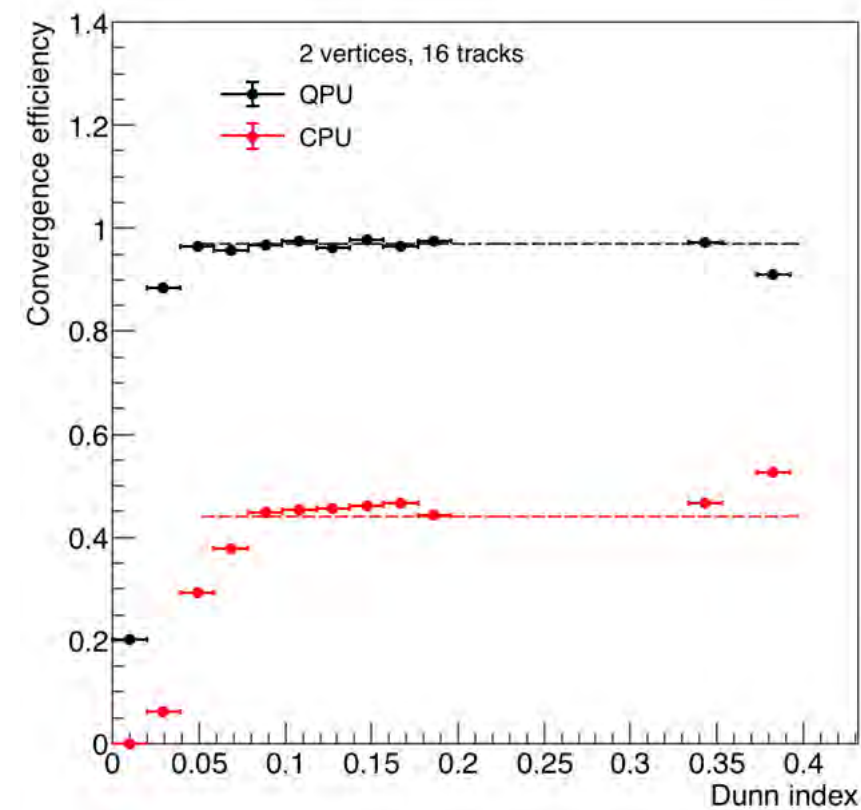
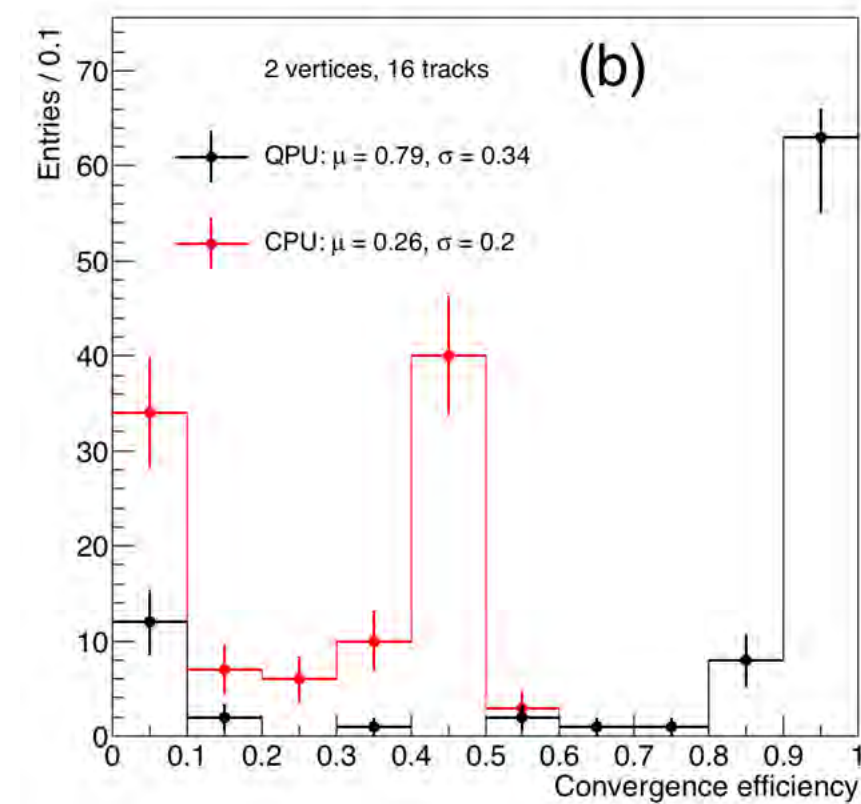
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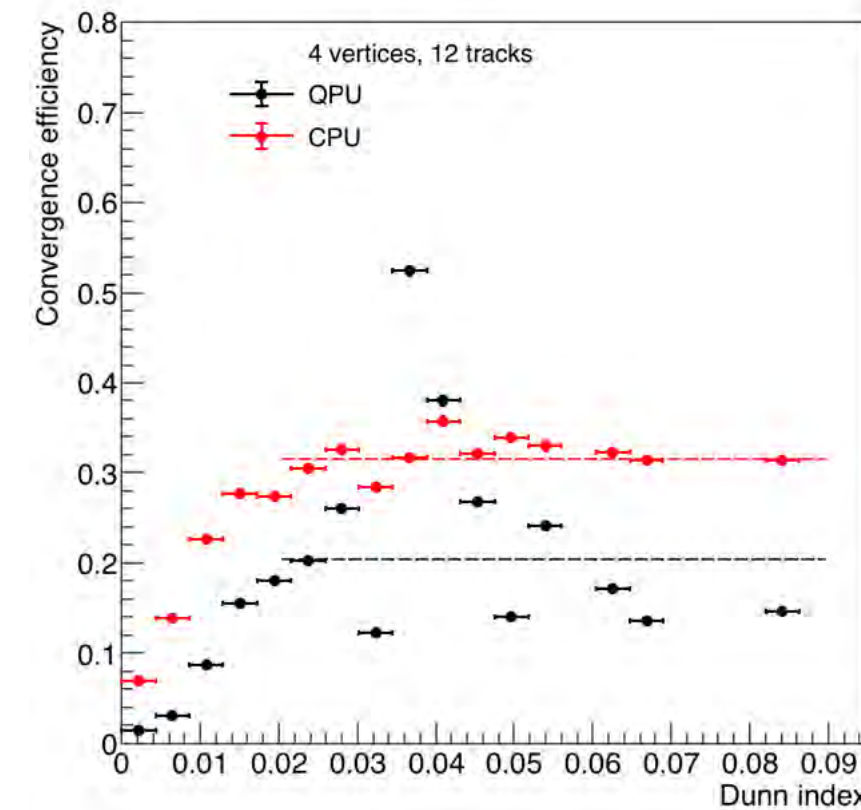
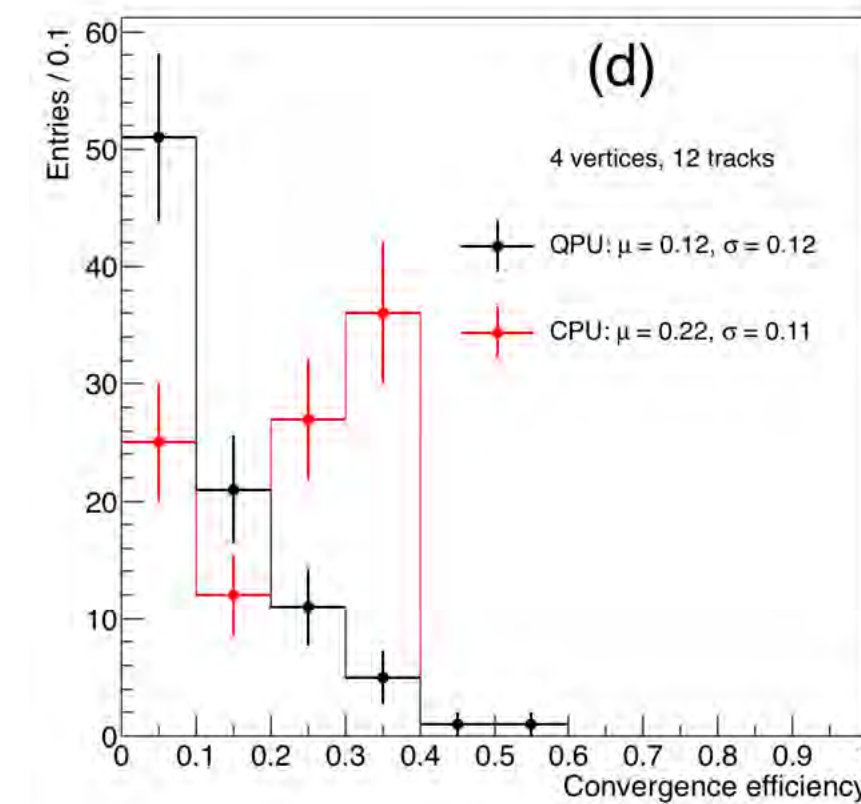
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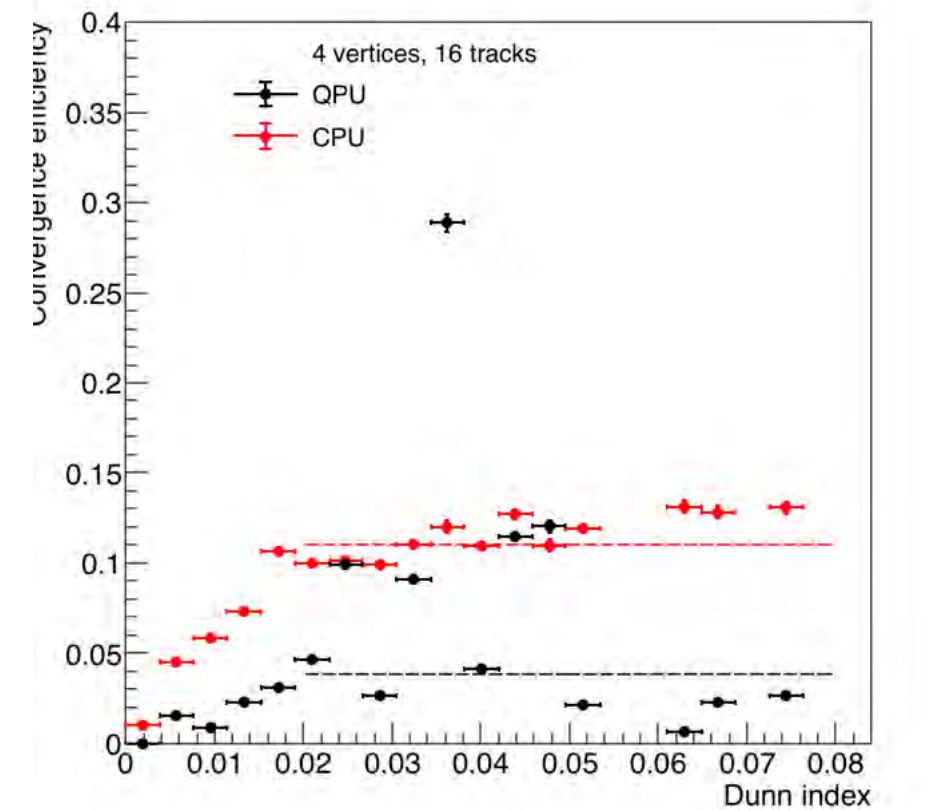
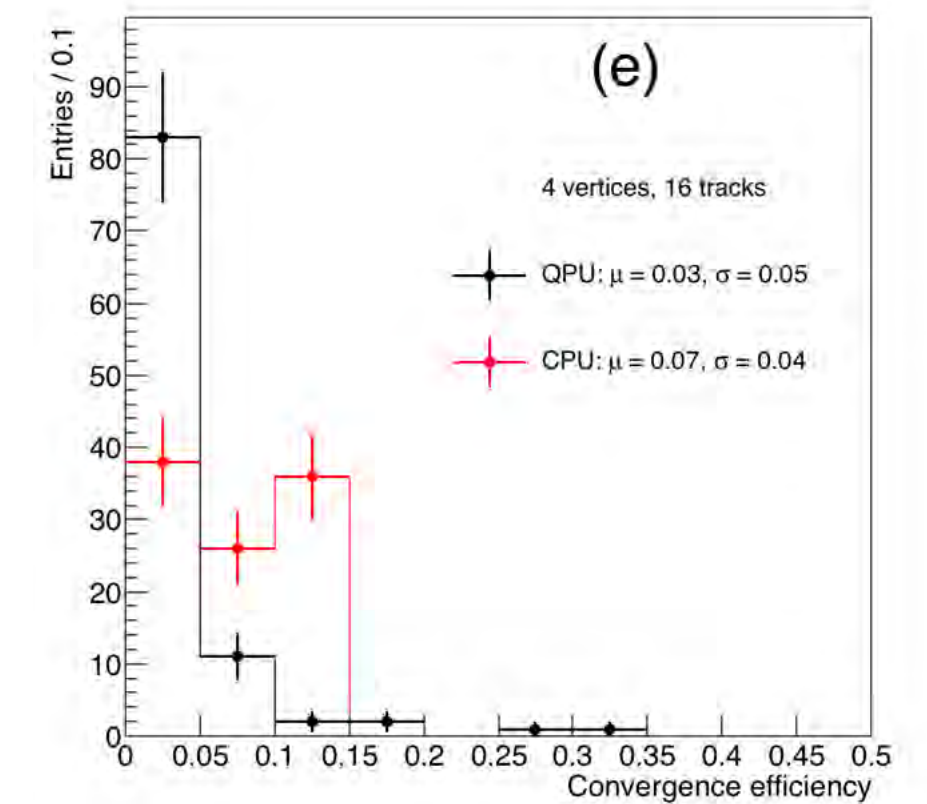
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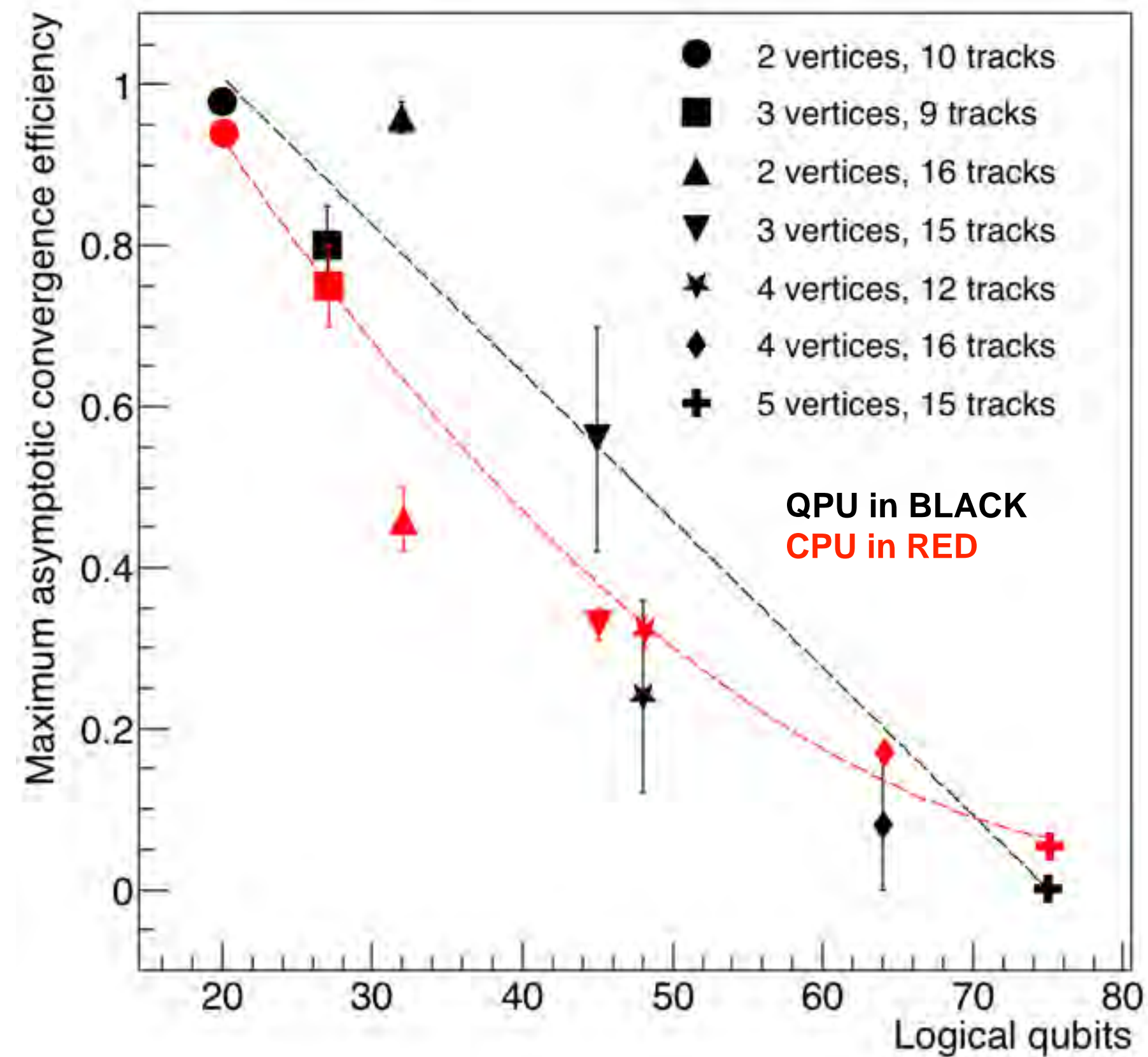
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Maximum convergence efficiencies

• QPU has advantage at low complexity. Why?

• Can any measure highlight the tunneling advantage?

# QPU vs CPU scaling with event complexity



- One measure of complexity: Number of logical qubits used = number of collisions × number of tracks
- Trend: **Asymptotic maximum of convergence efficiency** plotted against logical qubits
  - **Spread of maximum convergence efficiency** represented by uncertainty bars

**QPU performance comparable to a modern CPU**  
**QPU running may be further optimized**

# Conclusions and outlook

## Conclusions

- The D-Wave 2000Q\_2\_1 QPU can reconstruct p-p collision positions at hadron colliders in a limited capacity
  - ➔ The Tevatron had ~ 3 p-p collisions per event. Would have been possible with D-Wave
- QPU implementation comparable to Simulated Annealing on MacBook CPU for equal time
- QPU implementation to be optimized for LHC complexities: 50 to 200 p-p collision per event

## Outlook

Two research directions to improve QPU implementation:

- Improve convergence efficiency:
  - ➔ Understand how distortion functions like  $g(x; m)$  work
  - ➔ Use annealing offsets
  - ➔ Tune annealing time, re-thermalization delay
  - ➔ Try reverse annealing
  - ➔ Optimize chain lengths and weights
- Fit larger problems on QPU:
  - Customized embedding
  - Solve larger problems with hierarchical clustering

## Acknowledgements

Support and useful discussions with D-Wave: **Joel Gottlieb**, Mark Johnson, Alexander Condello  
 Support from the Department of Energy grant DE-SC0007884 and the Purdue Research Foundation

Track clustering with a quantum annealer for primary vertex reconstruction at hadron colliders

S. Das, A. J. Wildridge, S. B. Vaidya, A. Jung  
 Department of Physics and Astronomy, Purdue University

### Abstract

Clustering of charged particle tracks along the beam axis is the first step in reconstructing the positions of hadronic interactions, also known as primary vertices, at hadron collider experiments. We use a 2048 qubit D-Wave quantum annealer to perform track clustering in a limited capacity on artificial events where the positions of primary vertices and tracks are drawn from distributions measured by the Compact Muon Solenoid experiment at the Large Hadron Collider. The algorithm, which is not a classical-quantum hybrid but relies entirely on quantum annealing, is tested on a variety of event topologies from 2 primary vertices and 10 tracks to 5 primary vertices and 15 tracks. It is benchmarked against simulated annealing run on a modern CPU constrained to the same processor time per anneal as time in the physical annealer, and performance is found to be comparable. We chart three research directions to improve the performance of quantum annealers for this class of problems.

### 1. Introduction

Hadron colliders circulate counter-rotating beams of hadrons in closely packed bunches that cross at designated interaction points. These interaction points are instrumented with experiments that detect particles produced at hadron-hadron collisions when the bunches cross. Reconstructing the positions of these collisions within a bunch crossing, also known as primary vertices, from the trajectories of charged particles detected by the apparatuses is of paramount importance for physics analyses. The Large Hadron Collider (LHC) is a high luminosity collider that produces an average of 20 proton-proton (p-p) collisions at each bunch crossing, distributed in one dimension along the beam axis. At one of the LHC interaction points, the Compact Muon Solenoid experiment (CMS) reconstructs the paths of charged particles from p-p collisions as tracks detected by its silicon tracker [1]. Track reconstruction uncertainties obscure which tracks originated together at a primary vertex. Thus, primary vertex reconstruction begins with a one-dimensional clustering of tracks by their positions along the beam axis where they approach it most closely, also known as the tracks'  $z_0$ . In this paper, we demonstrate a method of performing this clustering on a D-Wave quantum annealer and report preliminary results benchmarked against simulated annealing on a classical computer.

The D-Wave 2000Q quantum computer, available from D-Wave Inc., performs computations through quantum annealing [2, 3, 4]. The quantum processing unit (QPU)

has 2048 RF-SQUID flux qubits implemented as superconducting niobium loops [5]. Each qubit has a programmable external magnetic field to bias it. The network of qubits is not fully connected and programmable couplings have been implemented between 6016 pairs of qubits. A computational problem is defined by setting the biases ( $h_i$ ) and couplings ( $J_{ij}$ ) such that the ground state of the qubits' Hamiltonian corresponds to the solution. We call this the "problem Hamiltonian" ( $H_p$ )

$$H_p = \sum_i h_i \sigma_z^i + \sum_i \sum_{j>i} J_{ij} \sigma_z^i \sigma_z^j, \quad (1)$$

where  $\sigma_z^i$  is a spin projection observable of the  $i^{\text{th}}$  qubit with eigenvalues +1 and -1. (This  $z$  direction is not related to the beam axis at CMS.) It may be trivially mapped to a bit observable  $q_i$  with eigenvalues 0 and 1 through the shift  $2q_i = \sigma_z^i + I$ , where  $I$  is the identity matrix. The problem Hamiltonian may then be expressed for quadratic unconstrained binary optimization (QUBO) as

$$H_p = \sum_i a_i q_i + \sum_i \sum_{j>i} b_{ij} q_i q_j, \quad (2)$$

notwithstanding energy offsets that are irrelevant for optimization. The D-Wave 2000Q programming model allows us to specify a problem in QUBO form by specifying  $a_i$  and  $b_{ij}$ .

At the beginning of a typical annealing cycle in the QPU, a driver Hamiltonian puts all qubits in a superposition of the computational basis states by introducing a global energy bias in the transverse  $x$ -direction. Annealing proceeds by lowering this driver Hamiltonian while si-

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**S. Das, A. J. Wildridge, S. B. Vaidya, A. W. Jung,**  
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<https://arxiv.org/abs/1903.08879>