

A precise α_s determination from the R-improved QCD Static Energy

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Introduction

- α_s most important parameter in QCD \rightarrow should be determined with high precision.
- Our strategy \rightarrow comparing the QCD Static Energy obtained in lattice simulations with highly accurate perturbative results.
- $V_{\text{QCD}} \propto \alpha_s \rightarrow$ very sensitive.
- We improve this method building on previous analyses [A. Bazavov, N. Brambilla, X. Garcia i Tormo et al, 2012, A. Bazavov, N. Brambilla, X. Garcia i Tormo et al, 2014] in several ways:
 - Leading renormalon subtraction \rightarrow short-distance scheme (MSR).
 - Resummation of associated large logs with R-evolution.
 - Profiles functions for the renormalization scales.
- We can fit (for the first time) lattice data up to $r \sim 0.5$ fm.

Static Energy and Static Potential

- The Static Energy is defined as the potential energy between an infinitely massive quark anti-quark pair at a distance r , corrected by ultra-soft effects. In pNRQCD

$$E_s(r) = V_s(r, \mu) + \delta_{\text{us}}(r, \mu).$$

- The Static Potential is the basic object to understand the behavior of non-relativistic QCD:

$$V_s(r, \mu) = V_s^{\text{soft}}(r, \mu) + V_s^{\text{us}}(r, \mu),$$
$$V_s^{\text{soft}}(r) = -C_F \frac{\alpha_s(\mu)}{r} \sum_{i=0}^3 \left[\frac{\alpha_s(\mu)}{4\pi} \right]^i \sum_{j=0}^i a_{ij} \log^j(r\mu e^{\gamma_E}).$$

- Coefficients a_{i0} are known to three loops. $a_{ij \geq 0}$ obtained with RGE. $\alpha_s(\mu) = \alpha_s^{(n_\ell=3)}(\mu)$.
- At $\mathcal{O}(\alpha_s^4)$ ultra-soft contributions show up for the first time

$$V_s^{\text{us}}(r, \mu) = -\frac{C_A^3 C_F}{12\pi} \frac{\alpha_s^4(\mu)}{r} \log(\mu r e^{\gamma_E}).$$

Ultra-soft term and resummation

- μ_{us} appears both in the ultrasoft static potential V_s^{us} and in the matrix element δ_{us}

$$\delta_{us}(\mu_s, \mu_{us}) = -\frac{C_A^3 C_F}{12\pi} \frac{\alpha_s^3(\mu_s) \alpha_s(\mu_{us})}{r} \log \left[\frac{C_A \alpha_s(\mu_s) e^{-5/6}}{\mu_{us} r} \right]$$

$$V_s^{us}(r, \mu_s, \mu_{us}) = -\frac{C_A^3 C_F}{12\pi} \frac{\alpha_s^3(\mu_s) \alpha_s(\mu_{us})}{r} \log(\mu_{us} r e^{\gamma_E})$$

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- One can see the necessity of resummation with these two (incompatible) choices that minimize logs.
 - $\mu_{us} \sim \alpha_s/r$ [use this one: no large logs in δ_{us}]
 - $\mu_{us} \sim 1/r$ [discarded: sum up logs in V_s^{us}]

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- We perform resummation using pNRQCD RGE

$$\mu \frac{dV_s(r, \mu_s, \mu)}{d\mu} = -\frac{2C_F C_A^3}{24r} \frac{\alpha_s(\mu)}{\pi} \left[1 + B \frac{\alpha_s(\mu)}{\pi} \right] \alpha_s^3(\mu_s) \left\{ 1 + 3 \frac{\alpha_s(\mu_s)}{4\pi} \left[a_{1,0} + 2\beta_0 \log(r\mu_s e^{\gamma_E}) \right] \right\}.$$

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- Integrating from $\mu = \mu_s \sim 1/r$ to $\mu = \mu_{us} \sim \alpha_s/r$

$$V_s(r, \mu_s, \mu_{us}) = V_s(r, \mu_s) + U_{us}(r, \mu_s, \mu_{us}),$$
$$U_{us}(r, \mu_s, \mu_{us}) = \frac{C_A^3 C_F}{6\beta_0 r} \alpha_s^3(\mu_s) \left\{ \left(1 + 3 \frac{\alpha_s(\mu_s)}{4\pi} \left[a_{1,0} + 2\beta_0 \log(r\mu_s e^{\gamma_E}) \right] \right) \log \left[\frac{\alpha_s(\mu_{us})}{\alpha_s(\mu_s)} \right] + \left(B - \frac{\beta_1}{4\beta_0} \right) \left[\frac{\alpha_s(\mu_{us})}{\pi} - \frac{\alpha_s(\mu_s)}{\pi} \right] \right\}$$

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- Replacing in the static energy \rightarrow no large logs for $\mu_s \sim 1/r$ and $\mu_{us} \sim \alpha_s/r$

$$E_s(r) = V_s^{\text{soft}}(r, \mu_s) + \frac{C_A^3 C_F}{12} \frac{\alpha_s^3(\mu_s)}{r} \left\{ \frac{2}{\beta_0} \left(B - \frac{\beta_1}{4\beta_0} \right) \left[\frac{\alpha_s(\mu_{us})}{\pi} - \frac{\alpha_s(\mu_s)}{\pi} \right] \right. \\ \left. + \frac{2}{\beta_0} \left(1 + 3 \frac{\alpha_s(\mu_s)}{4\pi} \left[a_{1,0} + 2\beta_0 \log(r\mu_s e^{\gamma_E}) \right] \right) \log \left[\frac{\alpha_s(\mu_{us})}{\alpha_s(\mu_s)} \right] \right. \\ \left. - \frac{\alpha_s(\mu_{us})}{\pi} \log \left[\frac{C_A \alpha_s(\mu_s) e^{-5/6}}{r\mu_{us}} \right] - \frac{\alpha_s(\mu_s)}{\pi} \log(r\mu_s e^{\gamma_E}) \right\}.$$

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- To cancel the renormalon we need a short-distance mass scheme \rightarrow MSR mass [A. H. Hoang, A. Jain, C. Lepenik, V. Mateu, I. Scimemi and I. W. Stewart, 2018]

$$\begin{aligned} \delta m_Q^{\text{MSR}}(\mu, R) &\equiv m_Q^{\text{pole}} - m_Q^{\text{MSR}}(R) = R \sum_{n=1}^{\infty} \delta_n^R \left[\frac{\alpha_s(R)}{4\pi} \right]^n \\ &= R \sum_{n=1}^{\infty} \left[\frac{\alpha_s(\mu)}{4\pi} \right]^n \sum_{j=0}^{n-1} \delta_{nj}^R \log^j \left(\frac{\mu}{R} \right). \end{aligned}$$

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- **New type of large logs** \rightarrow resummation using R-evolution [A. H. Hoang, A. Jain, I. Scimemi and I. W. Stewart, 2008].

Previous works

- To see the advantages of R-evolution we go back to the first work that used the static energy to obtain α_s [A. Bazavov, N. Brambilla, X. Garcia i Tormo et al, 2012]

$$E_s(r) = V_s(r, \mu_s, \mu_{us}) + \delta_{US}(r, \mu_s, \mu_{us}) + RS(\rho),$$

- $V_s(r, \mu_s, \mu_{us}) \rightarrow \ln(r\mu_{us})$
- $RS(\rho) \rightarrow \ln(\rho/\mu_{us})$
- There is no right choice for $\mu_{us} \rightarrow$ fits only possible for small values of r .
- The force was used in [TUMQCD collaboration, 2019]

$$F_s(r) = \frac{dE_s(r)}{dr}$$

- Force avoids explicit subtraction.
- Using fully canonical scales $\sim 1/r \rightarrow$ limits fit-range to $r \sim 0.076$ fm.

R-improvement, renormalon subtractions

- The next goal is to obtain a renormalon-free potential.

$$E_{Q\bar{Q}}(r) = 2m_Q^{\text{MSR}}(R_0) + 2\delta m_Q^{\text{MSR}}(R_0, \mu) + E_s(r) \equiv 2m_Q^{\text{MSR}}(R_0) + V_s^{\text{MSR}}(r, \mu, R_0).$$

- We sum up large logs of μ/R_0 .

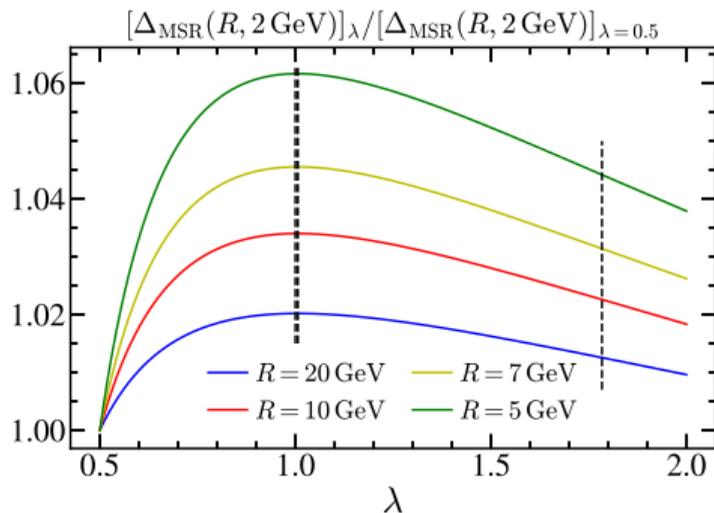
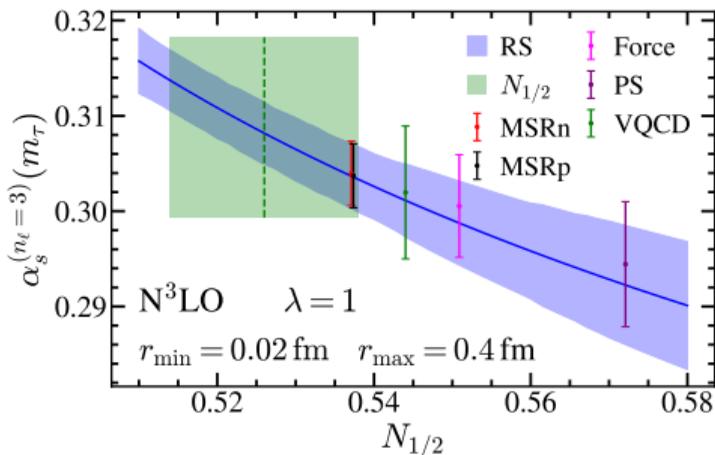
$$\begin{aligned}\delta m_Q^{\text{MSR}}(R_0) &= \delta m_Q^{\text{MSR}}(R_0) + \delta m_Q^{\text{MSR}}(R) - \delta m_Q^{\text{MSR}}(R) \\ &= \delta m_Q^{\text{MSR}}(R) + m_Q^{\text{MSR}}(R) - m_Q^{\text{MSR}}(R_0) \\ &= \delta m_Q^{\text{MSR}}(R) + \Delta^{\text{MSR}}(R, R_0).\end{aligned}$$

- $\Delta^{\text{MSR}}(R, R_0)$: solution to MSR mass R-RGE \rightarrow sums up logs of R/R_0 .
- We have $\delta m_Q^{\text{MSR}}(R) \sim \log(\mu/R)$. By choosing $\mu \sim R \rightarrow$ no large logs.
- We define the R-improved static potential:

$$V_s^{\text{MSR}}(r, \mu, R_0) = V_s(r, \mu) + 2\delta^{\text{MSR}}(R, \mu) + 2\Delta^{\text{MSR}}(R, R_0).$$

Dependence on λ

- We don't know the R-evolution kernel up to infinite order \rightarrow can estimate the truncation error with a dimensionless parameter λ , its variation account for higher-order remnants (similar to renormalization scale variation).
- α_s is sensitive to $N_{1/2}$ (also depends on λ), we have studied the R-evolution dependence with λ .



- The canonical value of $\lambda = 1$ is clearly biased.
- We pick $\lambda = 1.784$ and vary it from 1.5 to 2.1.

Profiles

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- Solution: use profile functions that ensure series convergence:

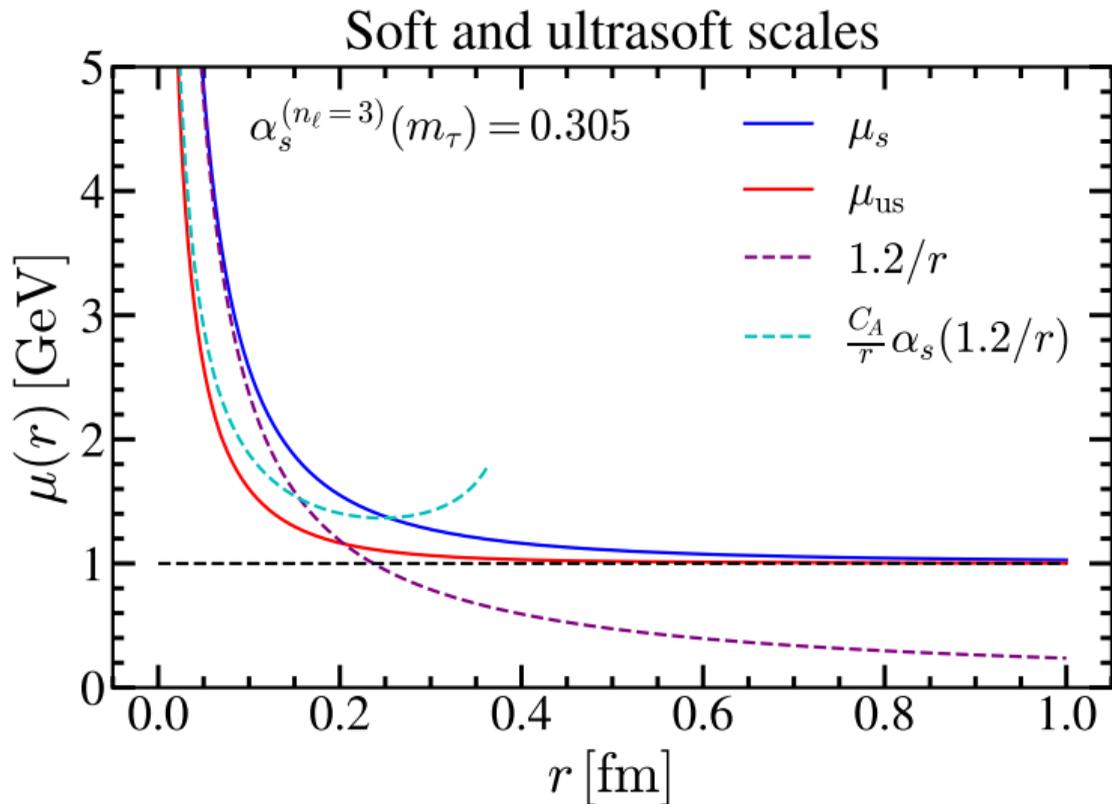
$$\mu_s(r, \xi, \mu_\infty, \Delta, b) = \sqrt{\left(\frac{\xi}{r}\right)^2 + \frac{b}{r} + (\mu_\infty - \Delta)^2} - \Delta = \begin{cases} \frac{\xi}{r} & \text{for } r \rightarrow 0 \\ \mu_\infty & \text{for } r \rightarrow \infty \end{cases}$$

$$R(r, \beta, R_\infty, \Delta, b) = \mu_s(r, \beta, R_\infty, \Delta, b)$$

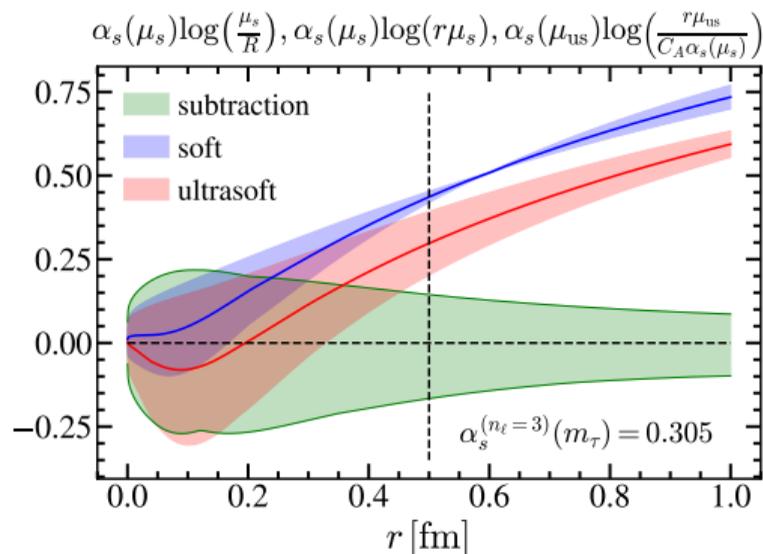
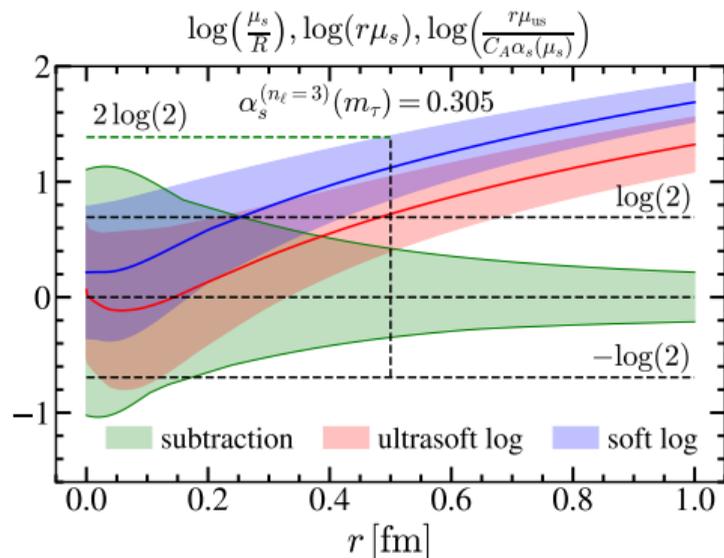
$$\mu_{\text{us}}(r, \xi, \kappa, \mu_\infty, \Delta, b) = C_A \left\{ \mu_s(r, \kappa\xi, \mu_\infty, \Delta, b) \alpha_s[\mu_s(r, \kappa\xi, \mu_\infty, \Delta, b)] - \mu_\infty \alpha_s(\mu_\infty) \right\} + \mu_\infty$$

with $\xi = \mathcal{O}(1)$, $\mu_\infty \sim 1 \text{ GeV}$ and $\mu_s \gg \mu_{\text{us}}$ for small distances. This makes sure the series is stable and convergent over the entire spectrum.

Profiles

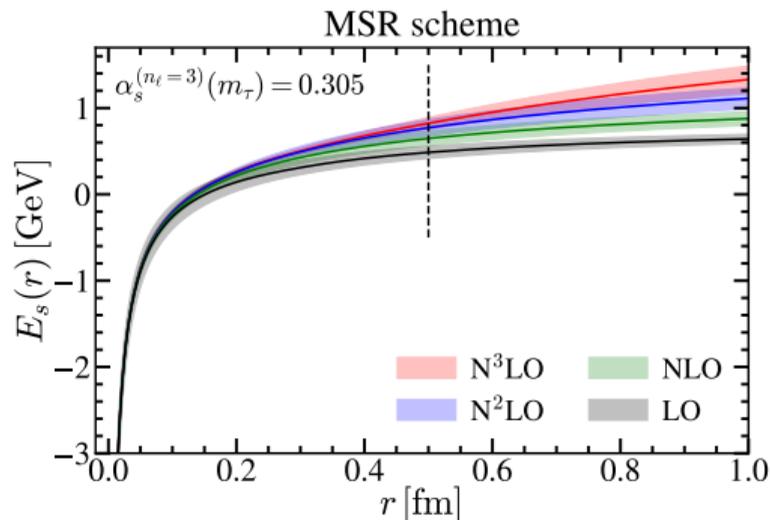
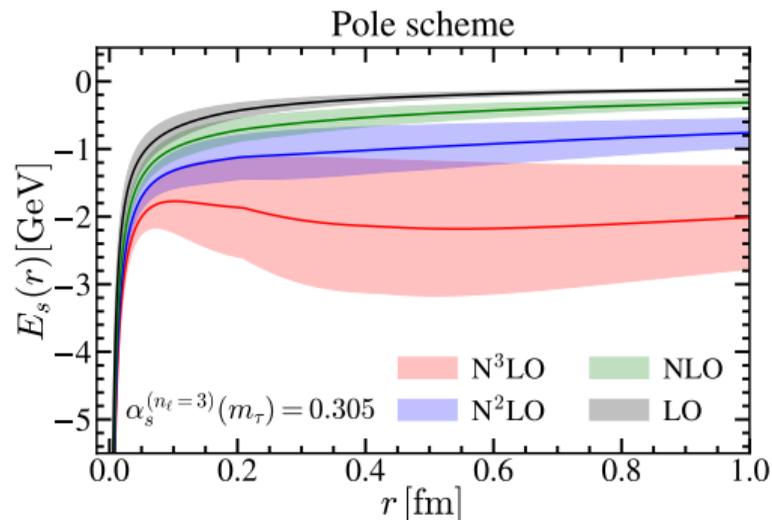


Profiles



- Ultrasoft logarithm smaller than the soft one and $\mathcal{O}(1)$ in the fit range.
- The product of $\alpha_s(\mu) \times \log(\mu)$ is smaller than 1.

Perturbative analysis



At $r = 0.3$ fm we have the following uncertainties for LO, NLO, N²LO and N³LO:

- Pole scheme \rightarrow [0.16, 0.31, 0.71, 1.86] GeV.
- MSR scheme \rightarrow [0.17, 0.15, 0.13, 0.07] GeV.

Lattica data

- Lattice QCD data from HotQCD. $(2 + 1)$ flavor simulations. We have 9 lattices sets adding up to 2512 data points.
- Only diagonal part of statistic covariance matrix publicly available.
- Small- r data super precise but, potentially affected by large discretisation errors.
- To change from lattice to physical units we use $r_1 = 0.3093(20)$ fm, average of [A. Bazavov et al, 2014] and [R. Larsen, S. Mukherjee, P. Petreczky et al, arXiv 2502.08061].

Fits

- Each set n has a different origin of the static energy \rightarrow offset A_n , included in the χ^2 function:

$$\chi^2(\alpha_s, \{A_k\}) = \sum_{k=1}^{N_s} \chi_k^2(\alpha_s, A_k) \quad \chi_k^2(\alpha_s, A_k) = \sum_{i=1}^{n_{\text{data}}^k} \frac{[V_{ik}(\alpha_s) + A_k - V_{ik}^{\text{exp}}]^2}{[\sigma_k^2]_i}$$

- We can marginalize analytically first with respect to the offsets.

$$\begin{aligned} \frac{\partial \chi^2}{\partial A_k} = 0 &\implies \tilde{A}_k(\alpha_s) = \frac{\sum_{i=1}^{n_{\text{data}}^k} [\sigma_k^{-1}]_i^2 [V_{ik}^{\text{exp}} - V_{ik}(\alpha_s)]}{\sum_{i=1}^{n_{\text{data}}^k} [\sigma_k^{-1}]_i^2}, \\ &\implies \tilde{\chi}_k^2(\alpha_s) = \chi_k^2(\alpha_s, \tilde{A}_k(\alpha_s)) - \frac{\left\{ \sum_{i=1}^{n_{\text{data}}^k} [\sigma_k^{-1}]_i [V_{ik}(\alpha_s) - V_{ik}^{\text{exp}}] \right\}^2}{\sum_{i=1}^{n_{\text{data}}^k} [\sigma_k^{-1}]_i}. \end{aligned}$$

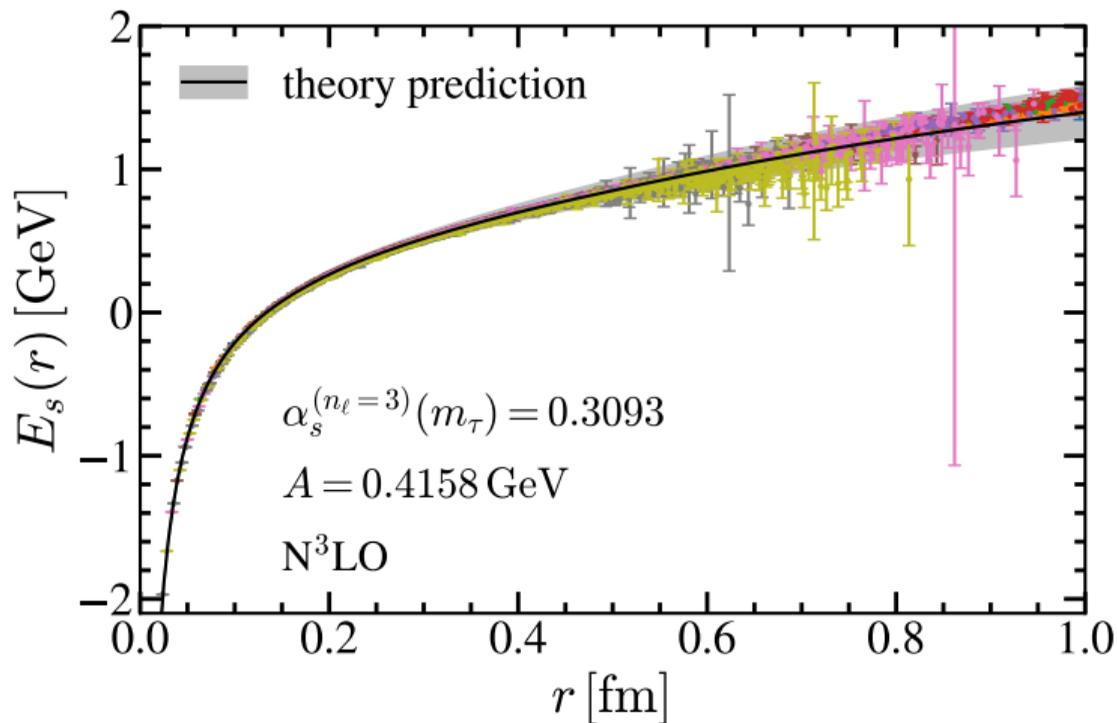
Fits

- Minimizing we obtain α_s^{BF} that verifies $\tilde{\chi}^2(\alpha_s^{\text{BF}}) = \chi^2(\alpha_s^{\text{BF}}, A^{\text{BF}}) = \chi_{\min}^2$.
- The fit uncertainty is defined as $\tilde{\chi}^2(\alpha_s^{\text{BF}} \pm \Delta_{\text{fit}}\alpha_s) = \chi_{\min}^2 + 1$.
- Theory uncertainties are estimated through a random scan.
- We perform the fit for 500 random profiles $\rightarrow \alpha_s^{\text{BF}}$ and $\Delta_{\text{fit}}\alpha_s$ for all of them.
- The theory (or perturbative) uncertainty is computed as

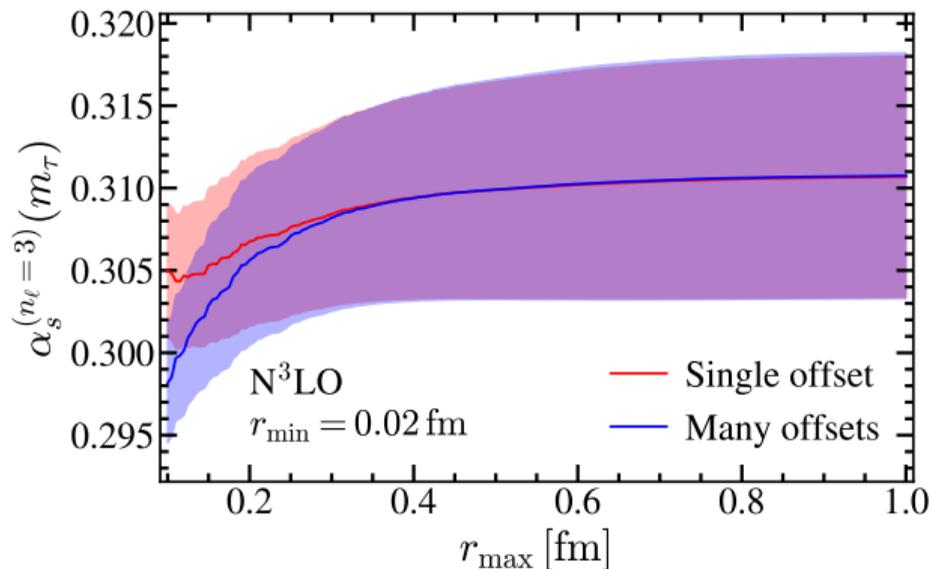
$$\Delta_{\text{pert}}\alpha_s = (\alpha_{\max} - \alpha_{\min})/2$$

- The final result is the mean of the 500 α_s^{BF} and $\Delta_{\text{fit}}\alpha_s$.
- We obtain $\mathcal{O}(100)$ χ^2 values \rightarrow inflate lattice error by $\sqrt{\chi_{\min}^2/\text{d.o.f.}}$.
- Fit for $r \in [r_{\min}, r_{\max}]$. $r_{\text{data}} \geq 0.023$ fm.
- $r_{\min} \in [0.02, 0.045]$ fm and $r_{\max} \in [0.1, 1]$ fm with 0.05 fm spacing \rightarrow 1086 different datasets.

Static Energy

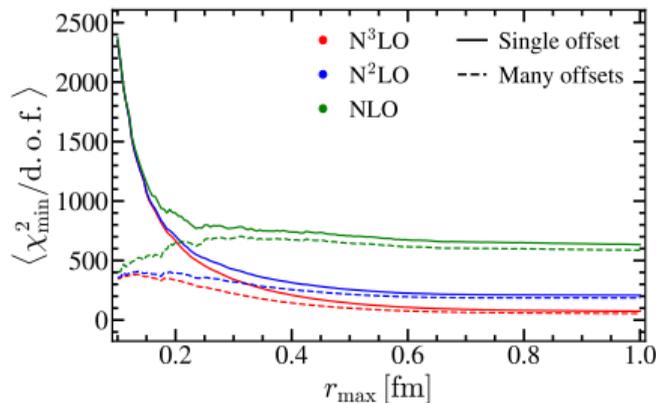
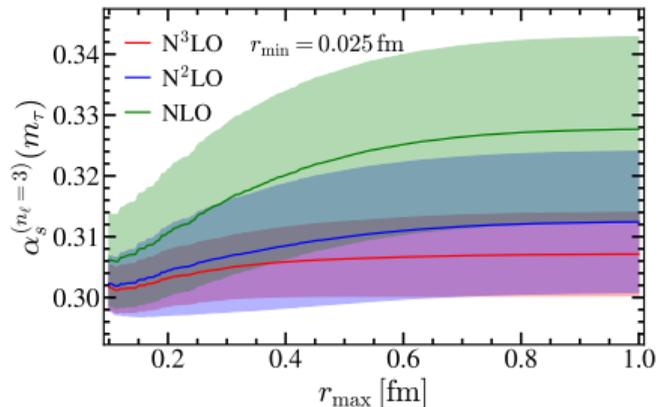
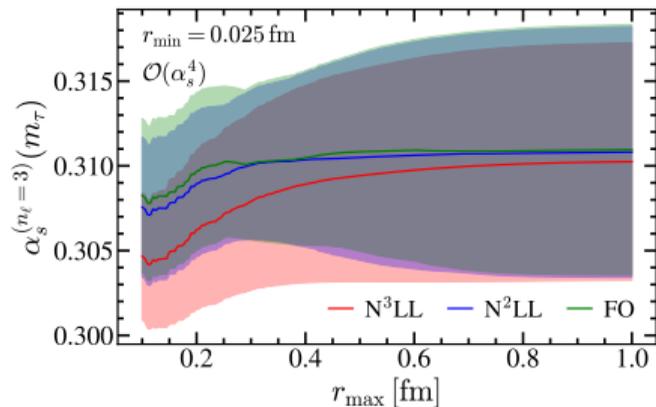


Offset approach



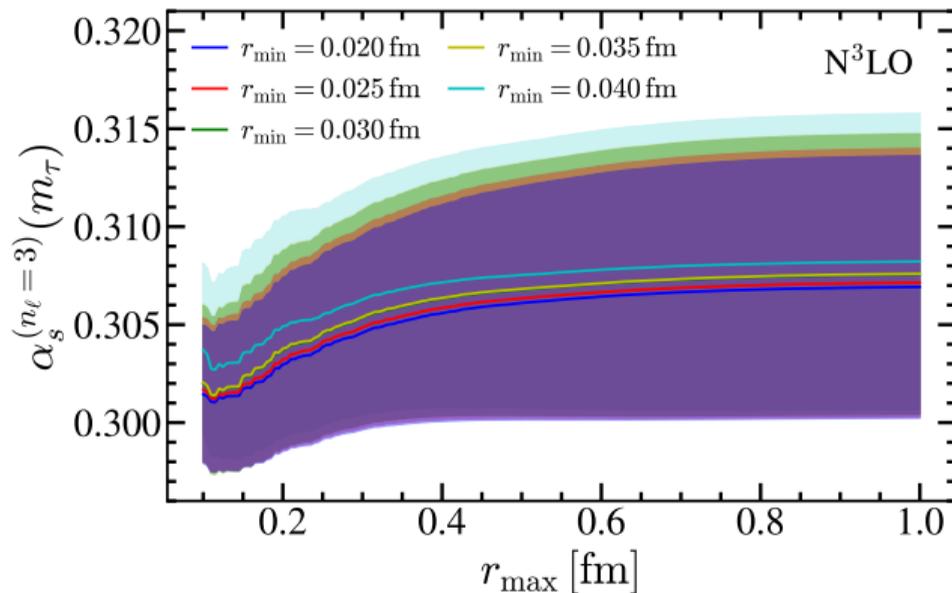
- We consider two offsets approaches:
 - Each ensemble has a different offset.
 - Common offset for all the ensembles.
- Both agree for $r_{\max} > 0.3$ fm.
- Overparametrization of χ^2 at small distances for the many offset case.
- We consider $r_{\max} \geq 0.35$ fm.

Order-by-order agreement



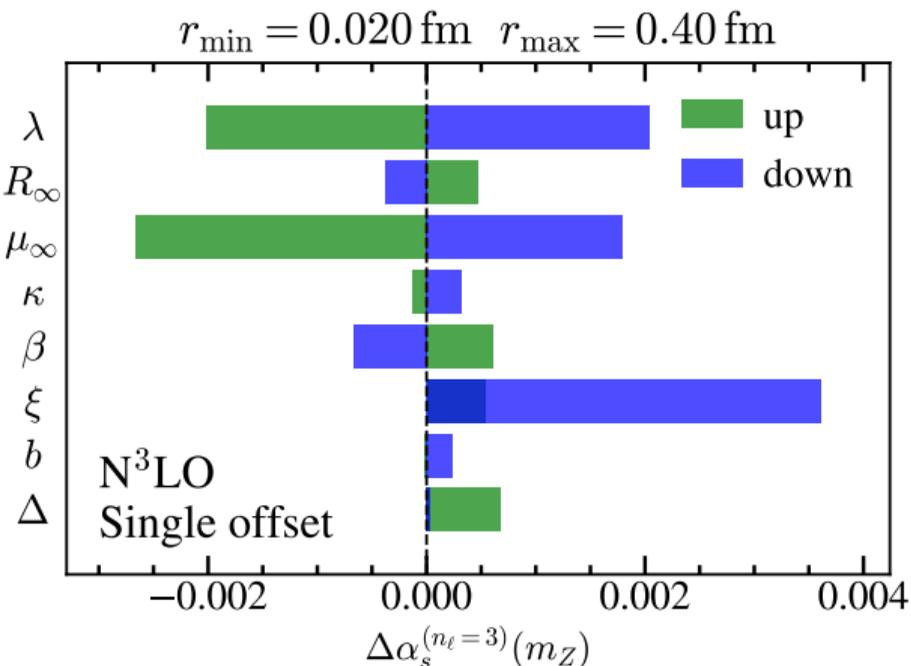
- Uncertainty bands nested for $r_{\max} < 0.5$ fm.
- We choose $r_{\max} \leq 0.45$ fm

Study of the dataset



- We consider the range $0.02 \leq r_{\min} \leq 0.04$ fm .

Variation of profiles parameters one at a time



Parameter	Default	Range
ξ	$1.2 \times \hbar c$	$[0.7, 2.2] \times \hbar c$
β	$1.2 \times \hbar c$	$[0.7, 2.2] \times \hbar c$
b	0	$[-0.3, 0.3] \times \hbar c$
Δ	0	$[-0.6, 0.6] \text{ GeV}$
μ_∞	1 GeV	$[0.9, 1.1] \text{ GeV}$
R_∞	1 GeV	$[0.9, 1.1] \text{ GeV}$
κ	1	$[0.8, 1.2]$
λ	1.8	$[1.5, 2.1]$

Final results

- Requirements: $r_{\min} \in [0.02, 0.04]$ fm and $r_{\max} \in [0.35, 0.45]$ fm \rightarrow 105-element dataset.
- We perform a fit in all of them:

$$\alpha_s^{(n_f=3)}(m_\tau) = 0.3093 \pm 0.00001_{\text{lattice}} \pm 0.0061_{\text{th}} \pm 0.0011_{\text{set}} \pm 0.0011_{r_1}$$

\downarrow

$$\alpha_s^{(n_f=3)}(m_\tau) = 0.3093 \pm 0.0063$$

$$\alpha_s^{(n_f=5)}(m_Z) = 0.1170 \pm 0.0008_{\text{th}} \pm 0.0001_{\text{set}} \pm 0.0001_{r_1} \pm 0.0003_{\mu_c} \pm 0.0002_{\mu_b}$$

\downarrow

$$\alpha_s^{(n_f=5)}(m_Z) = 0.1170 \pm 0.0009$$

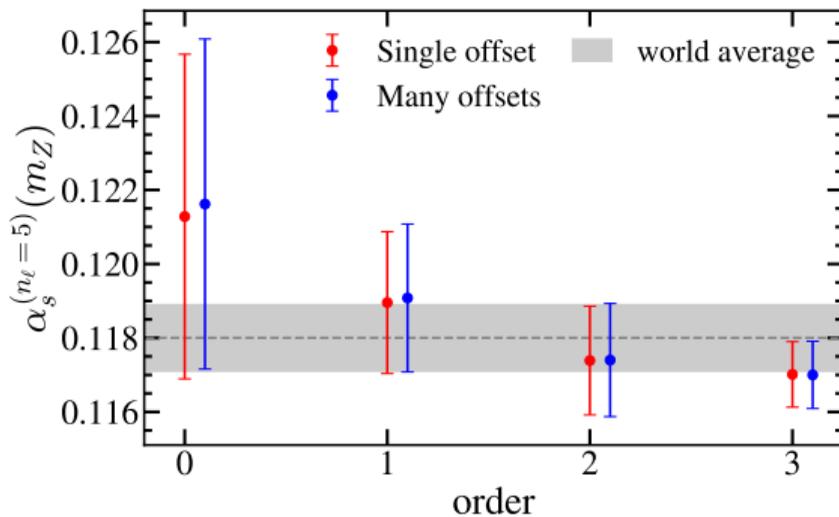
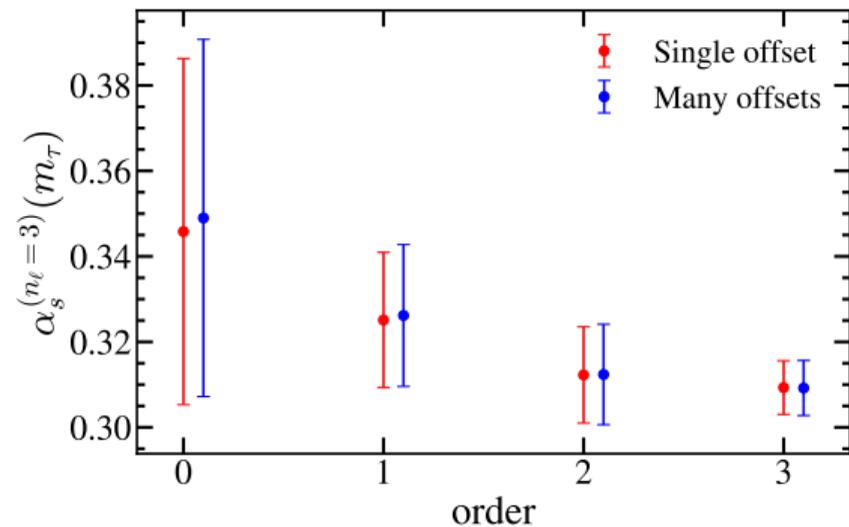
Competitive with the w.a. $\alpha_s^{(5)}(m_Z) = 0.1180 \pm 0.0009$ and compatible at 0.5- σ level.

Comparing with previous analyses:

$$\alpha_s^{(n_f=5)}(m_Z) = 0.11660_{-0.00056}^{+0.00110} \rightarrow \alpha_s^{(n_f=5)}(m_Z) = 0.1162(8) \text{ [TUMQCD collaboration, 2019]}$$

$$\alpha_s^{(n_f=5)}(m_Z) = 0.1181 \pm 0.0009 \rightarrow \alpha_s^{(n_f=5)}(m_Z) = 0.1177(9) \text{ [C. Ayala, X. Lobregat, A. Pineda, 2020]}$$

Final results



Conclusions

- Used the static energy to determine α_s , building (and improving) on previous analyses.
- Performed ultra-soft large-log resummation up to N³LL.
- Employed the MSR mass and R-evolution to improve the static potential.
- Designed profile functions to increase the validity of the potential up to $r \sim 0.5$ fm.
- Studied carefully the dependence on the dataset.
- Carried out fits to lattice data to obtain a very competitive result for α_s .

$$\alpha_s^{(n_f=5)}(m_Z) = 0.3093 \pm 0.0063 \quad \alpha_s^{(n_f=5)}(m_Z) = 0.1170 \pm 0.0009$$

BACKUP

Force-type subtractions

- Integrating the Force is equal to perform R-evolution.
- The renormalon doesn't depend on $r \rightarrow$ we can subtract the potential at r_0 .

$$\begin{aligned} E_s^F(r, r_0) &\equiv E_s(r) - E_s(r_0) = E_s(r) - E_s(r_1) + [E_s(r_1) - E_s(r_0)] \\ &= E_s(r) - E_s(r_1) + \int_{r_0}^{r_1} dr' F_s(r') \equiv E_s(r) - E_s(r_1) + \Delta_F(r_0, r_1) \end{aligned}$$

- [A. Bazavov, N. Brambilla, X. Garcia i Tormo et al, 2014] chooses $r_1 = r \rightarrow$ only Δ_F left.
- Connecting with R-evolution, the subtraction term is (choosing $R = 1/r_1$ and $\mu = R$):

$$E_s(r_1) \equiv \delta_{\text{soft}}^F(R) = \frac{1}{2} V_s^{\text{soft}}\left(\frac{1}{R}\right) = -2\pi C_F R \sum_{i=1} \left[\frac{\alpha_s(R)}{4\pi} \right]^i \sum_{j=0}^{i-1} a_{i-1,j} \gamma_E^j \equiv R \sum_{i=1} \left[\frac{\alpha_s(R)}{4\pi} \right]^i \delta_i^F.$$

- We can express Δ_F as an R-evolution integral:

$$\begin{aligned} \gamma_{\text{soft}}^F(R) &= -\frac{1}{2} \left[r^2 F_s^{\text{soft}}(r) \right]_{r=1/R}, \\ \Delta_F^{\text{soft}}(r_0, r_1) &= \int_{r_0}^{r_1} \frac{dr'}{(r')^2} \left[(r')^2 F_s^{\text{soft}}(r') \right] = \frac{1}{2} \int_{1/r_0}^{1/r_1} dR' \gamma_{\text{soft}}^F(R'). \end{aligned}$$

- It inherits the infrared sensitivity of the Static Potential.

Potential-type subtractions

- If we choose $\mu = R$ and $R = \frac{e^{-\gamma E}}{r}$

$$\delta_{\text{soft}}^V(R) \equiv V_s^{\text{soft}}\left(\frac{e^{-\gamma E}}{R}\right) = -2\pi e^{\gamma E} C_F R \sum_{i=1} \left[\frac{\alpha_s(R)}{4\pi}\right]^i a_{i-1,0} \equiv R \sum_{i=1} \left[\frac{\alpha_s(R)}{4\pi}\right]^i \delta_i^V.$$

- We can obtain their anomalous dimensions

$$\gamma_{\text{soft}}^V(R) = -\frac{1}{2} \left[r^2 V_s^{\text{soft}}(r) \right]_{r=1/R}$$
$$\Delta_V^{\text{soft}}(r_0, r_1) = \int_{r_0}^{r_1} \frac{dr'}{(r')^2} \left[(r')^2 V_s^{\text{soft}}(r') \right] = \frac{1}{2} \int_{1/r_0}^{1/r_1} dR' \gamma_{\text{soft}}^V(R')$$

Subtraction schemes: PS scheme

- Defined from its relation to the pole mass: $m_p - m^{\text{PS}}(R) \equiv \delta_{\text{PS}}(R)$

$$\delta_{\text{soft}}^{\text{PS}}(R) = -\frac{1}{2} \int_{|\vec{q}| < R} \frac{d^3 \vec{q}}{(2\pi)^3} \tilde{V}_s^{\text{soft}}(q, R) \equiv C_F R \frac{\alpha_s(R)}{\pi} \sum_{i=0} \left[\frac{\alpha_s(R)}{4\pi} \right]^i c_i,$$

$$c_i = \sum_{j=0}^i a_{i,j} h_j, \quad \text{with} \quad h_j = j! \sum_{\ell=0}^j \sum_{k=0}^{\text{floor}[\frac{\ell}{2}]} \kappa_{\ell-2k} \left(\frac{\pi}{2}\right)^{2k} \frac{(-1)^k}{(2k)!}.$$

Subtraction schemes: ultrasoft terms

- Since all these three schemes come from the potential, they have ultrasoft terms in their R-evolution.

$$\gamma^F(R) = \gamma_{\text{soft}}^F(R) - \frac{C_A^3 C_F}{24\pi} [\alpha_s(R)]^4 \log \left[C_A \alpha_s(R) e^{\gamma_E - 5/6} \right],$$

$$\gamma^V(R) = \gamma_{\text{soft}}^V(R) - \frac{C_A^3 C_F}{24\pi} e^{\gamma_E} [\alpha_s(R)]^4 \log \left[C_A \alpha_s(R) e^{\gamma_E - 5/6} \right],$$

$$\gamma^{\text{PS}}(R) = \gamma_{\text{soft}}^{\text{PS}}(R) - \frac{C_F C_A^3}{12\pi^2} [\alpha_s(R)]^4 \log \left[C_A \alpha_s(R) e^{\gamma_E - 5/6} \right].$$

- They inherit the infrared problem of the static potential.

RS scheme

- It is defined from the pole mass by subtracting its leading asymptotic behavior [Antonio Pineda, 2001].

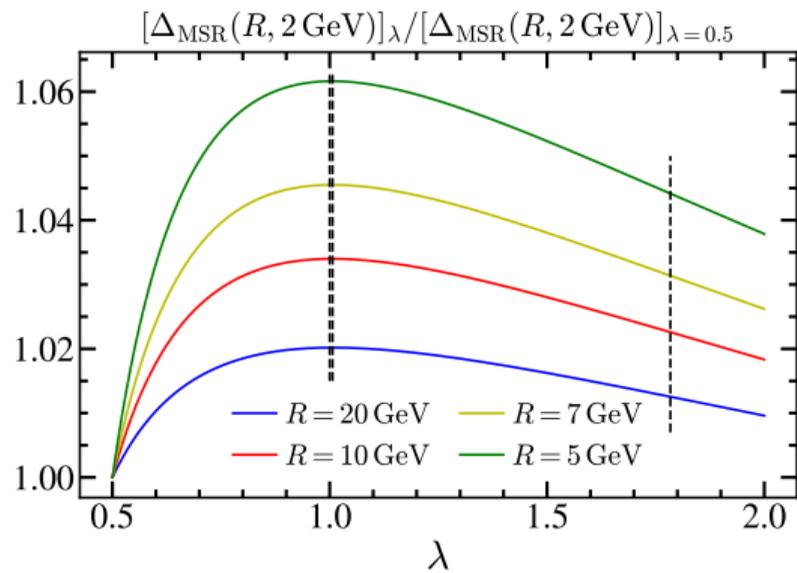
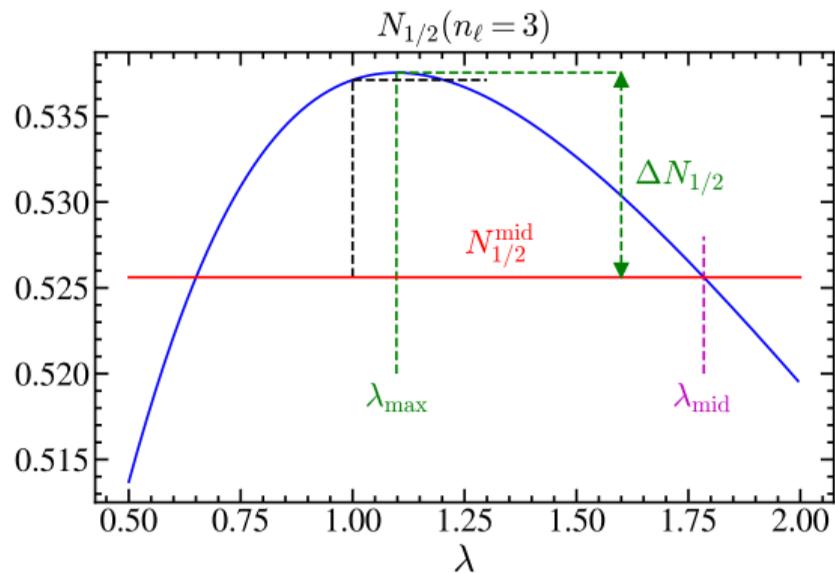
$$m_Q^{\text{pole}} - m_Q^{\text{RS}}(R) = \frac{2\pi}{\beta_0} RN_{1/2} \sum_{n=1}^{\infty} \left[\frac{\beta_0 \alpha_s(R)}{2\pi} \right]^n \sum_{\ell=0}^{\infty} g_\ell \left(1 + \hat{b}_1\right)_{n-1-\ell},$$

- $N_{1/2}$ is the normalization of the leading renormalon

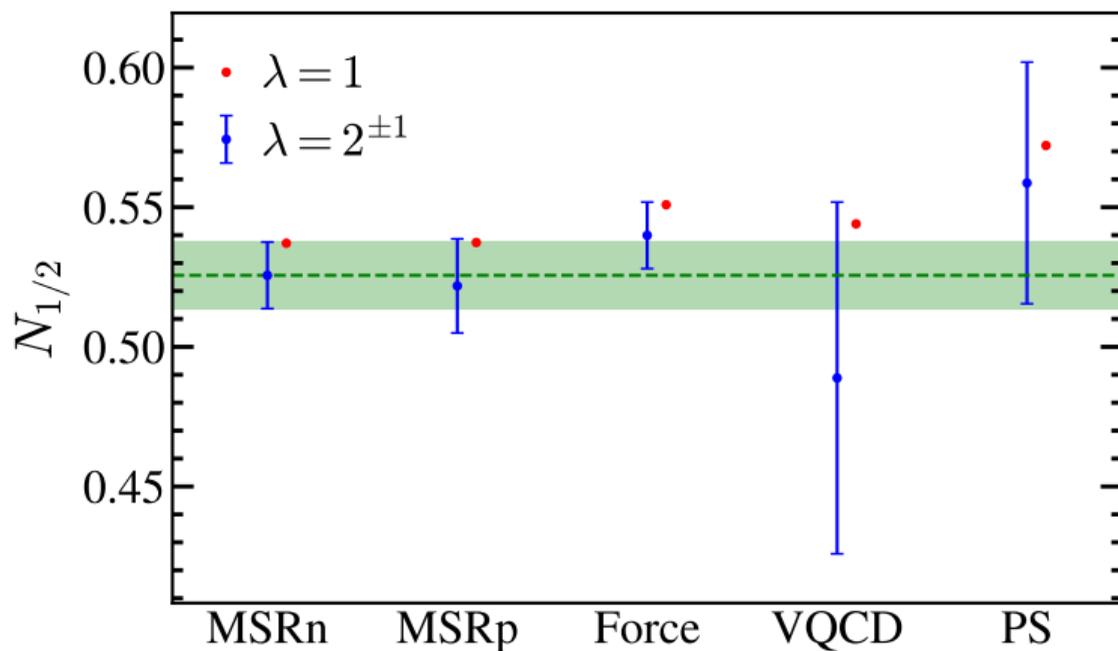
$$N_{1/2}^{(n)} = \frac{\beta_0}{2\pi} \sum_{k=0}^n \frac{S_k}{\left(1 + \hat{b}_1\right)_k} \quad S_k = \sum_{k=0}^j \tilde{\gamma}_k^R \sum_{i=0}^{j-k} (-1)^i \tilde{b}_i^N \tilde{g}_{j-i-k}^N,$$

- $N_{1/2}$ depends on λ , it reshuffles higher perturbative orders. We vary it to estimate $N_{1/2}$ (similar to scale variation).
- λ is also used to estimate R-evolution uncertainty.

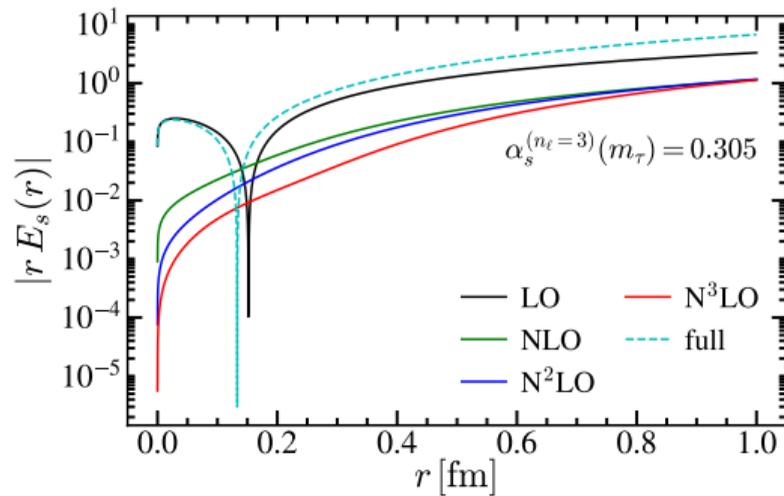
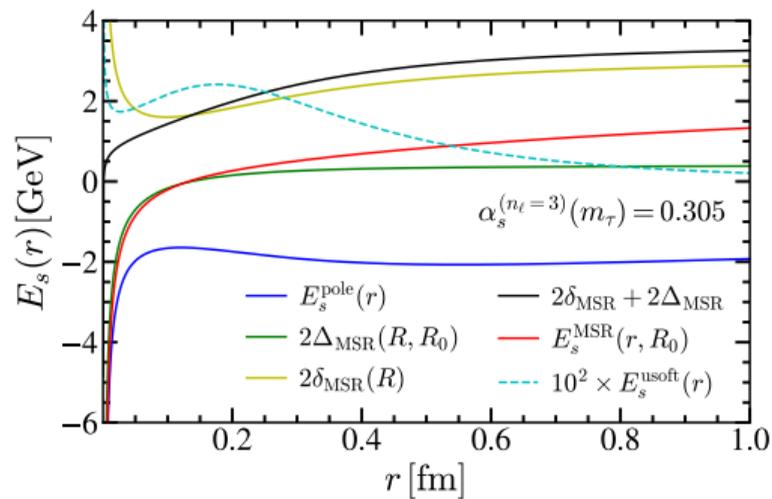
$N_{1/2}$



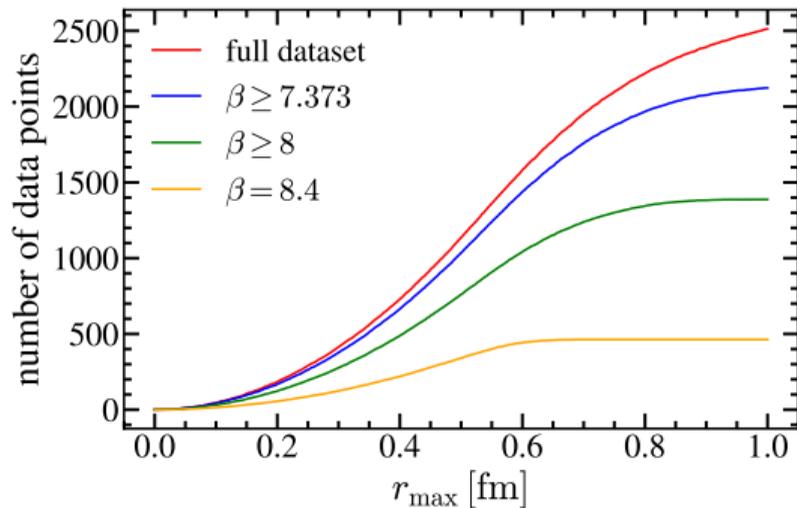
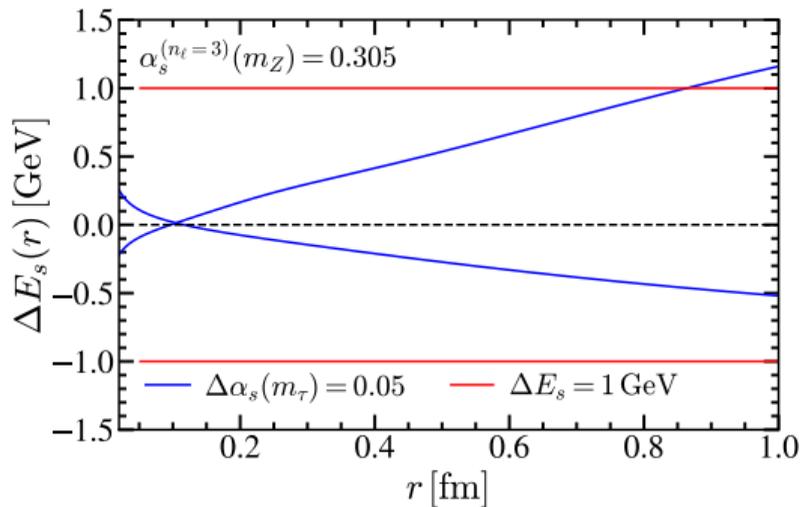
Renormalon subtraction



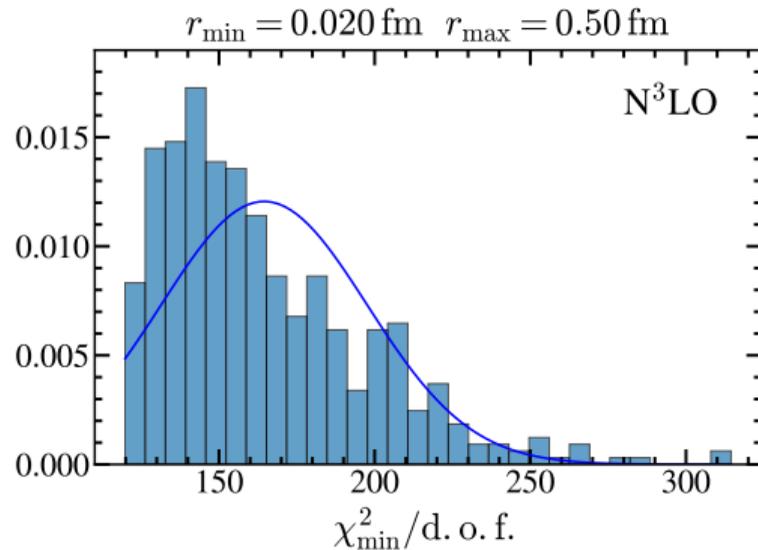
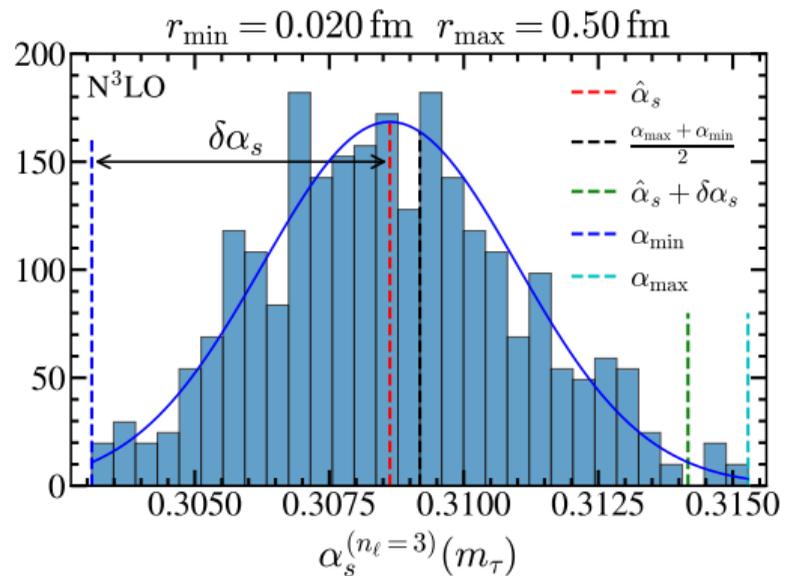
Perturbative analysis



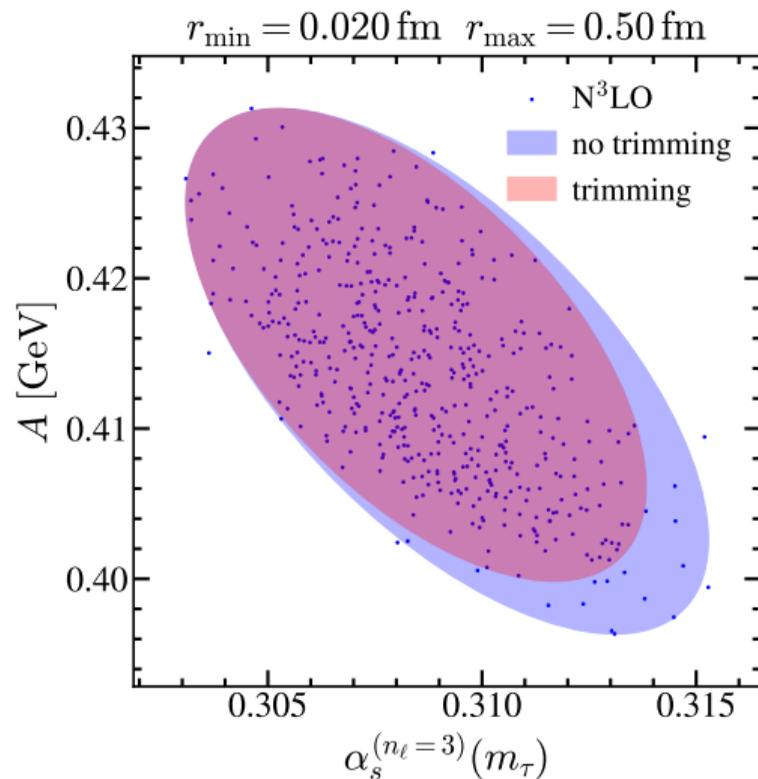
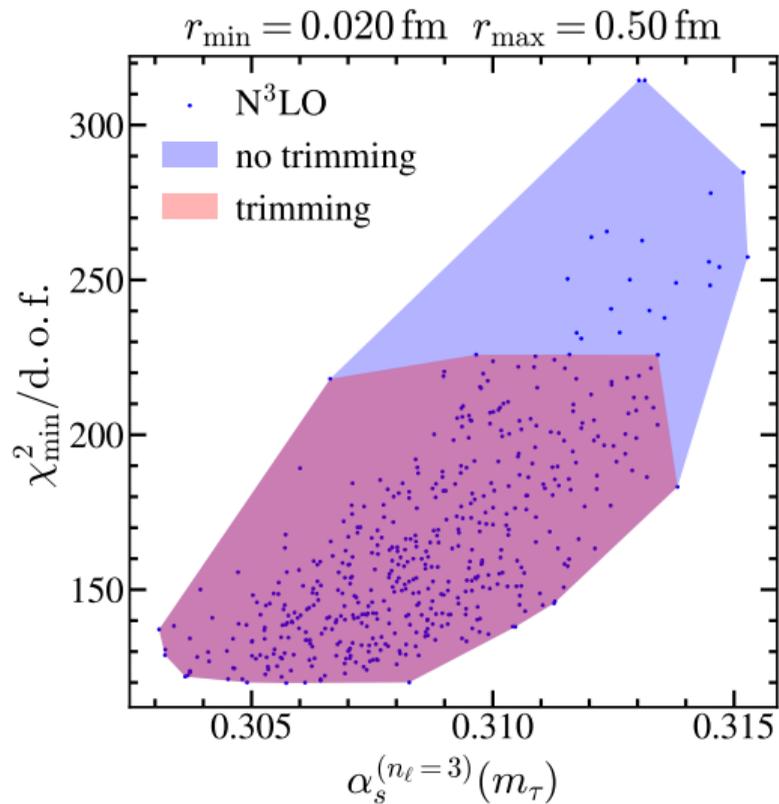
Auxiliar figures



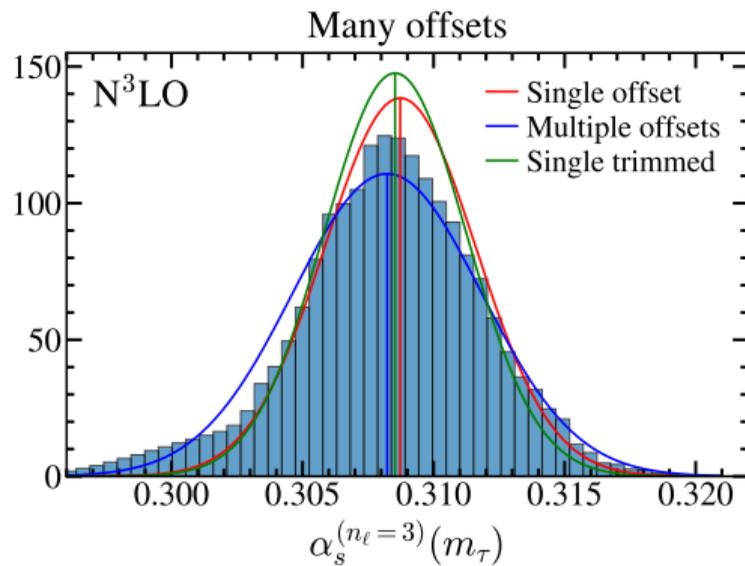
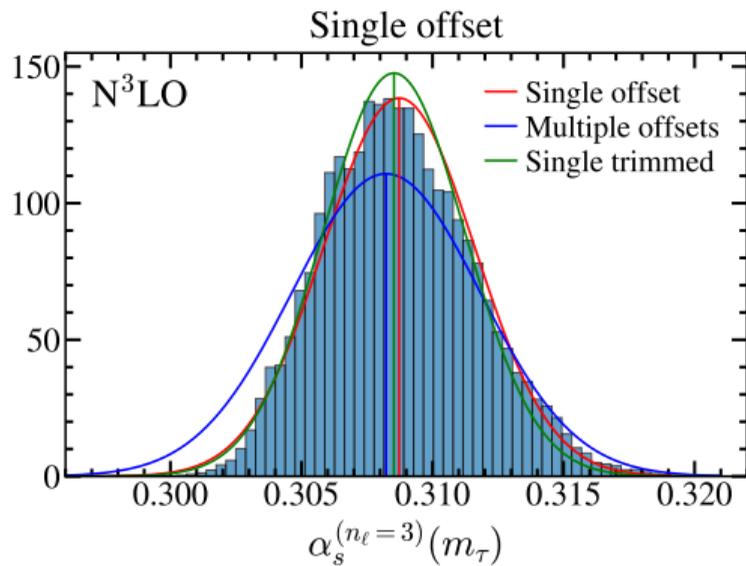
Random scan histograms



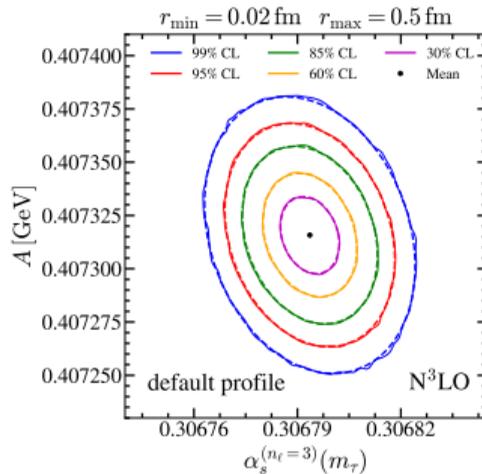
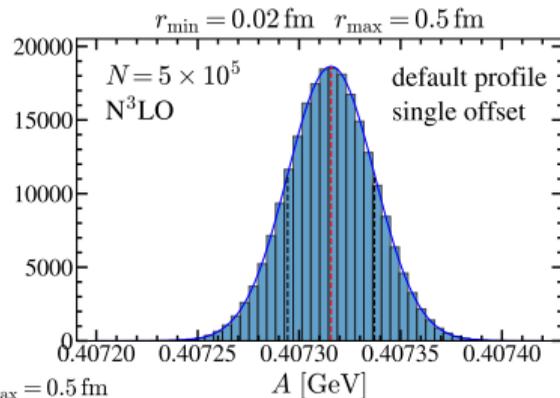
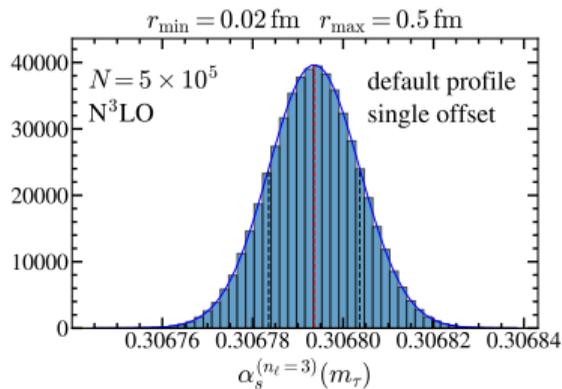
χ^2 trimming



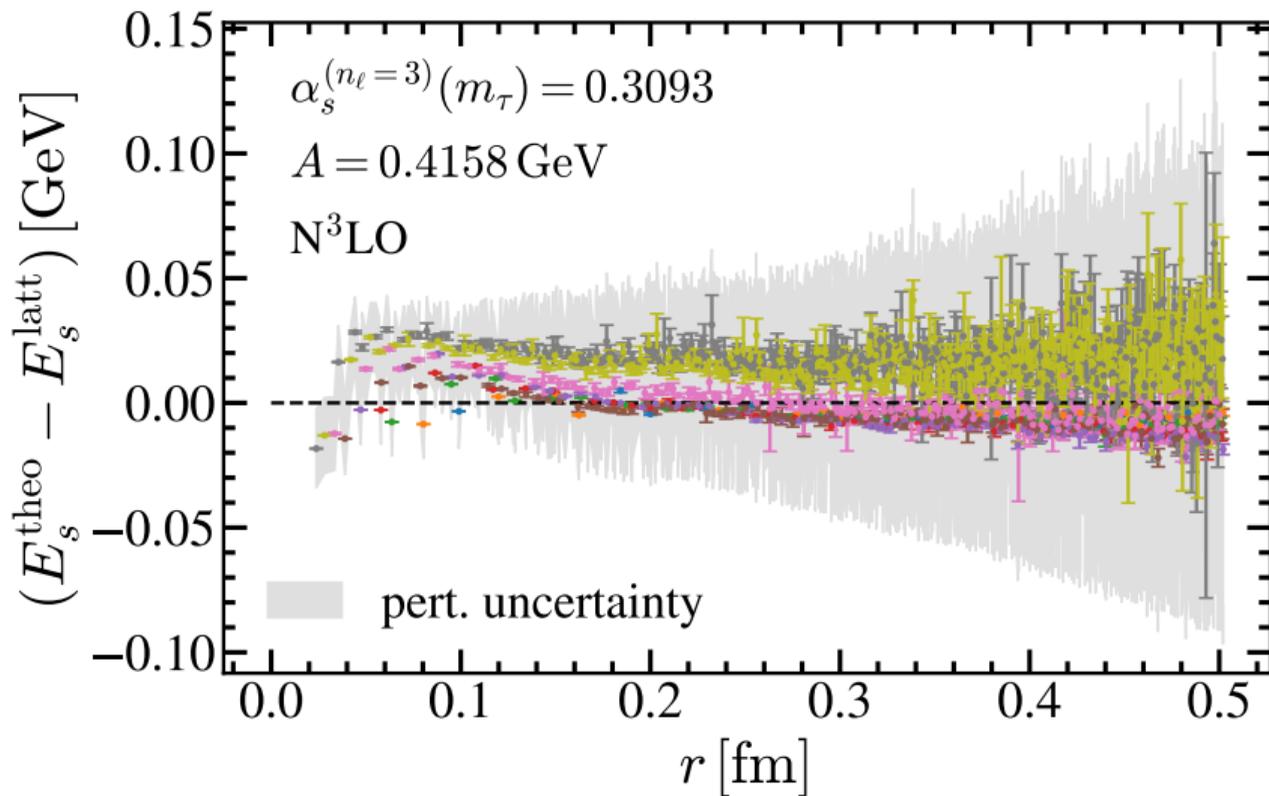
Fits histograms



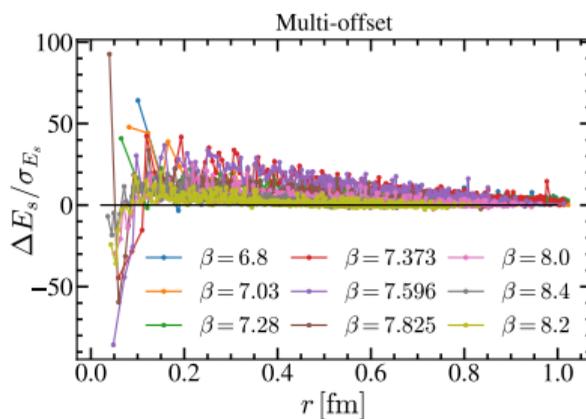
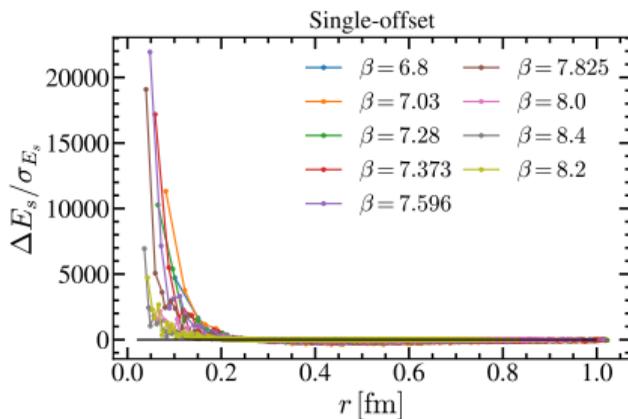
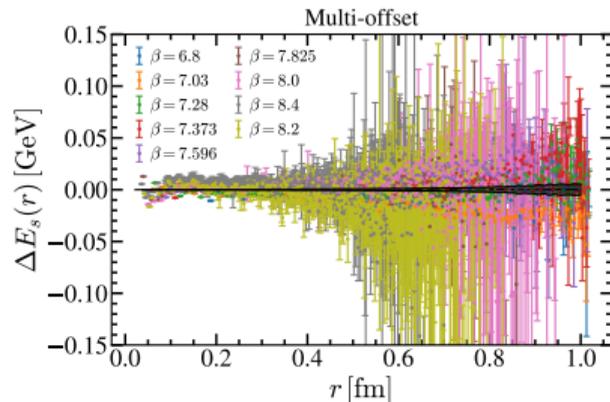
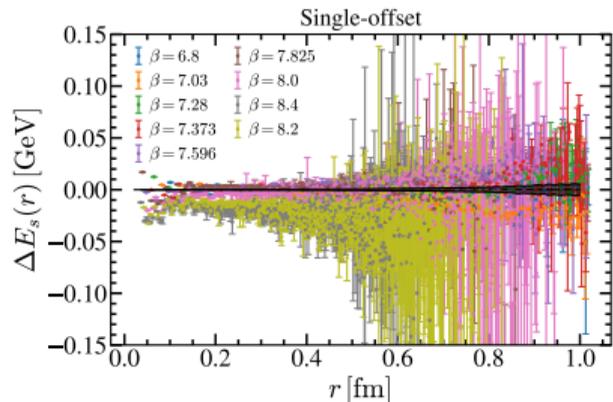
Replica method



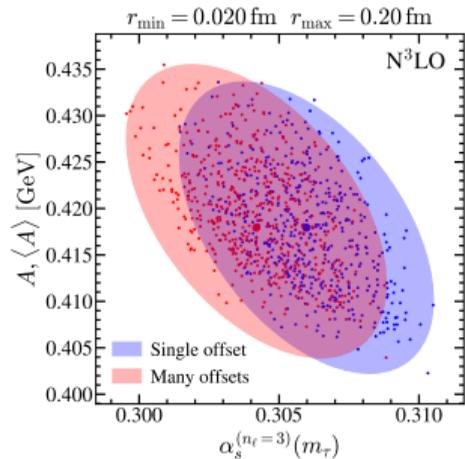
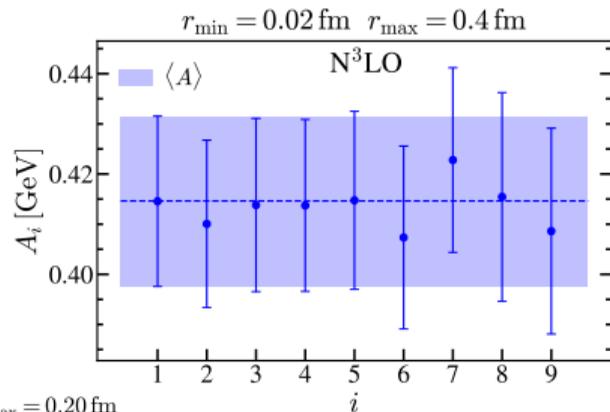
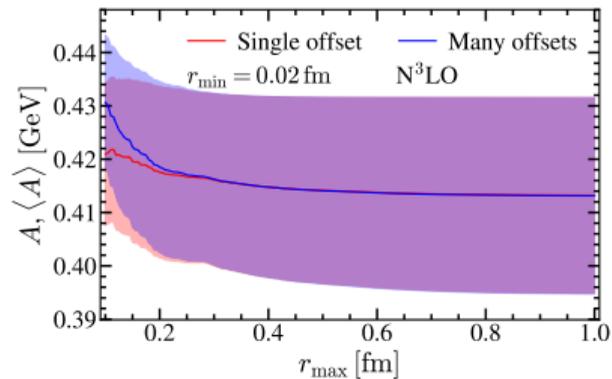
Lattice data



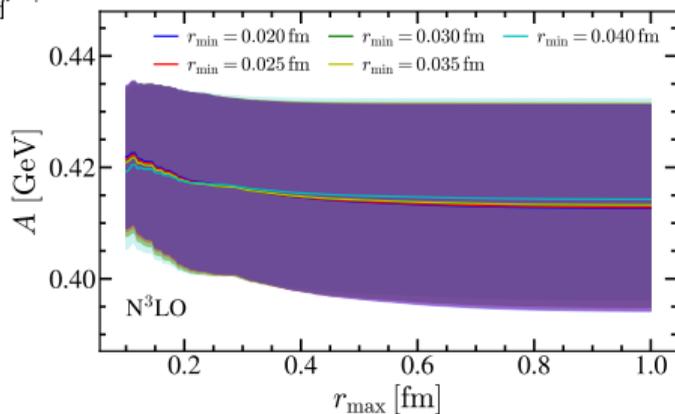
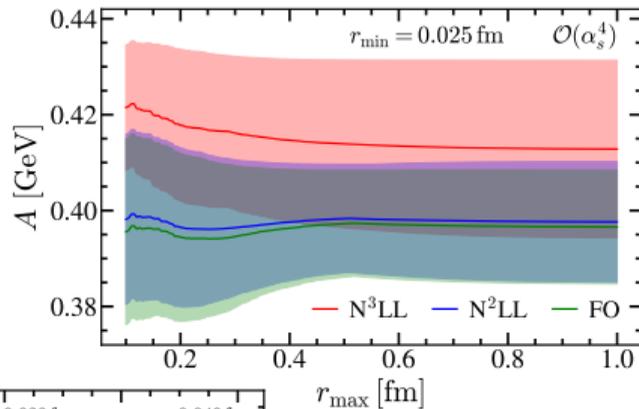
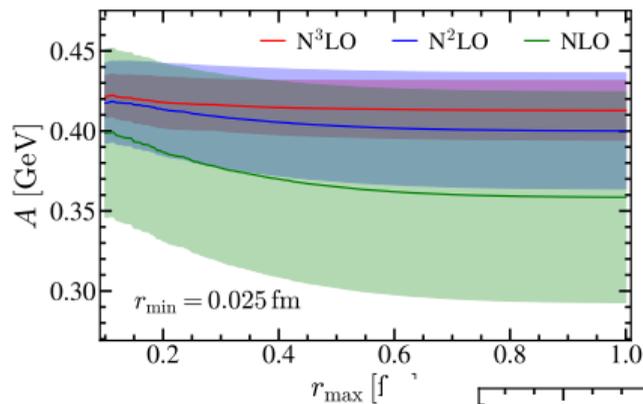
Lattice data



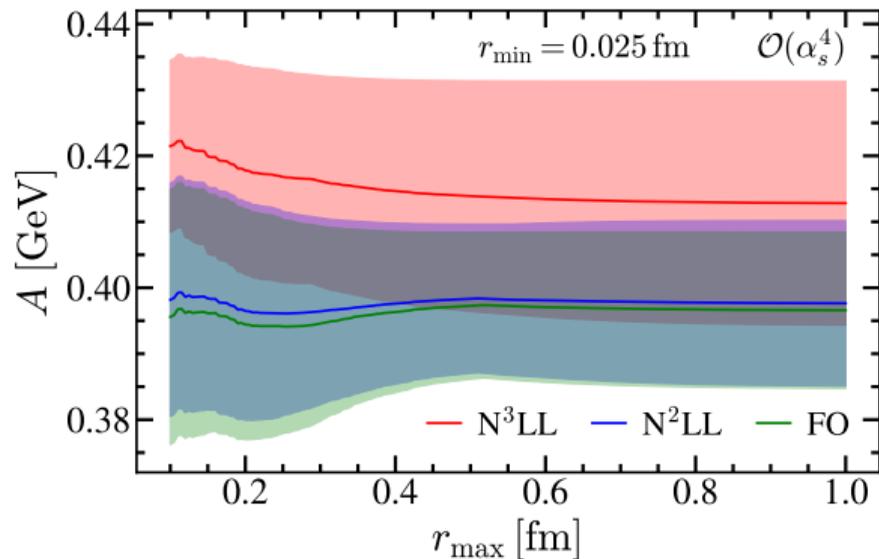
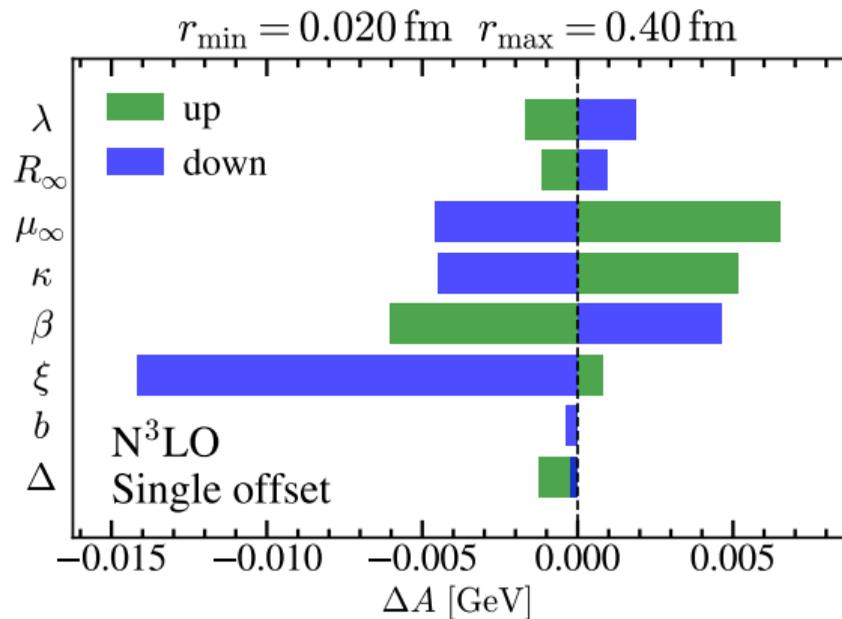
Offset analysis



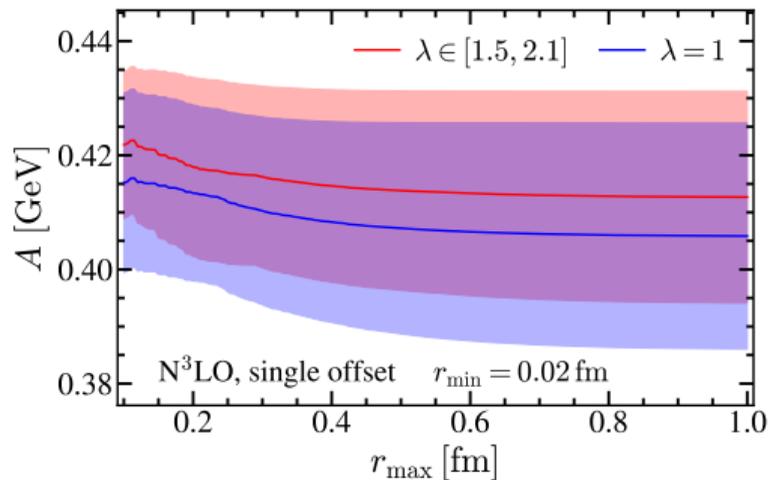
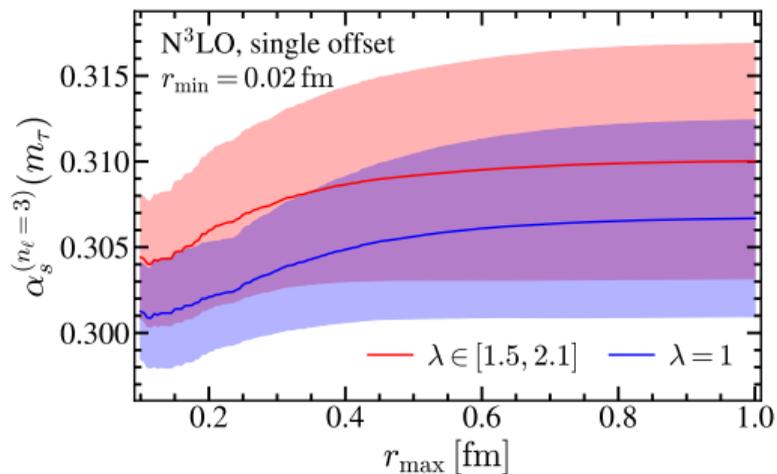
Offset analysis



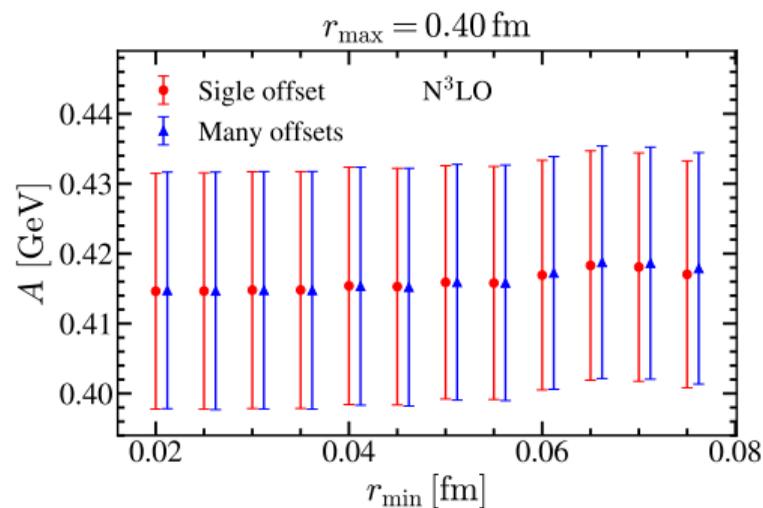
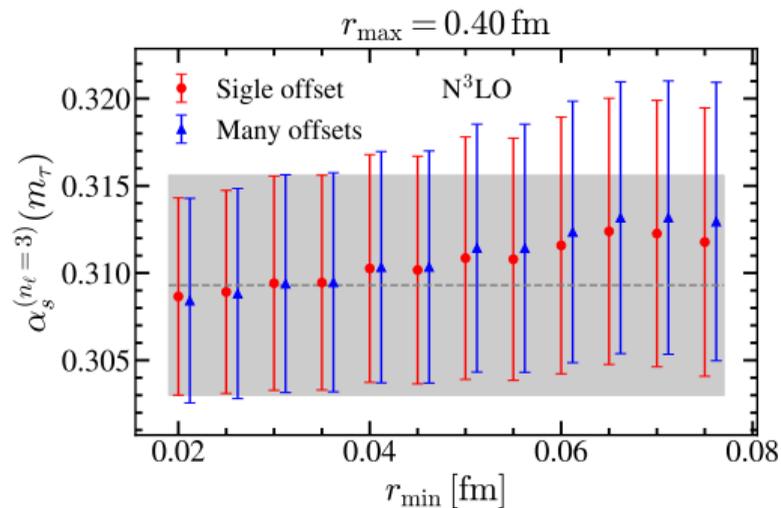
Offset analysis



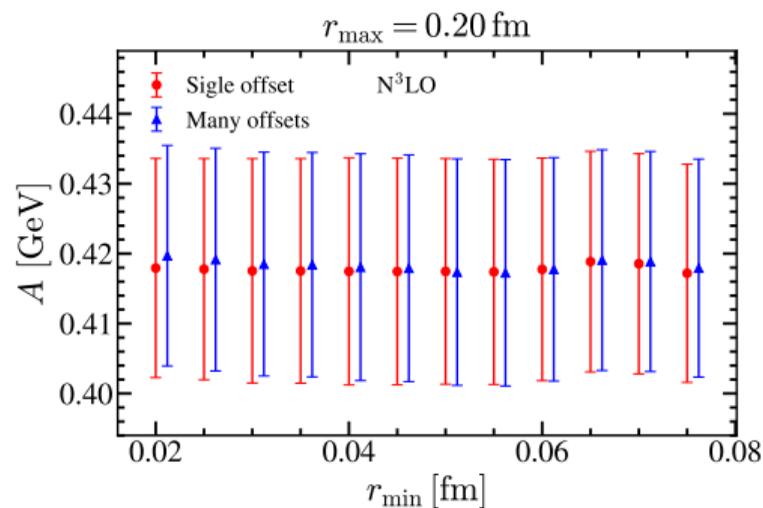
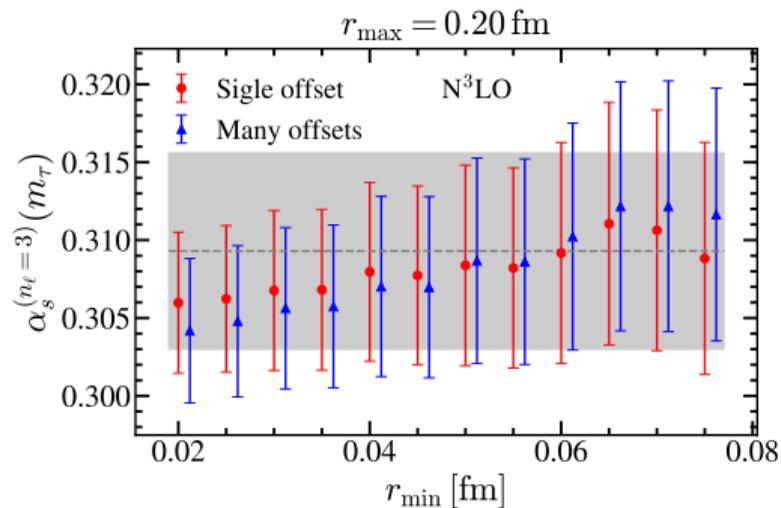
λ variation



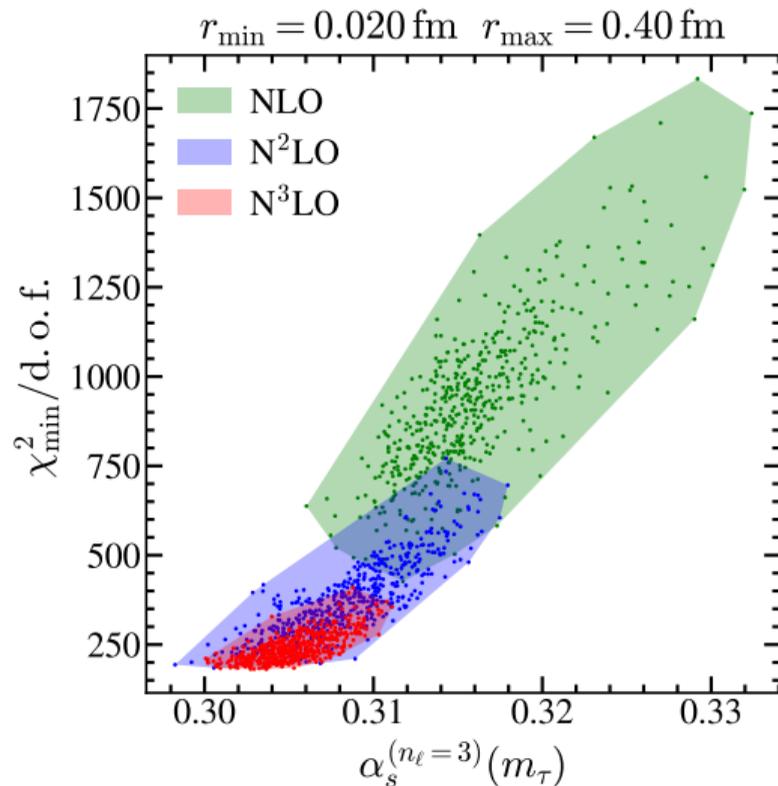
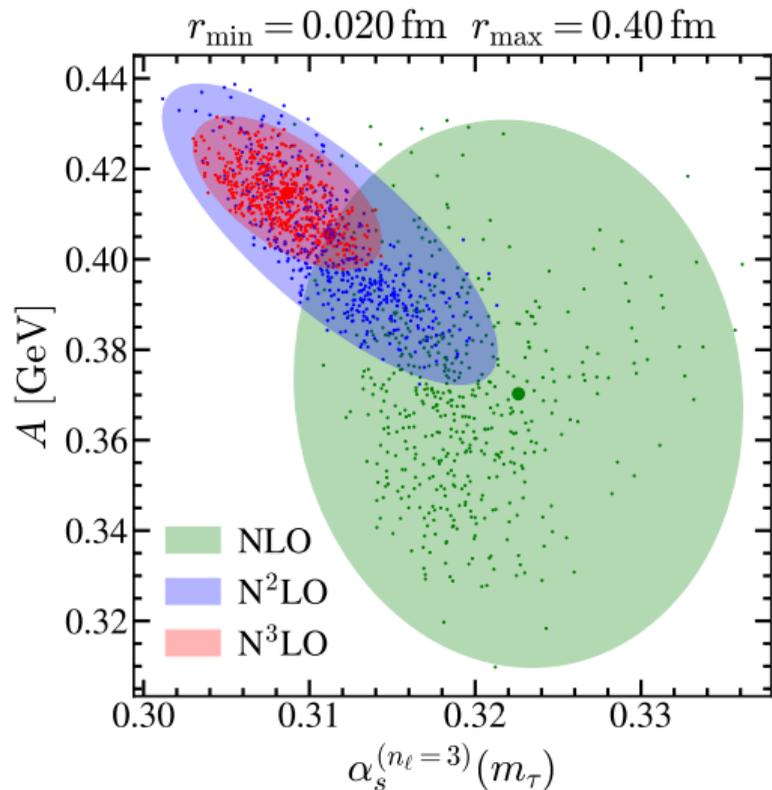
Further dependence check on r_{min}



Further dependence check on r_{min}



Scattered fit points



Additional fits

	full	$\beta \geq 7.737$	$\beta \geq 8$	$\beta = 8.4$
$\alpha_s^{(n_\ell=3)}(m_\tau)$	0.3093 ± 0.0063	0.3089 ± 0.0061	0.3089 ± 0.0070	0.3080 ± 0.0072
$\alpha_s^{(n_\ell=5)}(m_Z)$	0.1170 ± 0.0009	0.1170 ± 0.0009	0.1170 ± 0.0010	0.1169 ± 0.0010

	$\mu_c = \bar{m}_c$	$\lambda = 1$	FO	many offsets
$\alpha_s^{(n_\ell=3)}(m_\tau)$	0.3093 ± 0.0063	0.3055 ± 0.0049	0.3108 ± 0.0057	0.3092 ± 0.0064
$\alpha_s^{(n_\ell=5)}(m_Z)$	0.1174 ± 0.0009	0.1165 ± 0.0007	0.1172 ± 0.0008	0.1170 ± 0.0009