

Precision physic at colliders: introducing reSolve

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What is reSolve ?

- **reSolve** is a MC code which aims to add analytic transverse momentum resummation to **generic** processes
- **Generic** doesn't mean quite **any** – limited by the formalism: means (for now at least) **no colour** in the final state
- **Work in progress**, however first trial version already available: → <https://github.com/fkhorad/reSolve>
(many intended features still missing¹)

¹Still young (< 1 year) & small (2 of us)! 

What is reSolve ?

- reSolve is a spiritual successor to the family of codes DYqT, DYres, HqT, Hres, 2gres
- It shares (almost) no lines of code with its predecessors: completely reimaged and reimplemented in C++
- Two updates already in the pipeline: one in the next few days + another on a “soonish” timescale

What is reSolve ?

- Underlying philosophy: **modular**, **transparent**, easily **customisable** code
- Set it up in way which allows (relatively) easy **interfacing** with other codes
- Build it with **parallelisation** already in mind (more on this later)

On q_T resummation

- ... surely does not need a lengthy introduction in this context
- Long story short, logarithmic contributions of the form

$$\sim \log(q_T^2/M^2)^n \alpha_s^m$$

appear to all orders in perturbation theory for $q_T \ll M$.

- They are associated with multiple, (relatively) soft radiation emission and have to be included to get a realistic description of the q_T spectrum of F

On q_T resummation

- A standard way to take this into account is through Parton Showers. **reSolve** uses instead an analytic approach, based on [Catani,Cieri,De Florian,Ferrera,Grazzini('14)]
- History of the formalism is long – dates back to [Dokshitzer,Diakonov,Troian('78)] , [Parisi,Petronzio('79)] , [Kodaira,Trentadue('82)] , [Collins,Soper,Sterman('85)] – much work happened between now and then
- I'll jump straight to the final **master formula** and highlight some of its characteristics, relevant to describe what **reSolve** can potentially do

q_T resummation: Master Formula

- As already specified, F is a **generic non-coloured** final state

$$\frac{d\sigma^{F(res)}(s, q_T, M, y, \Omega)}{d^2q_T dM^2 dy d\Omega} = \frac{M^2}{s} \int \frac{d^2b}{(2\pi)^2} e^{ib \cdot q_T} S_c(M, b) \int_{x_1}^1 \frac{dz_1}{z_1} \int_{x_2}^1 \frac{dz_2}{z_2} d\tilde{\sigma}_{c\bar{c}}^{F; h_1 h_2 \lambda_1 \lambda_2} \cdot C_{ca_1}^{h_1 \lambda_1}(z_1, \alpha_s(b_0^2/b^2)) C_{\bar{c}a_2}^{h_2 \lambda_2}(z_2, \alpha_s(b_0^2/b^2)) f_{a_1/h_1}(x_1/z_1, b_0^2/b^2) f_{a_2/h_2}(x_2/z_2, b_0^2/b^2)$$

with

$$S_c = \exp \left[- \int_{\frac{b_0^2}{b^2}}^{M^2} \frac{dq^2}{q^2} \left(A_c(\alpha_s(q^2)) \log(M^2/q^2) + B_c(\alpha_s(q^2)) \right) \right]$$

$$d\tilde{\sigma}_{c\bar{c}}^{F; h_1 h_2 \lambda_1 \lambda_2} = d\hat{\sigma}_{c\bar{c}}^{F(0)} H_c^{F; h_1 h_2 \lambda_1 \lambda_2}; \quad x_{1,2} = \frac{M}{\sqrt{\hat{s}}} e^{\pm y}$$

$$K_{c,\dots}(\alpha_s) = \sum_n \left(\frac{\alpha_s}{\pi} \right)^n K_{c,\dots}^{(n)}$$

q_T resummation: Master Formula

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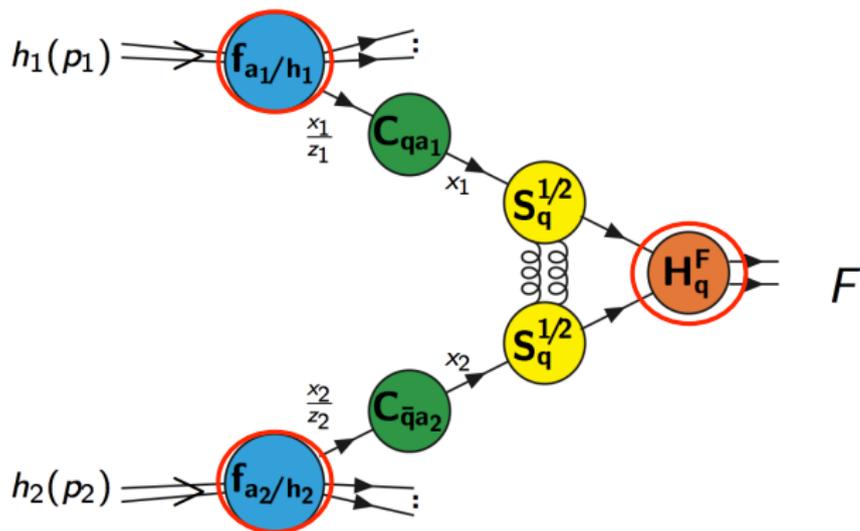
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q_T resummation: Master Formula

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- As highlighted, most factors in master formula are **universal** – which really begs for a **general** numerical implementation
- Exactly as **standard factorisation**, the formula is **inclusive** in unresolved radiation, but **fully differential** in all F internal variables
- The formula is a **small q_T approximation**: it formally holds up to $\mathcal{O}(q_T^2/M^2)$ corrections. A **matching** to fixed order can be defined in a general way in order to handle the **full q_T spectrum** (+ any other observable) in a single formalism.

q_T : matching and subtraction

- The matching strategy – which I won't describe in detail – essentially involves **using a fixed-order expansion of the master formula as a counterterm** to avoid double counting
- Nontrivial additional benefit of the matching is automatically defining a **NNLO subtraction mechanism**. This is **nonlocal** (and obviously only works for final states without colour), but is **immediate to implement** once you have resummation

back to reSolve – usage

- Installation as simple as typing “make” (only needs reasonably recent gcc and gfortran)
- Currently needs the legacy **minuit** library (sources+makefile provided with the code) for PDF fit
- Optionally uses the **CUBA**_[Hahn('04)] library for integration – if desired, has to be downloaded and installed separately
- Main code is C++11-standard compliant – in principle fully portable (tested on several linux distros + MacOSX)

resolve – usage

- Input read from a simple **flexibly-styled, customisable** input card – most settings are rather obvious (couplings, masses, cuts, # of events. . .), all are **documented**
- Straightforward to add more parameters as needed
- Straightforward to add **more processes** (has to be done by hand, though – some templates are provided)

resolve – usage

- Basic output: of course the total cross-section, and distributions as desired (specify on input file)
- Main output are however **LHA-style events**, which can then be analysed with the user's favourite tool(s) – a simple built-in histogrammer tool is also provided

Code structure

- Code is organised in **component modules**, which are as far as possible **independent**
 - ▷ **Main**: read input, select process, call integrations
 - ▷ **Histogrammer**
 - ▷ **Integral**: integration routines
 - ▷ **PDFfit**: currently needed for resummation, kept separate
 - ▷ **Process**: one subfolder per final state
 - ▷ **Diphoton**: SM background only (no Higgs)
 - ▷ **Drell-Yan**
 - ▷ **Resummation**: the main module
 - ▷ **Utility**: some basic general-purpose routines

On universality

- The most important and “unique” part of the code, the **resummation module** makes full use of the **universality** of the formula
- It does not need to know any detail about the underlying process
- Can be compiled and used **independently**
- In principle, this part of the code to be interfaced with any matrix element provider and integrator/event generator

On parallelisation

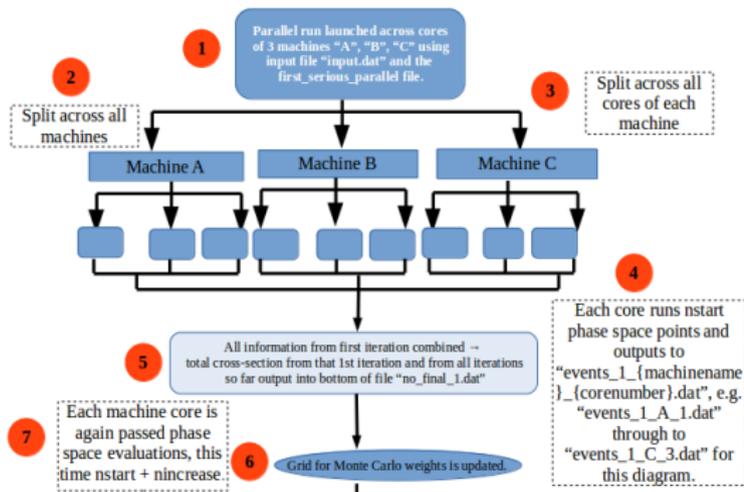
- I stated the code is **set-up for parallelisation** – what do I mean?
- ME are “agnostic” w.r.t. to the integrator: they only use **clearly specified** and **local** variables
- The code by default contains **CUBA**, which attempts to automatically parallelise on all available cores on the local machine (tested on Desktops with 4-8 cores)
- The code also has a dedicated Vegas implementation dubbed **k_vegas** – can also be used as a stand-alone integrator

On parallelisation

- `k_vegas` is just Vegas, but is set-up in order to be able to do **partial iterations**, save partial results for the integration grid, and recombine them at a later stage
- This allows for parallelisation on different **nodes** (→ machines). An example simple bash script provided with the code uses this to parallelise a run over an arbitrary number of computers which are reachable from the central node via **ssh**.
- Of course, much room for improvement (but not so easy if one wants to stay system-independent)

On parallelisation

- Schema of parallelisation using k_vegas + script

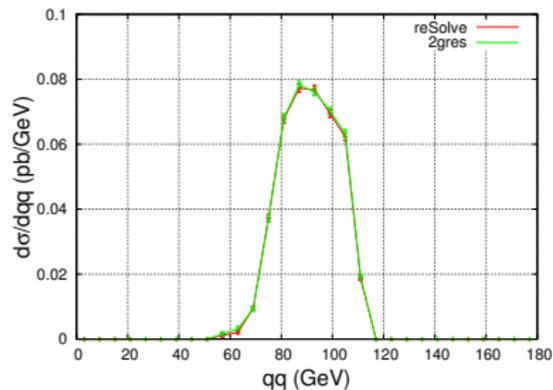
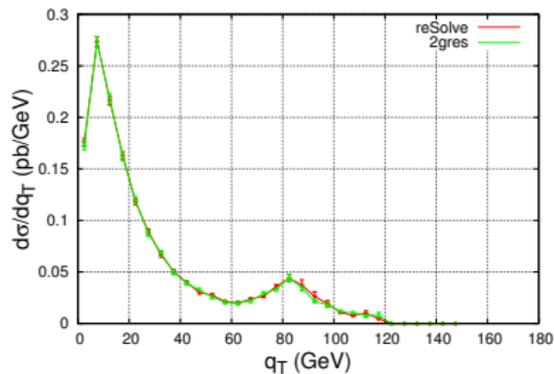


Current state & upgrades

- What the code does now: **diphoton** (SM background), no matching. It **is** however, the first public incarnation of 2gres (used for predictions in [Cieri,FC,De Florian('15)])
- What will definitely do very soon (\sim days): **Drell-Yan**, easier (still non-automated) implementation of new processes (+ improvements not obvious from a user p.o.v.)
- What it will do “soonish”: Higgs production, Higgs resonance-SM background interference in diphoton channel, matching, **QED resummation**

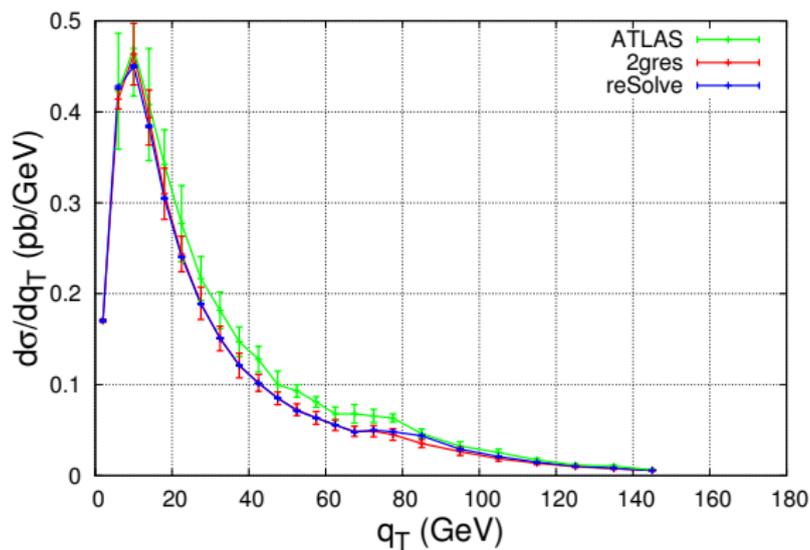
Some validation plots

- Sample comparison with 2gres



Some validation plots

- Comparison with data (includes **matching**, done using **2gNNLO** for the fixed-order part)



“The future” – potential directions

(essentially a personal wishlist)

- **OBVIOUSLY**: add more complete NNLL (or even N^3LL) processes (+ their f.o. for matching)
- Interface with existing ME generators: would allow **automatic NLL+NLO**, even for **BSM** – usual F class only, of course (easy)
- Interface in reverse order: produce a plug-in version to be used **inside** other MC codes

“The future” – potential directions

- Interface with existing **NLO** ME generators for **semi-automatic NNLL+NNLO** (not-so-easy)
- Extension to (at least some) Final States with colour (speculative)
- Improvements on q_T subtraction (speculative)

Conclusions

- Precision calculations for processes at hadron colliders are one of the important challenges which lie ahead the theoretical particle physics community
- We must have the right tools to face the challenge: I hope **reSolve** will be a nice addition to the phenomenologist's (and possibly experimentalists's) toolbox