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Doktoratskolleg Particles and Interactions





# A fast code for automated running and matching of couplings and masses in QCD

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Mini Workshop on Quark Masses, KIT/Online, 2020-10-26

#### Preamble

- **REvolver** is a code written in C++ (+MathLink +Python) to
  - Run quark masses (RGE)
    - Automatic mass decoupling at thresholds
  - Convert quark masses between renormalization schemes
    - Using R-Evolution to resum large logs
    - Including light flavor effects
  - Run QCD coupling (RGE)
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  - Run QCD coupling (RGE)
    - Automatic mass decoupling
  - → To be released very soon (beta available on request)



# Outline

- Motivation
  - Quark Mass Renormalization Schemes
    - Mass Schemes & Conversions: Massless lighter flavors
    - Mass Schemes & Conversions: Massive lighter flavors
- Operating Principle of REvolver
- Live Demo, Documentation Overview

#### **Motivation**

#### Quark Mass Schemes Massless lighter Flavors

# Quark Mass Schemes

• Quark masses:

![](_page_5_Picture_2.jpeg)

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- Confinement → Quark masses not physical observables
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![](_page_6_Picture_2.jpeg)

- Important parameters for SM predictions
- Confinement → Quark masses not physical observables
  - Parameter in QCD action
- Quark self energy UV divergent → Needs to be absorbed into mass
  - Additional finite contributions renormalization scheme dependent
  - Can choose appropriate mass scheme depending on application.
     → Need precise methods to convert between schemes

# Mass Scheme Examples

$$m_Q^{\text{pole}} - \overline{m}_Q = \overline{m}_Q \sum_{n=1}^{\infty} a_n^{\overline{\text{MS}}} \left( \frac{\alpha_s^{(n_\ell + 1)}(\overline{m}_Q)}{4\pi} \right)^n$$

 $[n_{\ell}\cdots \# \text{ massless quark flavors}]$ 

- $\overline{\mathrm{MS}}$  mass:
  - Running mass
  - Analogous to  $\alpha_s(\mu)$ : Absorbs only UV  $1/\varepsilon$  poles from self-energy
  - Intrinsic physical scale:  $\overline{m}_Q$
  - Standard scheme for most high energy applications
  - Only physically relevant for  $\mu \gtrsim \overline{m}_Q$ 
    - Lower scales: Virtual heavy quark effects should be integrated out

## Mass Scheme Examples

$$m_Q^{1S} - m_Q^{\text{pole}} = M_B \sum_{n=1}^{\infty} c_n \left(\frac{\alpha_s^{(n_\ell)}(M_B)}{4\pi}\right)^n$$

• 1S mass:

- Example of low scale mass
- Defined as half of heavy quarkonium spin triplet ground state mass
   → Observable, well defined mass scheme
- Intrinsic physical scale: Inverse Bohr radius  $M_B \equiv C_F \alpha_s m_Q^{
  m pole}$

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  m pole}$
- Many more well defined mass schemes: PS, RS, kinetic, jet, ...

$$m_Q^{\text{pole}} - m_Q^{\text{MSR}}(R) = R \sum_{n=1}^{\infty} a_n \left(\frac{\alpha_s^{(n_\ell)}(R)}{4\pi}\right)^n$$

- MSR mass:
  - Natural extension of  $\overline{\rm MS}$ -mass for scales  $\ll m_Q$
  - Defined directly from pole- $\overline{\mathrm{MS}}$  relation
    - Heavy DOF integrated out  $\rightarrow$  matching to  $\overline{\mathrm{MS}}$
  - Intrinsic physical scale: Adjustable momentum cut-off R

 $\left[a_n = a_n^{\overline{\mathrm{MS}}}(n_h = 0)\right]$ 

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  - Low-scale short-distance mass with direct relation to self-energy diagrams

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- Lighter massive quarks can be incorporated systematically via matching → flavor-number dependent RG evolution → Later

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  - MSR mass contains self-energy corrections only for scales larger than R
- Pole mass: Formal limit  $R \rightarrow 0$ 
  - Absorbs all contributions from quark self energy
     → sensitivity to non-perturbative regions
  - Suffers from  $\mathcal{O}(\Lambda_{QCD})$  renormalon in QCD

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- FOPT: Common scale  $\mu$  and flavor number  $n_f$  have to be used for all  $\alpha_s \rightarrow$  renormalon cancellation
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## Mass Scheme Conversion

- FOPT: Common scale  $\mu$  and flavor number  $n_f$  have to be used for all  $\alpha_s \rightarrow$  renormalon cancellation
  - Potentially large logs when converting mass schemes with different characteristic scales
- Solution: MSR / R-Evolution
  - Utilize freely adjustable intrinsic scale R
  - As intermediate step RG running between scales

## R-Evolution

$$R\frac{\mathrm{d}m_Q^{\mathrm{MSR}}(R)}{\mathrm{d}R} = -R\gamma^R(\alpha_s(R)) = -R\sum_{n=0}^{\infty}\gamma_n^R\left(\frac{\alpha_s(R)}{4\pi}\right)^{n+1}$$

- RGE in IR scale R, relating MSR masses at different scales
  - Sums systematically renormalon series and large logs
  - Linear dependence on R

#### R-Evolution

• MS-scheme running: Large scale ( $\overline{\mathrm{MS}}$ ) and Low scale (MSR)

![](_page_18_Figure_2.jpeg)

## R-Evolution

Fixed order vs. R-Evolution

![](_page_19_Figure_2.jpeg)

![](_page_20_Picture_0.jpeg)

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- Many applications: Lighter massive quarks set massless
  - Needed in high precision calculations
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  - Needed in high precision calculations
- Corrections known for some schemes
- MSR scheme: Systematic treatment via flavor matching

![](_page_23_Figure_5.jpeg)

- Scheme conversions & light massive flavors
  - Impact at higher orders
    - → Massive virtual quark loops act as IR-cutoff
    - → Change renormalon structure

![](_page_24_Picture_5.jpeg)

- Scheme conversions & light massive flavors
  - Impact at higher orders
    - → Massive virtual quark loops act as IR-cutoff
    - → Change renormalon structure
- Consistency is crucial
  - when converting to/from pole mass
  - when converting between schemes

#### REvolver

#### **Features**

## REvolver Features

- All implemented in **REvolver**:
  - MSR scheme as low-scale extension of  $\overline{\mathrm{MS}} \rightarrow$  "MS-scheme"
    - R-Evolution
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    - using R-Evolution to resum logs
    - including light massive flavor contributions (if known)
    - Various schemes supported: MS, 1S, PS, RS, kinetic, pole, ...
- Highest available orders implemented for running/matching
  - 5 loop  $\alpha_s$ , 5/4 loop  $\overline{\mathrm{MS}}/\mathrm{MSR}, \dots$

#### **REvolver**

#### **Operating Principle**

 $n_f^{\text{tot}}, \alpha_s^{(n_\alpha)}(\mu_\alpha), \{m_i^{(n_{m_i})}(R_i)\}, \text{optional parameters}$ 

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Core

 $n_{f}^{\text{tot}}, n_{f}^{m=0}$  $m_{0}^{(n_{0})}(\overline{m}_{0}), m_{0}^{(n_{1})}(\overline{m}_{1}), \dots$  $m_{1}^{(n_{0})}(\overline{m}_{0}), m_{1}^{(n_{1})}(\overline{m}_{1}), \dots$ 

 $\alpha_s^{(n_0)}(\overline{m}_0), \, \alpha_s^{(n_1)}(\overline{m}_1), \dots$ optional parameters  $\Lambda_{\text{QCD}}^{(n_0)}, \dots$ perturbative coefficients, ...

 $n_f^{\text{tot}}, \alpha_s^{(n_\alpha)}(\mu_\alpha), \{m_i^{(n_{m_i})}(R_i)\}, \text{optional parameters}$ 

Request quantity via member function

Core  $n_f^{\text{tot}}, n_f^{m=0}$  $m_0^{(n_0)}(\overline{m}_0), m_0^{(n_1)}(\overline{m}_1), \ldots$  $m_1^{(n_0)}(\overline{m}_0), m_1^{(n_1)}(\overline{m}_1), \ldots$  $\alpha_{s}^{(n_{0})}(\overline{m}_{0}), \, \alpha_{s}^{(n_{1})}(\overline{m}_{1}), \, \dots$ optional parameters  $\Lambda^{(n_0)}_{\text{OCD}}, \dots$ perturbative coefficients, ...

 $n_f^{\text{tot}}, \alpha_s^{(n_\alpha)}(\mu_\alpha), \{m_i^{(n_{m_i})}(R_i)\}, \text{optional parameters}$ 

Core

Request quantity via member function

![](_page_34_Figure_3.jpeg)

perturbative coefficients, ...

 $\longrightarrow m_i^{(n_f)}(\mu)$ 

 $\blacktriangleright \alpha_s^{(n_f)}(\mu)$ 

 $\longrightarrow m_i^{\mathrm{PS}}(\mu_f)$ 

#### • Three ways to use the code:

#### - directly via C++ library

- Fast
- Most suitable for extensive, automated tasks
- Interaction with other C++ libraries

- via WSTP/MathLink in Mathematica
  - User friendly
  - Most suitable for interactive tasks
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