QUnfold: Quantum Annealing for Distributions Unfolding in High-Energy Physics

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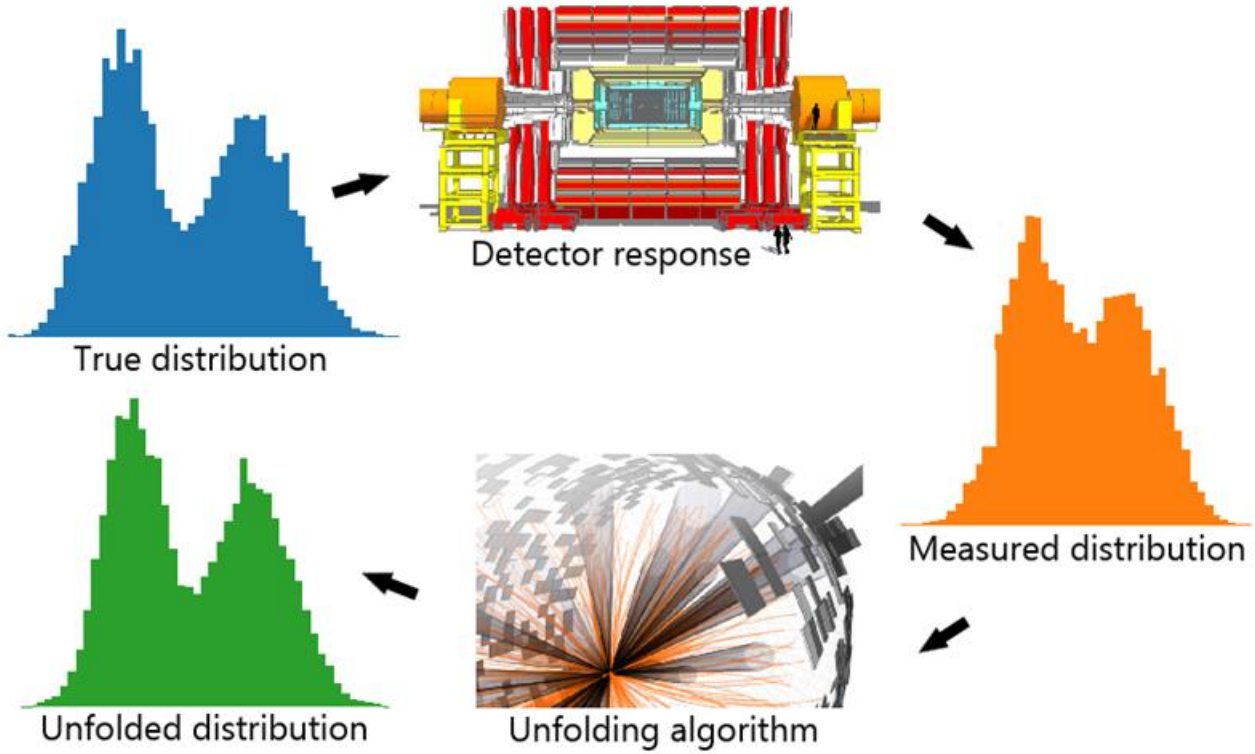
What is unfolding?

- In **High-Energy Physics** (HEP) experiments each measurement apparatus has a unique signature in terms of *detection efficiency*, *resolution,* and *geometric acceptance*
- The overall effect is that the distribution of some measured observable in a given physical process is *biased* and *distorted*
- **Unfolding** is the mathematical technique to correct for this distortion and recover the original distribution



$$\vec{\mu} = R\vec{x}$$

"reco" distribution

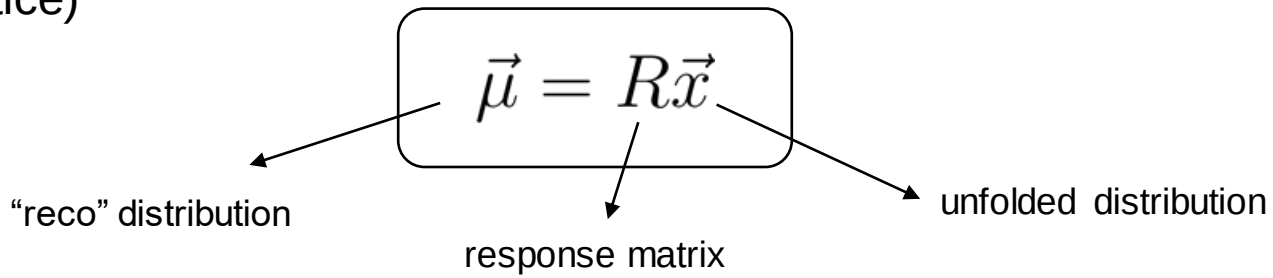
response matrix

unfolded distribution

Unfolding techniques

Classical unfolding methods in HEP:

- Standard matrix inversion (never used in practice)
- Bin-by-bin unfolding (never used in practice)
- Likelihood-based unfolding (SVD)
- Iterative Bayesian unfolding (IBU)



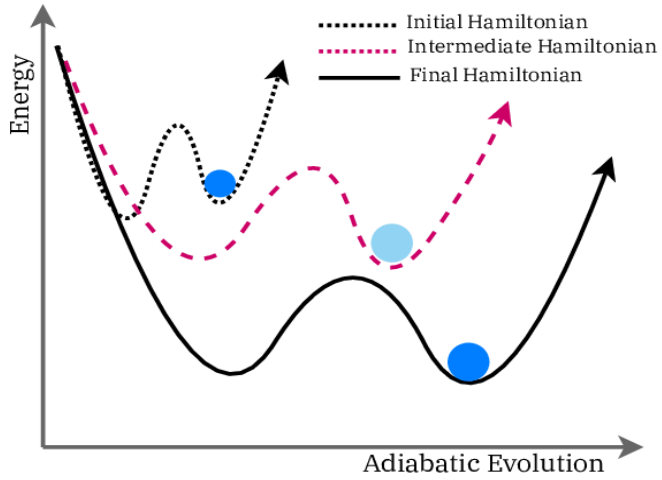
"Quantum" unfolding methods:

- First proof-of-concept by *R. Di Sipio* et al in 2019: the model worked only on really small-sized problems (very few bins and entries) using the D-Wave 2000Q quantum annealer machine
- Our open-source experimental proposal is *QUnfold*

What is quantum annealing?

- The quantum-mechanical system is prepared in the known ground-state of an **initial Hamiltonian** H_{init}
- The target solution is encoded in the ground-state of a **final Hamiltonian** H_{fin}, written as the energy/cost function of a Quadratic Unconstrained Binary Optimization problem (QUBO problem)
- The system evolution is controlled by the following time-dependent Hamiltonian:

$$H(t) = A(t)H_{init} + B(t)H_{fin} \qquad A(t)\downarrow \quad B(t)\uparrow$$



- **Quantum Adiabatic theorem**:
  «if the evolution is slow enough, the quantum-mechanical system stays close the ground-state of the istantaneous Hamiltonian»

QUBO problem $\Rightarrow$ $H(\vec{x}) = \sum_i a_i x_i + \sum_{i,j} b_{ij} x_i x_j \qquad x_i \in \{0,1\} \quad a_i, b_{ij} \in \mathbb{R}$

D-Wave quantum annealer

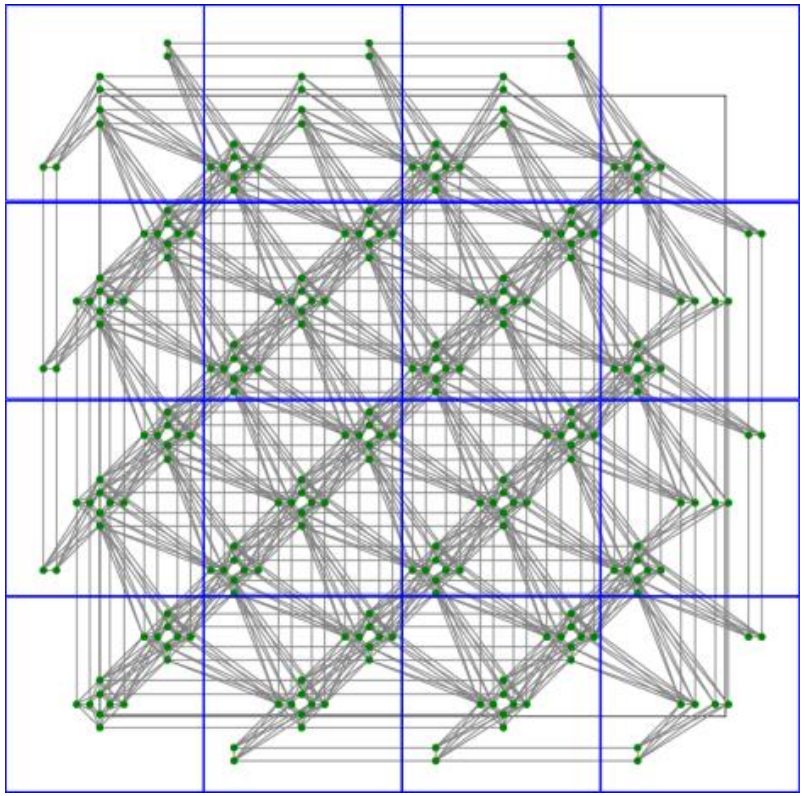
- The D-Wave company is the only commerical quantum annealing machines provider so far (1 min/month QPU access time for free)

- The D-Wave QPU is a **lattice of interconnected superconducting qubits** operating at around 15 mK and with a fixed limited topology

D-Wave Advantage is currently their best quantum annealer:

- 5000+ qubits
- 35000+ couplers
- *Pegasus* topology

Pegasus topology in D-Wave Advantage



D-Wave
The Quantum Computing Company™

QUnfold - Mathematical formulation

Log-likelihood maximization unfolding: $\max_{\vec{x}} \left(\log \mathcal{L}(\vec{\mu}|\vec{d}) + \lambda \mathcal{S}(\vec{\mu}) \right)$

measured distribution

regularization term

$\vec{\mu} = R\vec{x}$

"reco" distribution

detector response matrix

unfolded distribution

Quadratic model minimization: $\min_{\vec{x}} \left(||R\vec{x} - \vec{d}||^2 + \lambda ||G\vec{x}||^2 \right)$

Tikhonov regularization: discrete 2nd order derivative (Laplacian operator G)

- Classical optimization problem, **nothing quantum yet**!
- $\vec{x}$ is the vector of integer numbers representing the unfolded histogram

QUnfold - Mathematical formulation

$$\vec{a} = -2R^T\vec{d}$$
$$B = R^TR + \lambda G^TG$$

$$\Longrightarrow \quad H(\vec{x}) = \sum_i a_i x_i + \sum_{i,j} b_{ij} x_i x_j = \vec{a}\cdot\vec{x} + \vec{x}^T B\vec{x}$$

To get the QUBO model from this integer-variables quadratic problem, a “binarization” process based on the **logarithmic encoding** of the variables is needed

Their total number, which represents also the number of required logical qubits, scales as:

$$N_{\text{qubits}} \propto \sum_{i=1}^{N_{\text{bins}}} \log_2(N^i_{\text{entries}})$$

N_{bins} is the number of bins of the histogram
N_{entries} is the vector of the number of entries in each bin of the histogram

Example:
- Gaussian distribution
- 20 histogram bins
- 5M histogram entries

→ *N* qubits ≈ 350

QUnfold - Software package

- Implemented using *NumPy* and *D-Wave Ocean SDK* but fully compatible with ROOT
- Designed to address real-scale HEP applications
- Very simple and intuitive **Python** interface
- Public repository and documentation on GitHub
- Available on PyPI and easy to install via *pip*:

```
pip install Qunfold
```

Solver methods:

- Simulated annealing sampler (CPU only)
- Hybrid sampler (CPU + QPU)
- Quantum annealing sampler (QPU only)

The package is still work in progress!

```python
# Import QUnfold base class and plotter
from QUnfold import QUnfoldQUBO
from QUnfold import QUnfoldPlotter

# Read numpy data from a file or sample them
truth = ... # truth distribution
measured = ... # measured distribution
response = ... # response matrix
binning = ... # binning of the distributions

# Run unfolding
unfolder = QUnfoldQUBO(response, measured, lam=0.1)
unfolder.initialize_qubo_model()
unfolded_SA, error_SA = unfolder.solve_simulated_annealing(num_reads=10)

# Plot results
plotter = QUnfoldPlotter(
    response=response,
    measured=measured,
    truth=truth,
    unfolded=unfolded_SA,
    error=error_SA,
    binning=binning,
)
plotter.plotResponse()
plotter.plot()
```

Free regularization parameter

QUnfold - Tests on simulated data

Dataset

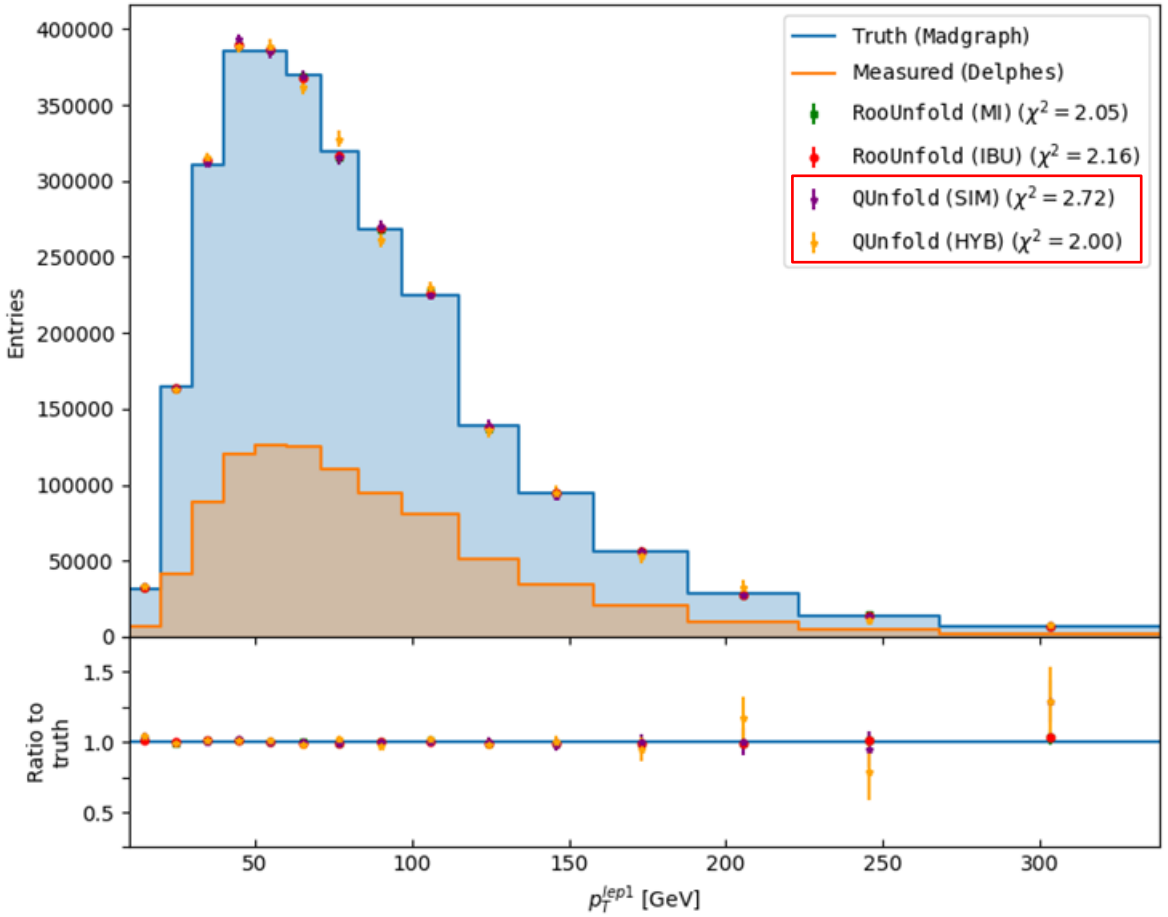
- $t\bar{t}$ **process** in the *dileptonic channel* (2 leptons and at least 2 *b*-jets required in the final state)
- ≈ 2.5M **truth-level** events generated using the *MadGraph* generator (*truth* distribution)
- **Detector-level** data generated using the *Delphes* simulator (*measured* distribution)

Technique

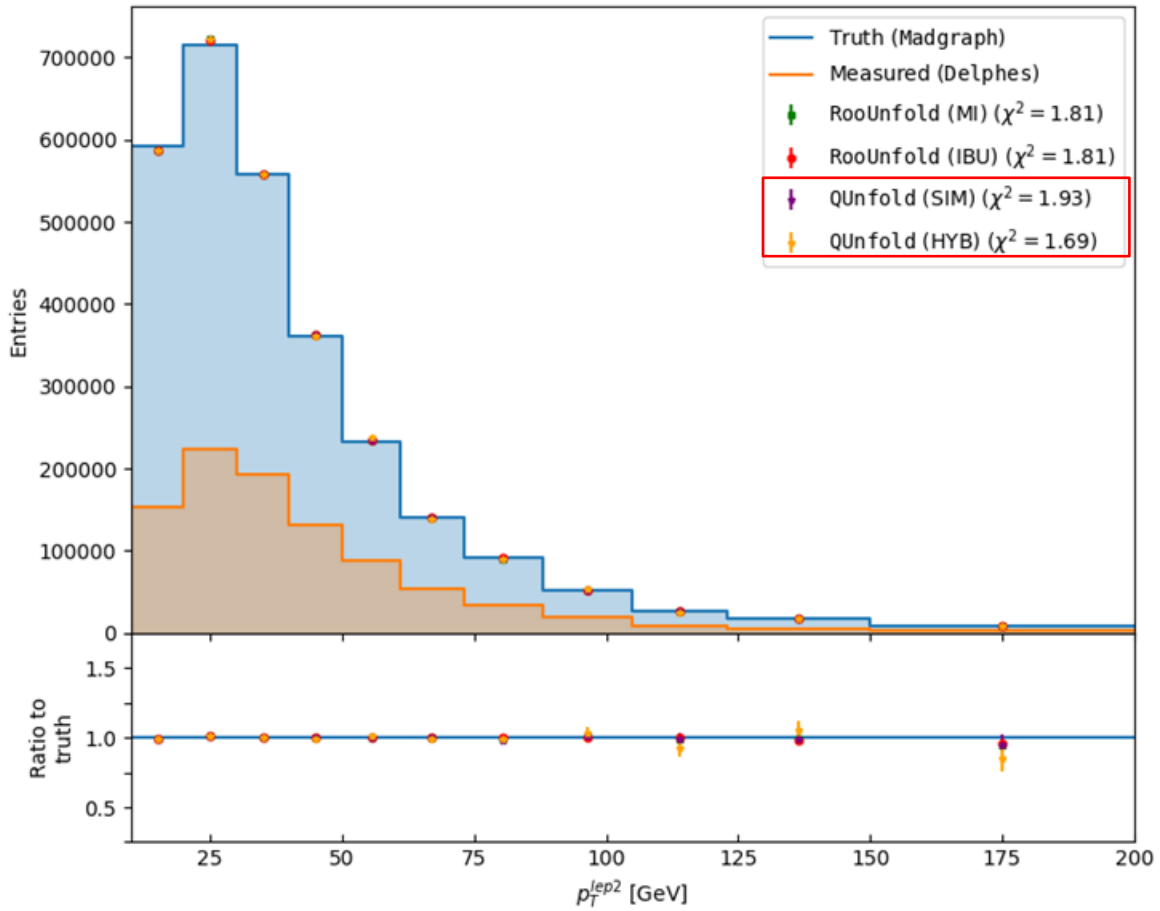
- Simulated annealing and hybrid solvers are used (quantum annealing solver is work in progress)
- Results are compared to the classical HEP unfolding methods *MI* and *IBU* (RooUnfold framework)
- **Toy Monte Carlo experiments** are run to compute the covariance matrix for evaluating the quality of the result (X^2 test) and estimating the statistical errors associated to the unfolding method

QUnfold - Preliminary results

Leading lepton p_T

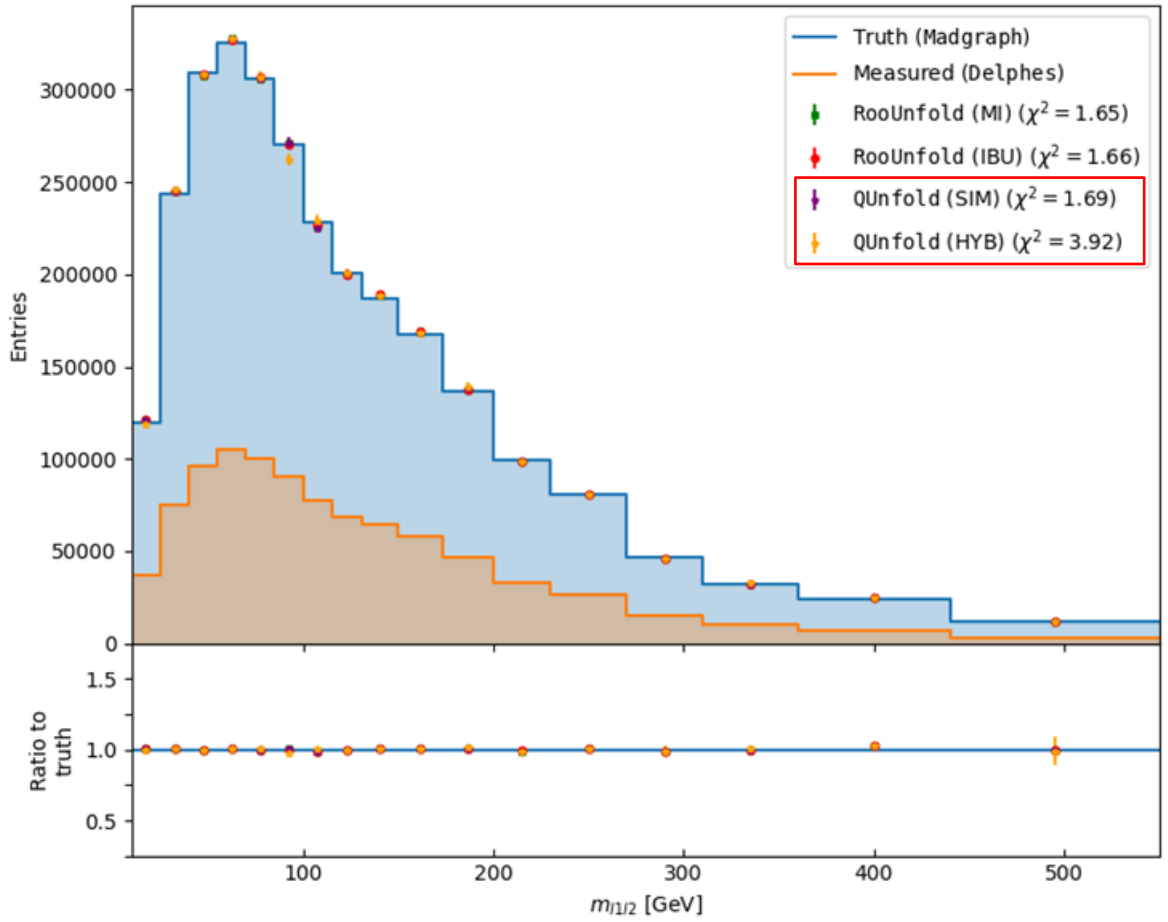


Subleading lepton p_T

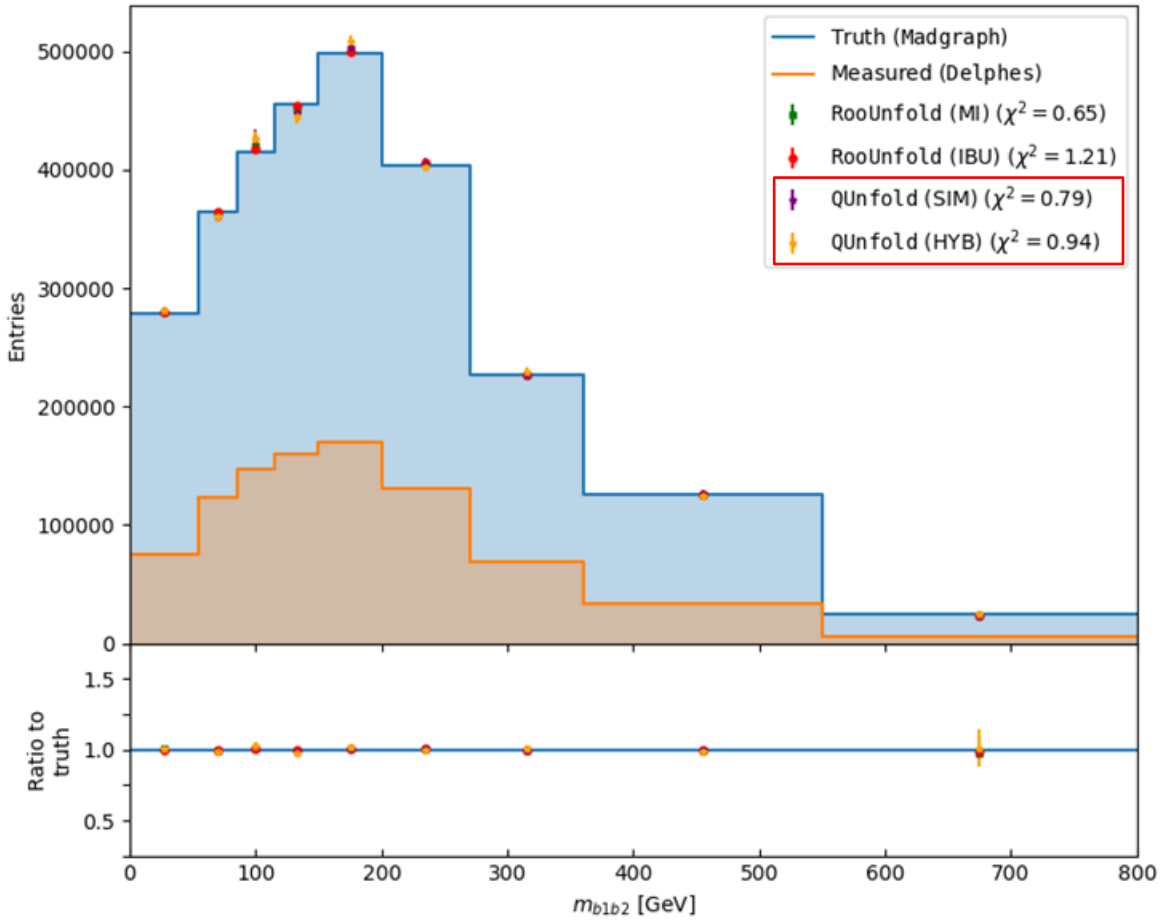


QUnfold - Preliminary results

Leptons invariant mass



b-jets invariant mass



Conclusion

- New unfolding approach based on the **QUBO formulation** of the problem and ***quantum annealing***
- Model implemented and tested in the **QUnfold** Python package, very easy to install and start using

Future steps

- Further optimize the algorithm (integer model *binarization*, QUBO matrix *pre-conditioning*, etc.)
- Perform more experiments on real quantum hardware (D-Wave resources by CINECA)
- Develop PyXSec: a new framework to measure differential cross-sections of HEP processes

Thanks for the attention!

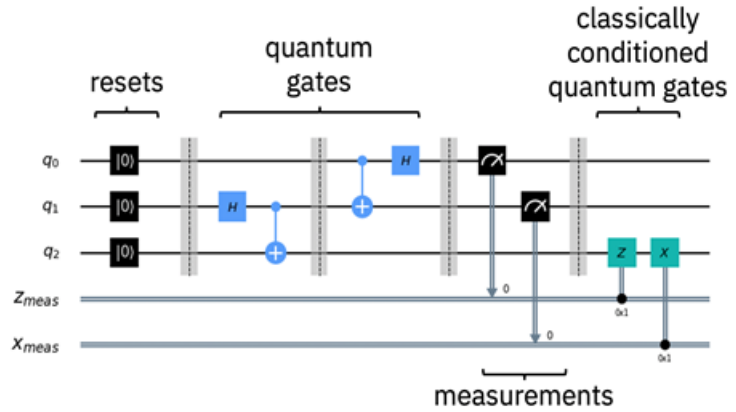
https://github.com/JustWhit3/QUnfold

Backup

Quantum computing

Key concepts:

- Quantum computing: science based on *information theory* and *quantum mechanics*
- Unit of measurement of quantum information is the qubit (mix of 0 and 1 states)
- Quantum algorithms: algorithms that exploit properties from quantum mechanics
- Quantum computing is performed by creating quantum circuits
- Quantum algorithms need to operate through *quantum computers*

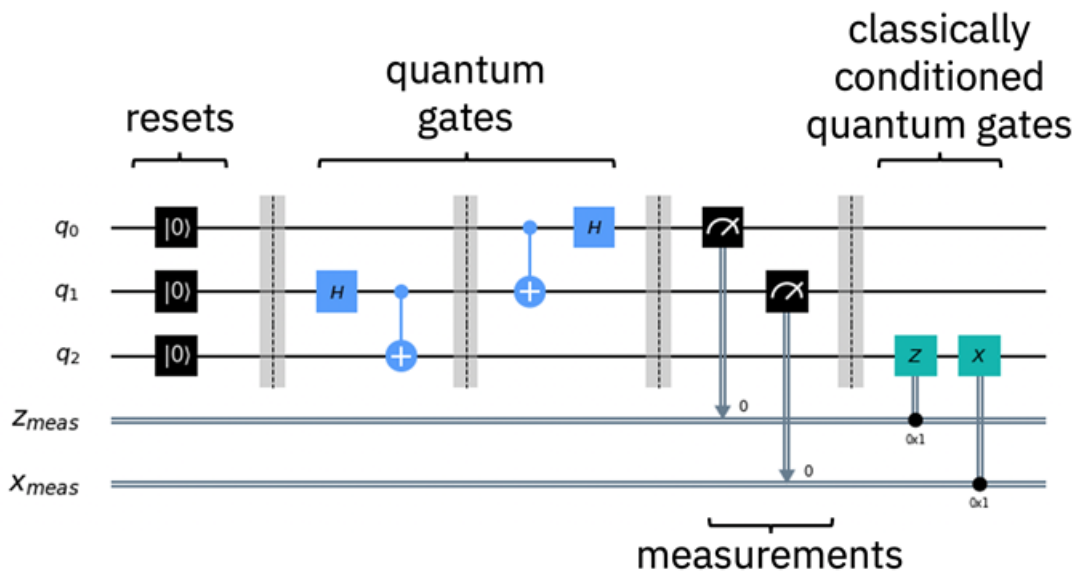


Quantum computing is based on **3 fundamental quantum concepts**:

- *Superposition principle*
- *Quantum entanglement*
- *Tunneling effect*

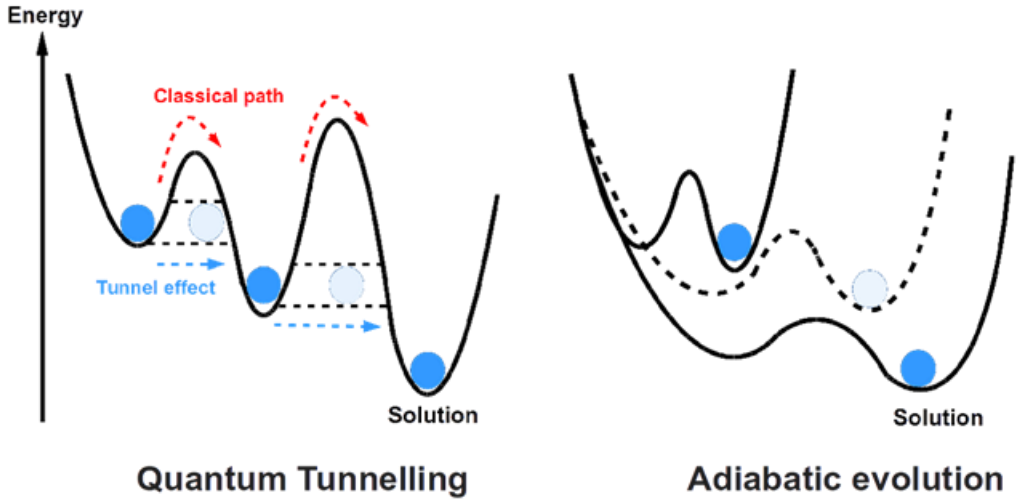
Quantum computing and quantum annealing

Gate-based quantum computing
(eg: IBM, Google)



Universal

Quantum-annealing-based quantum computing
(eg: D-Wave)



Only quantum annealing can be performed

Measurement of differential cross-sections

Current standard method used in ATLAS and CMS:

- Particle/parton-level phase space
- Iterative Bayesian unfolding (IBU)
- In ATLAS TTbarUnfold is used (written in C++ by Marino)

$$\frac{d\sigma^{\text{fid}}}{dX^i} \equiv \frac{1}{\mathcal{L} \cdot \Delta X^i} \cdot \frac{1}{\epsilon^i} \cdot \sum_j M^{-1} \cdot f_{\text{acc}}^j \cdot \left(N_{\text{obs}}^j - N_{\text{bkg}}^j\right)$$

$$\frac{d\sigma^{\text{norm}}}{dX^i} = \frac{1}{\sigma^{\text{fid}}} \cdot \frac{d\sigma^{\text{fid}}}{dX^i}$$

Our idea:

- Working on a new general framework called PyXSec (open-source on GitHub), based on TTbarUnfold but written in Python
- Add full support to cross-sections measurements by using both **RooUnfold** classical methods and **QUnfold** quantum algorithms

X^2 and errors computation

Covariance matrices and **errors** are computed through *MC pseudo-experiments*:

- A random *Poissonian smearing* is added to the measured distribution
- Unfolding is performed
- Procedure is repeated for *N* iterations (**toys**)
- Covariance matrix is computed considering the ensemble of the unfolding solution at each iteration:

$$c_{ij} = < (x_i - < x_i >)(x_j - < x_j >) >$$

- Errors are computed as the square-root of the diagonal of the covariance matrix

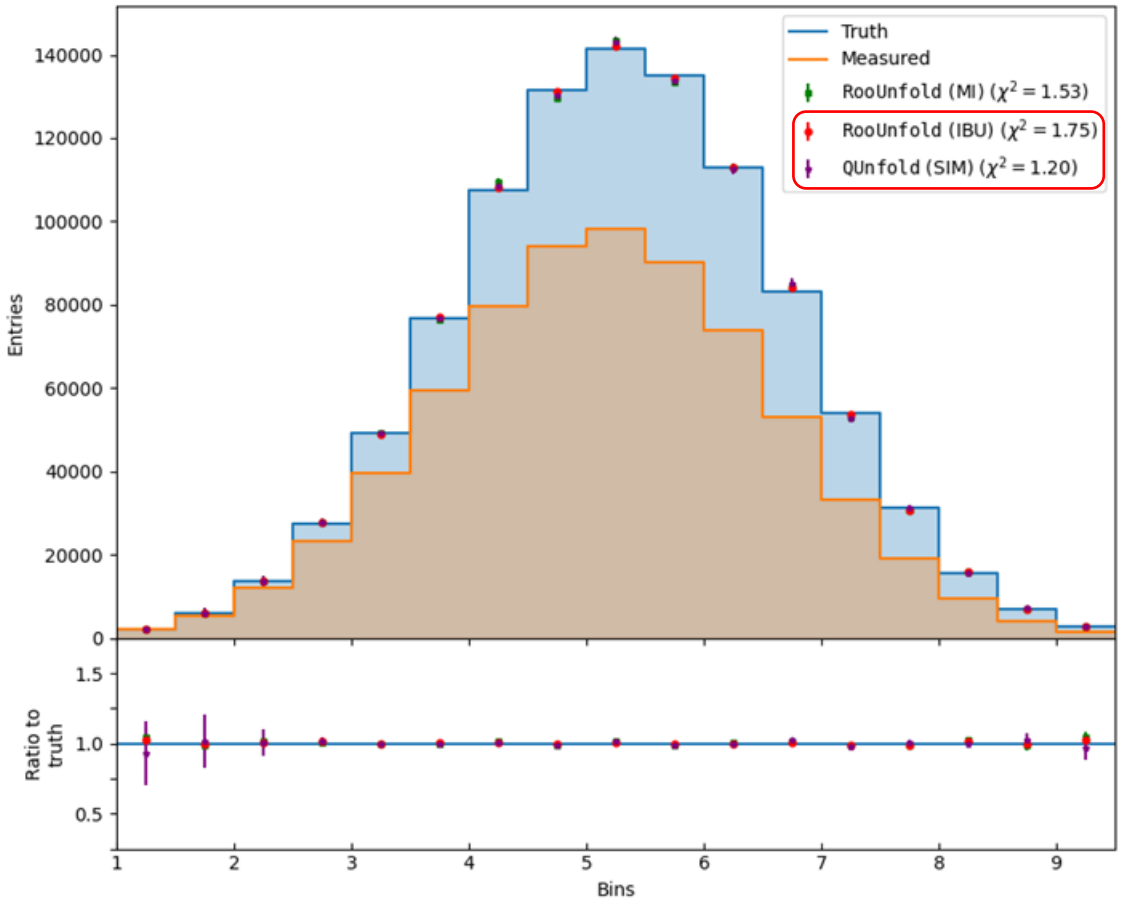
$\mathbf{X^2}$ are computed with:

$$X^2 = V^T \times Cov^{-1} \times V$$

Where V is the vector of *residuals*, defined as the difference between measurement and prediction

Preliminary results with Numpy

Normal



Exponential