

STATISTICAL METHODS FOR CROSS-SECTION MEASUREMENTS: PAST, PRESENT AND FUTURE

Lukas Koch

ECT* Workshop “Measuring neutrino

interactions for next-
generation oscillation
experiments” 2024



JOHANNES GUTENBERG
UNIVERSITÄT MAINZ

JG|U

DISCLAIMER: OPINIONS!

- Necessarily more familiar with some methods compared to others
 - Biased sample of methods previously/currently in use
 - **Very** biased sample of potential future developments
- If anything seems fishy, probably my fault
- Many subtleties at every step
 - I do not have the time to get into
- Open to Bayesian methods, but biased towards Frequentism
- Most probable answer in statistics: “It depends”

SOME NOTATION

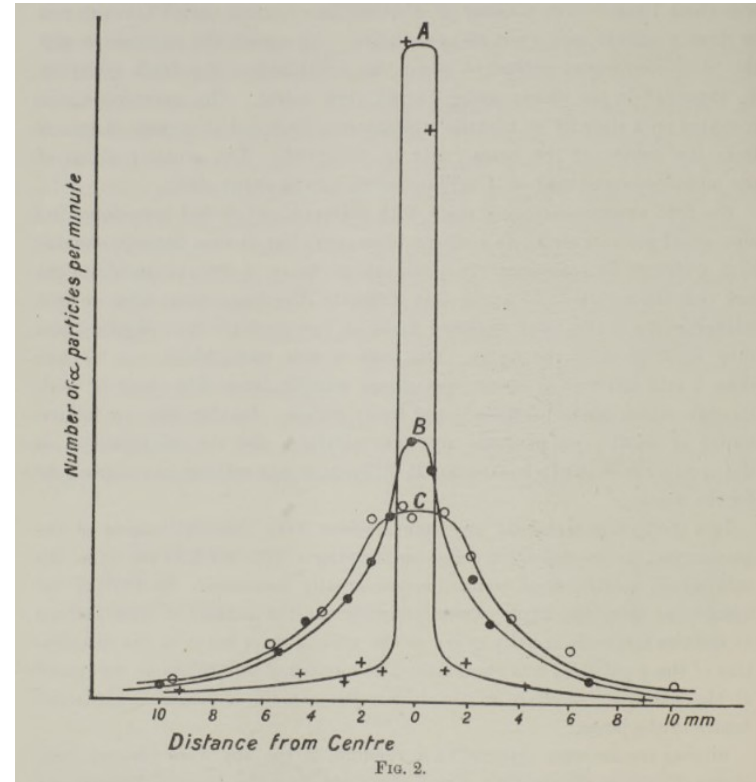
- $v_i = \sum_j R_{ij} \mu_j$
 - Expected number of observed events v_i in reco bin i
 - Expected number of true events μ_j in truth bin j
 - Response matrix R is $N \times M$ matrix
- Observed events:
 $n_i \sim \text{Poisson}(v_i)$
- True events:
 $m_j \sim \text{Poisson}(\mu_j)$
- Binned in multiple variables
- Not necessarily same physical meaning
 - $\text{track_length_reco} = R * \text{momentum_true}$
- Purely mathematical approach:
 $R = P(\text{event in reco } i \mid \text{event in truth } j)$
 $= S * \text{eff}$
- Background handling approaches
 - Subtract from observed events:
 $n_i = o_i - b_i$
 - “Breaks” Poisson statistics
 - Add to expectation
 $v_i = \zeta_i + \beta_i$

EVENT RATES VS CROSS SECTIONS

- $\mu_j = \sum_k T (d\sigma/dy)_{jk} \Phi_k \Delta y_j = T (d\sigma/dy)_{j,\Phi\text{-avg}} \Phi \Delta y_j$
 - For “thin” targets
 - For a neutrino, “thin” can mean a lightyear of lead
 - Assuming cross section is sufficiently constant over bin!
- Conceptual steps:
 - Measure $n_i \rightarrow$ Use it as proxy for ν_i
 - Unfold and efficiency correct to μ_j
 - Convert event rates to cross sections
- Uncertainties break neat factorisation
 - E.g. detector smearing depends on neutrino flux uncertainty?
- Details vary a lot: “It depends”

JUST LOOK AT RECO

- Implicitly compare n_i with μ_j
 - Pretend y_{reco} and y_{truth} are the same
- Ancient past: Don't even put error bars
 - Not as unreasonable as it sounds
 - n vs. ν
- Slight improvement: bin-by-bin efficiency correction: n_i / eff_i
 - Only does what you expect if R is diagonal → No smearing



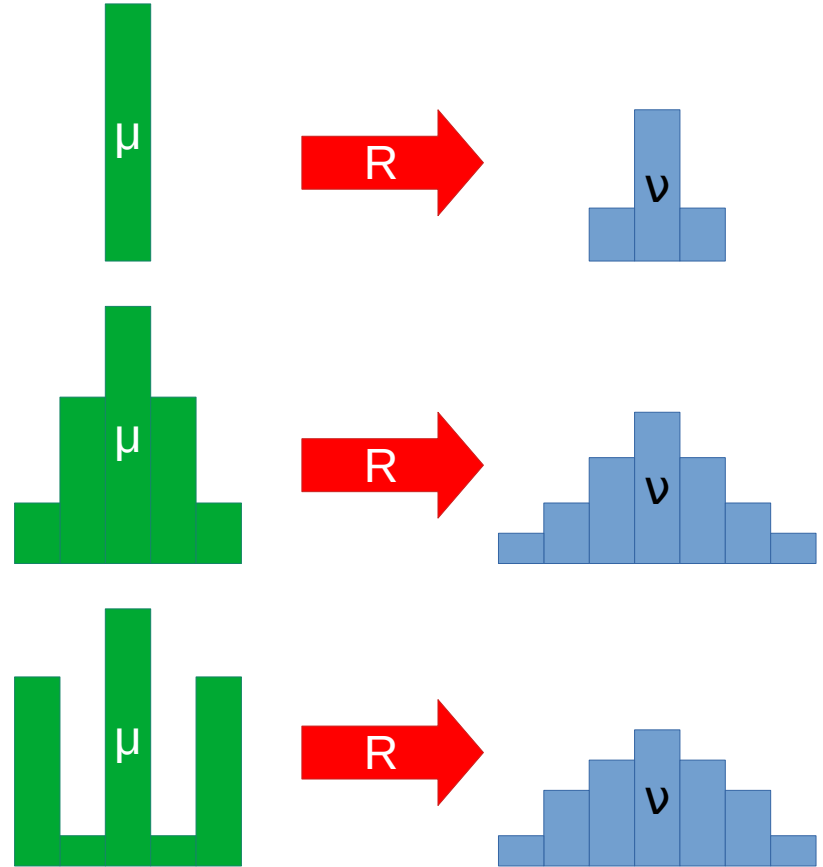
H. Geiger, On the scattering of the α -particles by matter, <https://doi.org/10.1098/rspa.1908.0067>

NAIVE APPROACH: JUST INVERT R

- Usually we have smearing
- $\mathbf{v} = \mathbf{R}\boldsymbol{\mu}$ so why not just calculate $\boldsymbol{\mu} = \mathbf{R}^{-1} \mathbf{v} \approx \mathbf{R}^{-1} \mathbf{n}$
- Possible when $N = M$
 - Choose suitable left-inverse when $N > M$
- Solves least squares problem:
 - Minimize $|\mathbf{v} - \mathbf{n}|^2 = |\mathbf{R}\boldsymbol{\mu} - \mathbf{n}|^2$
 - $\hat{\boldsymbol{\mu}} = (\mathbf{R}^T \mathbf{R})^{-1} \mathbf{R}^T \mathbf{n} = \mathbf{R}^{-1} \mathbf{n}$
 - Equivalent to maximum likelihood solution when uncertainties Gaussian with known variances
- Can lead to large variance and strong anticorrelations in result

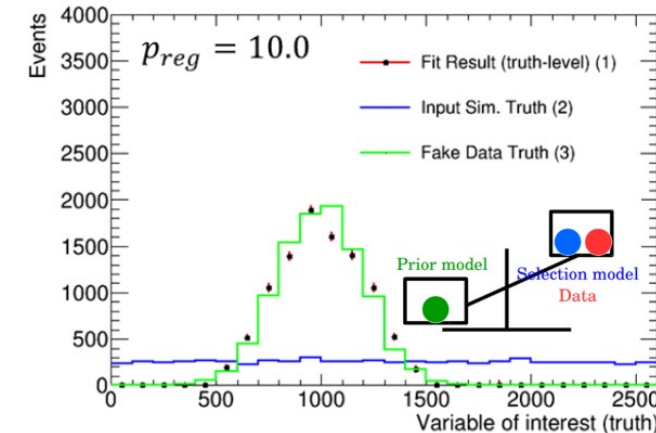
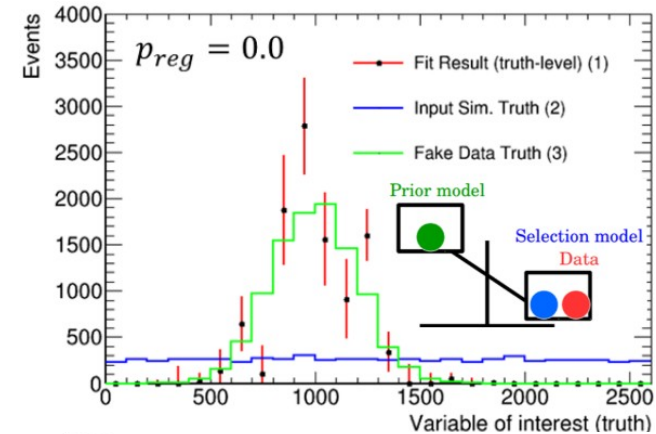
THE ILL POSED PROBLEM

- Strong correlations stem from fact that very different μ can lead to very similar ν
- Small fluctuations in n lead to large swings in “best guess” at μ
- Many different solutions are virtually indistinguishable
 - Pick a nicer looking one!
- Impose a slight preference for “nice looking” results
 - Can be interpreted as Bayesian prior or Frequentist external constraint



RIDGE REGRESSION / TIKHONOV REGULARISATION

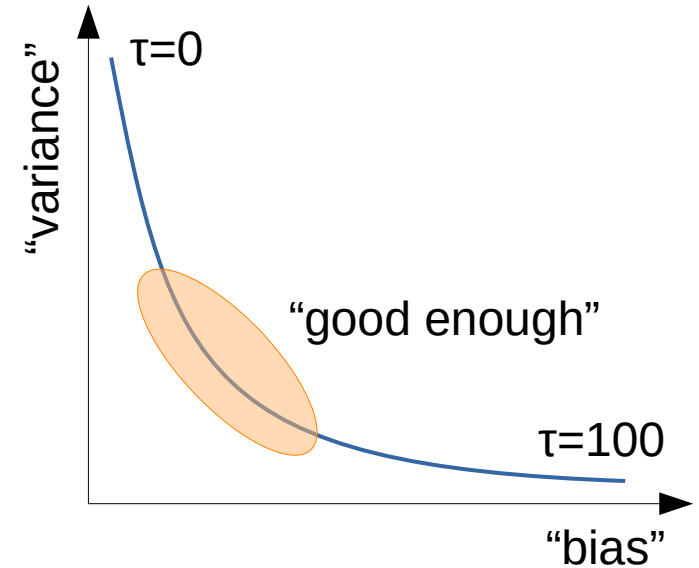
- Modify optimisation problem
 - Add a **penalty term** for “bad looking” solutions
 - Minimize $|\mathbf{R}\boldsymbol{\mu} - \mathbf{n}|^2 + |\mathbf{C}\boldsymbol{\mu}|^2$
 - $|\mathbf{C}\boldsymbol{\mu}|^2 = \boldsymbol{\mu}^T \mathbf{C}^T \mathbf{C} \boldsymbol{\mu} = \boldsymbol{\mu}^T \mathbf{Q} \boldsymbol{\mu}$
- Tikhonov matrix C, or penalty matrix Q
 - Notations vary
 - Choice of C/Q determines what is penalised and how strongly, e.g.
 - $\mathbf{Q} = \tau \mathbf{I} \rightarrow L_2$ norm of $\boldsymbol{\mu}$
 - $\boldsymbol{\mu}^T \mathbf{Q} \boldsymbol{\mu} = \tau \sum (\mu_j - \mu_{(j+1)})^2$
→ Squared differences of neighbouring bins
- New solution
 - $\hat{\boldsymbol{\mu}} = (\mathbf{R}^T \mathbf{R} + \mathbf{Q})^{-1} \mathbf{R}^T \mathbf{n}$
 - Adding Q makes $\mathbf{R}^T \mathbf{R}$ “less problematic” to invert



Borrowed from S. Dolan

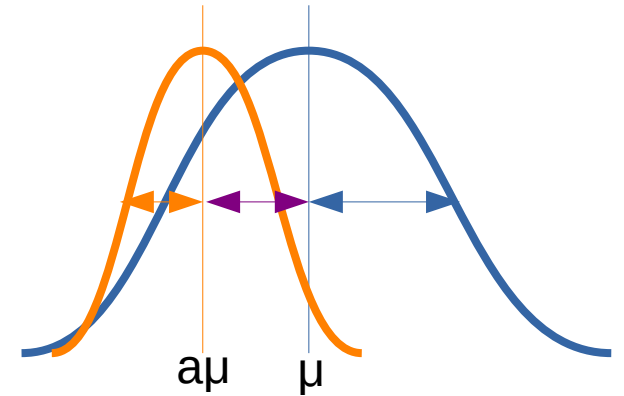
HOW STRONGLY TO REGULARISE

- Regularisation can be seen as prior/external constraint
 - Should be well defined
- Mostly it is introduced ad-hoc
 - Might know what we dislike, but not how much
 - Regularisation strength τ not known a priori
- Regularisation introduces bias
 - Also messes with coverage properties
- Usually some heuristic method to “balance” bias and variance of result
 - e.g. L-curve method
- Can define an objective function and optimize with respect to it
 - What should be optimized can be subjective



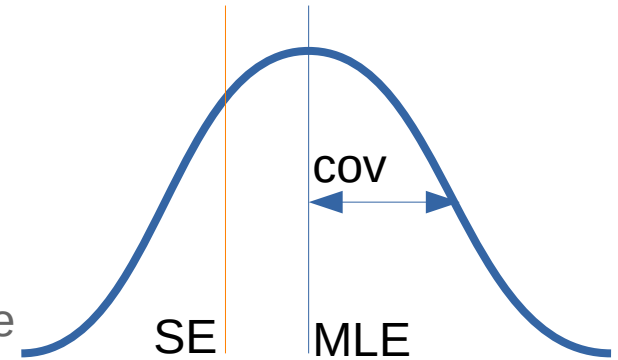
STATISTICAL SHRINKAGE

- Why is it reasonable to penalise large $|\mu|^2$?
- E.g. want to estimate mean value of normal distribution
- Single sample x from $N(\mu, \sigma)$
 - Maximum likelihood estimator (MLE): $\hat{\mu} = x$
 - $E[(x-\mu)^2] = \sigma^2$
- Multiply x by shrinkage factor a
 - Shrinkage estimator (SE): $\hat{\mu} = ax$
 - $E[(ax-\mu)^2] = (a-1)^2\mu^2 + a^2\sigma^2$
 - Minimal at $a = \mu^2 / (\sigma^2 + \mu^2) < 1$
- SE reduces expected squared deviation from true mean compared to MLE!
 - At cost of biasing point estimate towards 0
- Choosing a point estimator does not affect the likelihood function



POINT ESTIMATE VS LIKELIHOOD FUNCTION

- But all information of experiment is (should be) inside **likelihood function**
 - Often approximated as **MLE** and **covariance matrix**
 - It is what it is, even if we do not like how it looks
- Understand regularisation as shrinkage
 - Picking a “reasonable” **point estimate**
 - **Not** to regularise the **likelihood function**
- Regularised covariance just a visualisation tool?
 - Pick a subset of the allowed region around the point estimate
 - Less correlations, less confusing plots
- Need both for full picture
 - Unregularised data release for “undiluted” likelihood function
 - Regularised result as “better” point estimate
 - Consensus for long time that it would be good to publish likelihood functions
 - Used both in Bayesian and Frequentist analyses



WIENER SVD

- Singular Value Decomposition (SVD) can be used to get left inverse of R and solve the least squares problem
- Apply Wiener filter which maximises signal to noise ratio
 - Assuming a given signal shape
 - Inspired by signal processing
 - This is the regularisation
- No tunable regularisation strength
 - Already “optimized” for the signal to noise ratio

RELATION TO UNREGULARISED RESULT

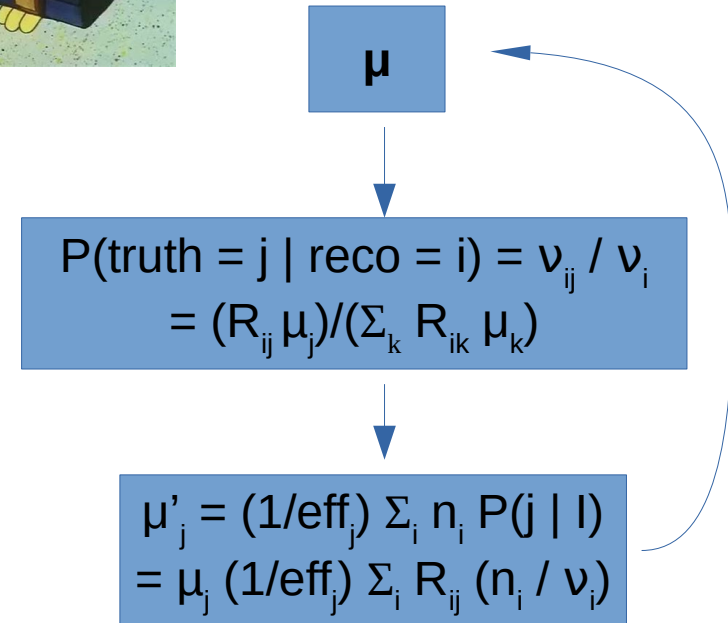
- Wiener SVD yields “additional smearing matrix” A
- It relates regularised result to unregularised one
 - $\boldsymbol{\mu}' = A \boldsymbol{\mu}$
 - Does this remind you of the shrinkage estimator?
 - $V' = AVA^T$
- No need to provide two separate results!
 - Just publish A together with either $(\boldsymbol{\mu}, V)$ or $(\boldsymbol{\mu}', V')$
- Better call A “regularisation matrix”?
 - Does not conserve event numbers and can have negative elements

ITERATIVE UNFOLDING / D'AGOSTINI METHOD



- Also known as Bayesian unfolding
 - Should we be calling it that?
 - It is Bayesian update of priors for 1 iteration
 - It approaches matrix inversion result for inf iterations (as long as all $\hat{\mu}$ are positive)
 - “Squeezing the data multiple times” for everything in between?
- # of iterations determines regularisation!
 - Low # → “remembers” first prior → strong regularisation
 - (# → inf) → “forgets” first prior → no regularisation
 - Assuming no smoothing in between iterations

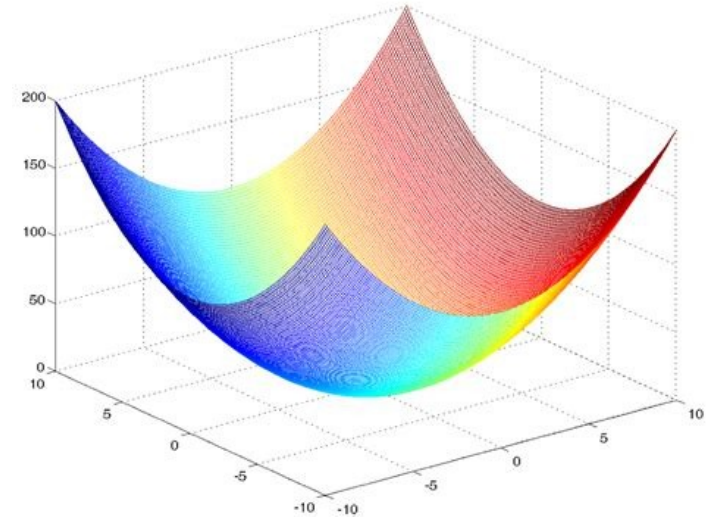
Simplified:



LIKELIHOOD FITTING

E.g. <https://arxiv.org/abs/2303.14228>

- Explicitly treat problem as parameter fit
 - Poisson likelihood in reco bins
 - Parameters of interest θ that scale cross section in truth bins
 - Systematic nuisance parameters φ
 - Constrained by “priors” = external constraints
 - “Just” need a function $-2 \log L(\theta, \varphi | \mathbf{n})$ and a minimizer
 - Get MLE & parabolic approximation (covariance)
- Add regularisation / penalty terms explicitly



FREQUENTIST FIT, BAYESIAN PROPAGATION?

- Result of fit contains many nuisance parameters
- Correlated uncertainties need to be propagated to XSECs
- Ideal Frequentist approach
 - For each M-dimensional XSEC, maximise likelihood over parameters
 - Profile likelihood
 - Not trivial
- Pragmatic approach
 - Throw parameters according to MLE & covariance
 - Calculate XSEC for each throw
 - Usually calculate central value and covariance from sample
 - Could also publish throws in case of non-Gaussian results

ADD REGULARISATION AFTER THE FACT?

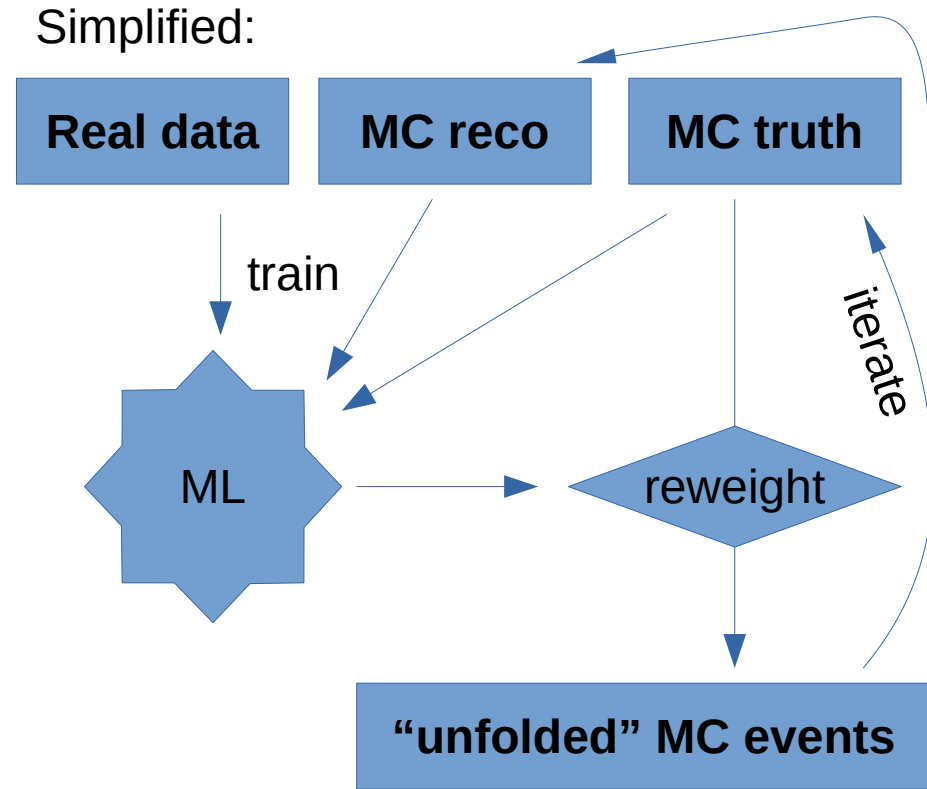
<https://doi.org/10.1088/1748-0221/17/10/P10021>

- Take inspiration from Wiener SVD
 - Apply regularisation as a matrix multiplication to the unregularised result
- Given any likelihood described as MLE & covariance, adding a Thikonov penalty term leads to a new result
- Can be applied to any unregularised result → post hoc
 - As long as regularised result is close to unregularised one
 - Parabola approximation of log likelihood stays valid

$$\begin{aligned} -2 \ln(L(\theta)) &\approx (\theta - \hat{\theta})^T V^{-1} (\theta - \hat{\theta}) + const. \\ P(\theta) &= \theta^T Q \theta \\ -2 \ln(L'(\theta)) &= -2 \ln(L(\theta)) + P(\theta) \\ &\approx (\theta - \hat{\theta}')^T V'^{-1} (\theta - \hat{\theta}') + const. \\ \hat{\theta}' &= A \hat{\theta} \\ V' &= A V A^T \\ A &= (V^{-1} + Q)^{-1} V^{-1} \end{aligned}$$

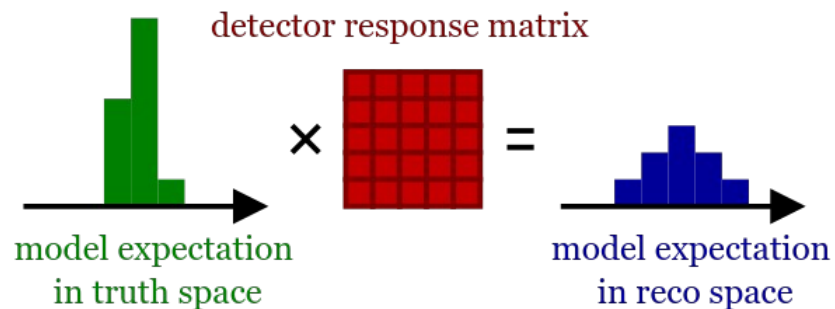
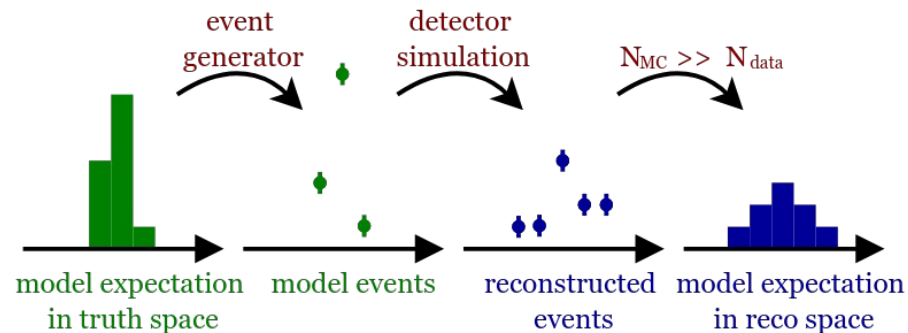
OMNIFOLD (AS I UNDERSTAND IT)

- Use Machine Learning (ML) techniques to create MC reweighter to match MC to measured reco data
 - Based on un-binned event properties
- Re-weighted MC is the “unfolded” result!
 - Can be binned in any way desired to report a XSEC
- Cutting edge research
 - Just about ready for production use?
 - We will hear more this week!



BACK TO THE ROOTS

- Possible to do science without unfolding
- Compare models with data in reco space
 - But consider detector effects: Forward folding
 - Allows full statistical analysis
 - The data is exactly what we saw: n is a perfectly known fixed number
 - Test whether models are compatible, i.e. the predicted v
- How to facilitate use of data by external consumers?
 - Not experts on the detector response
 - No access to (often complicated) simulation frameworks
 - Data needs low entry barrier to be used by many people

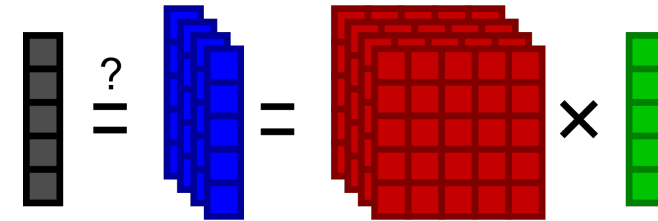


SOFTWARE AVAILABLE

- Delphes
 - <https://cp3.irmp.ucl.ac.be/projects/delphes>
 - Developed for collider experiments
- Rivet
 - <https://rivet.hepforge.org/>
 - Developed for collider experiments
- ReMU – Response Matrix Utilities
 - <https://remu.readthedocs.io>
 - Developed for neutrino interaction measurements
 - Builds response matrices and uncertainties from MC
 - Fully developed statistical model of detector, flux, and MC stat uncertainties



DELPHES
fast simulation



<https://iopscience.iop.org/article/10.1088/1748-0221/14/09/P09013>

DISCUSSION STARTERS

- Unregularised result is best approximation of Likelihood
 - e.g. for fits and statistical tests of models
- Regularisation should be used to pick a representative point estimate
 - e.g. for plots
- We should always make likelihood function available
 - Unregularised result or something more complicated
 - Wiener SVD and post-hoc regularisation make this trivially easy
 - Added bonus: regularised and unregularised result are directly related
- Include as many method details as possible in your papers
 - Lots of nuances, caveats, assumptions...
 - Not practical to spell out every single check/study/approximation
 - Or is it?
 - Dedicated method paper?
 - Have to take papers at face value
 - Trust in what is written
 - Assume the worst about what is not written?
 - Assume the best?
 - Hope for the best but expect the worst?



DISCUSSION STARTERS

- Unregularised result is best approximation of Likelihood
 - e.g. for fits and statistical tests of models
- Regularisation should be used to pick a representative point estimate
 - e.g. for plots
- We should always make likelihood function available
 - Unregularised result or something more complicated
 - Wiener SVD and post-hoc regularisation make this trivially easy
 - Added bonus: regularised and unregularised result are directly related
- Include as many method details as possible in your papers
 - Lots of nuances, caveats, assumptions...
 - Not practical to spell out every single check/study/approximation
 - Or is it?
 - Dedicated method paper?
 - Have to take papers at face value
 - Trust in what is written
 - Assume the worst about what is not written?
 - Assume the best?
 - Hope for the best but expect the worst?



“Good physicists do have priors and always use them!
(Only the perfect idiot has no priors.)

– G. D’Agostini
arXiv:1010.0632

“Note that venerable proverb:
Children and fools always speak the truth.

– Mark Twain
On the Decay of the Art of Lying

Thanks!

Backup

EXAMPLE PENALTY MATRICES

$$\tau Q_1 = \tau \begin{pmatrix} 1 & -1 & 0 & 0 & & \\ -1 & 2 & -1 & 0 & \dots & \\ 0 & -1 & 2 & -1 & & \\ 0 & 0 & -1 & 2 & & \\ & \vdots & & & \ddots & \end{pmatrix}$$

Penalise bin-to-bin differences

$$\tau Q_{1m} = \tau \begin{pmatrix} 1/m_1^2 & -1/(m_1 m_2) & 0 & 0 & & \\ -1/(m_1 m_2) & 2/m_2^2 & -1/(m_2 m_3) & 0 & \dots & \\ 0 & -1/(m_2 m_3) & 2/m_3^2 & -1/(m_3 m_4) & & \\ 0 & 0 & -1/(m_3 m_4) & 2/m_4^2 & & \\ & \vdots & & & \ddots & \end{pmatrix}$$

Penalise bin-to-bin model scaling differences

TWO WAYS OF INTERPRETING A

- Coordinate transformation
- New result describes exactly the same distribution, but with different axes
 - No information lost
- Intuitive in 2D
- Axes of histograms no longer make sense

- Modification of result
- Coordinate axes stay the same, but distribution changes
 - Change of result
- Axes and bin values retain same meaning

