



Towards EPPS21 nuclear PDFs

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DIS2021

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Short intro to EPPS framework

- Define nuclear PDFs in terms of

$$\underset{\text{bound-proton PDF}}{f_i^{p/A}}(x, Q^2) = \overset{\text{nuclear modification}}{R_i^{p/A}}(x, Q^2) \underset{\text{free-proton PDF}}{f_i^p}(x, Q^2)$$

- PDFs of the full nucleus are then constructed with

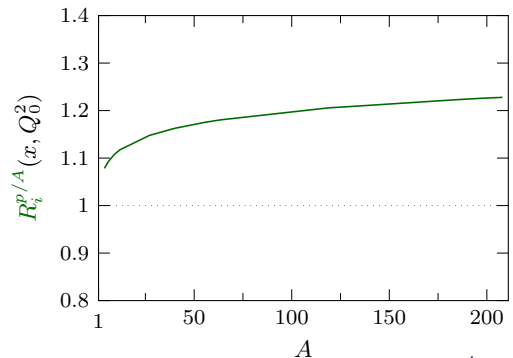
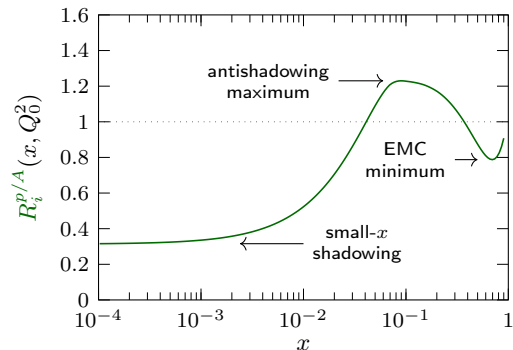
$$f_i^A(x, Q^2) = Z f_i^{p/A}(x, Q^2) + N f_i^{n/A}(x, Q^2),$$

and assuming $f_i^{p/A} \xleftrightarrow{\text{isospin}} f_j^{n/A}$

- Parametrize the x and A dependence of $R_i^{p/A}(x, Q_0^2)$ at $Q_0 = m_{\text{charm}} = 1.3 \text{ GeV}$

- Use a phenomenologically motivated piecewise function in x
- Use a power-law type function in A

We are currently exploring the best fit functions for our next 'EPPS21' fit



Comparison with other NLO nPDF global fits

	EPPS16	nNNPDF2.0	nCTEQ15WZ	nCTEQ15HIX	EPPS21 <i>prelim.</i>
IA NC DIS	✓	✓	✓	✓	✓
+ JLab NC DIS				✓	✓ <i>new!</i>
ν A CC DIS	✓	✓			✓
pA DY	✓		✓	✓	✓
π A DY	✓				✓
RHIC dAu/pp π^0	✓		✓		✓
LHC pPb dijet R_{FB}	✓				
→ dijet R_{pPb}					✓ <i>new!</i>
LHC pPb D^0					✓ <i>new!</i>
LHC pPb W,Z Run 1	✓	✓	✓		✓
+ Run 2 pPb W		✓	✓		✓ <i>new!</i>
Q cut in DIS	1.3 GeV	1.87 GeV	2 GeV	1.3 GeV	1.3 GeV
Data points	1811	1467	828	1564	2023 <i>prelim.</i>
Free parameters	20	256	19	19	24 <i>prelim.</i>
Error analysis	Hessian	Monte Carlo	Hessian	Hessian	Hessian
Error tolerance $\Delta\chi^2$	52	N/A	35	35	35 <i>prelim.</i>
Free-proton PDFs	CT14	NNPDF3.1	~CTEQ6M	~CTEQ6M	CT18A <i>prelim.</i>
HQ treatment	S-ACOT	FONLL	S-ACOT	S-ACOT	S-ACOT
Indep. flavours	6	6	5	4	6
Reference	EPJC 77, 163	JHEP 09, 183	EPJC 80, 968	arXiv:2012.11566	TBA

Need to mitigate free-proton PDF uncertainty

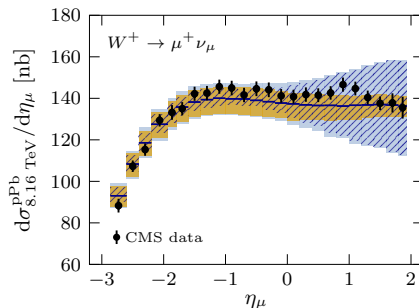
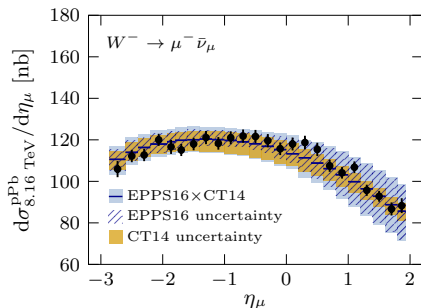
Absolute cross sections carry large proton-PDF uncertainty!

Should not be neglected when fitting the nPDFs!

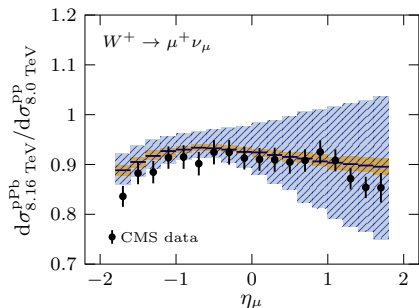
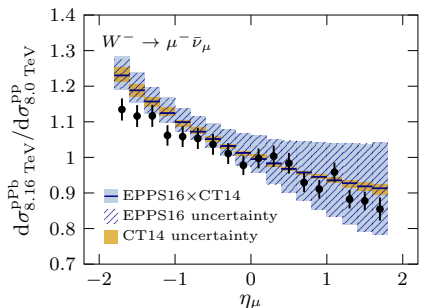
Wherever possible, we use nuclear modification ratios to cancel the free-proton PDF uncertainty

For Ws at 8.16 TeV, we formulate a mixed-energy nuclear modification ratio

$$R_{\text{pPb}} = \frac{d\sigma_{8.16 \text{ TeV}}^{\text{pPb}}/d\eta_{\mu}}{d\sigma_{8.0 \text{ TeV}}^{\text{pp}}/d\eta_{\mu}}$$

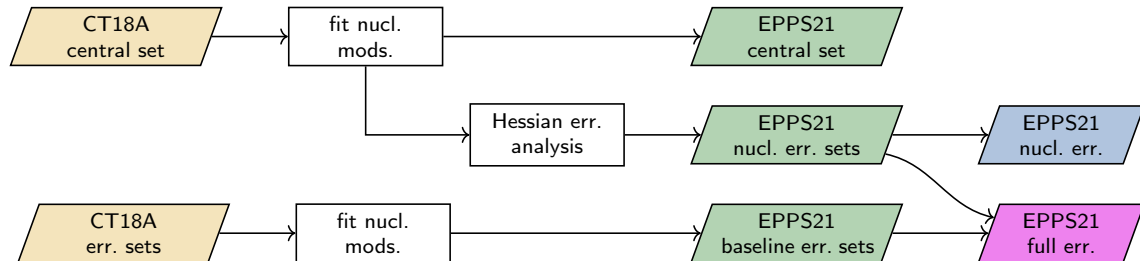


↓ Cancel proton-PDF uncertainty ↓

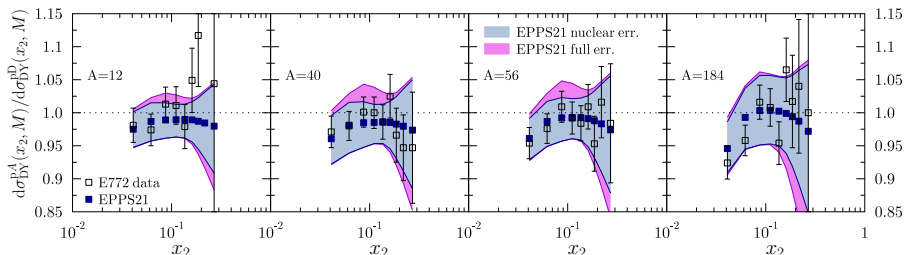


Charting the baseline proton-PDF uncertainty *new!*

We study baseline-PDF sensitivity by fitting nuclear modifications separately for each CT18A error set



Baseline error mostly subdominant in the observables we fit, but shows up e.g. in the fixed-target DY



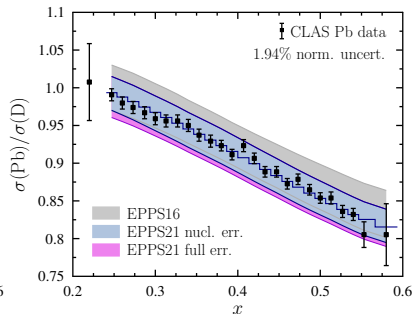
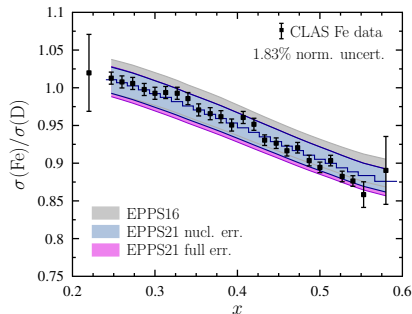
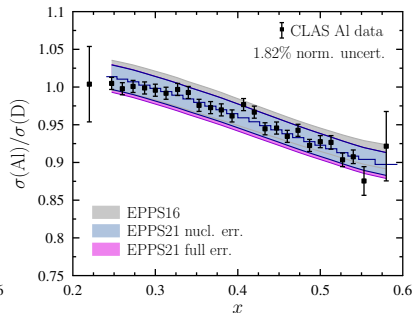
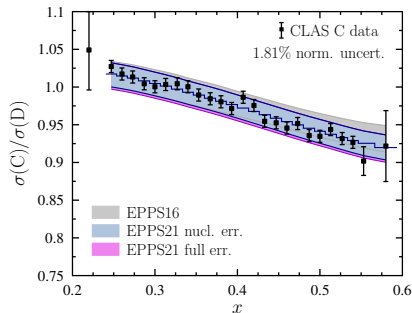
Excellent fit!

Results in line with
the reweighting study

[Paukkunen & Zurita,
Eur.Phys.J.C 80 (2020) 381]

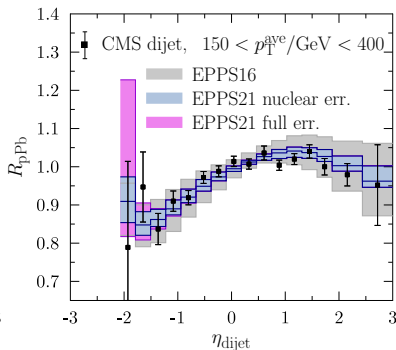
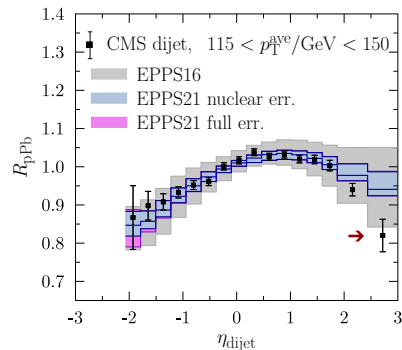
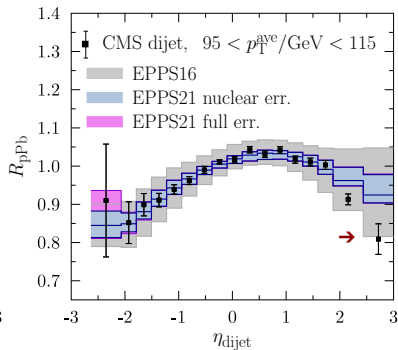
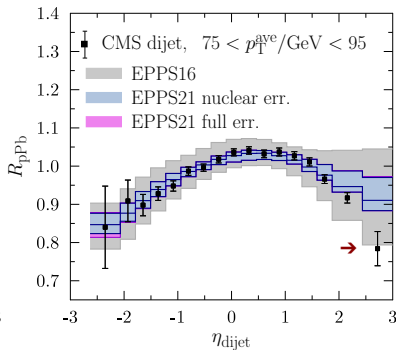
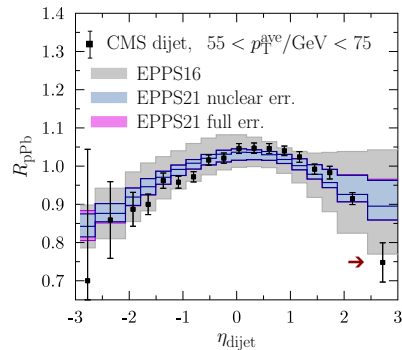
We take into account
the leading target-mass
corrections

No sign of isospin-dependence
in the bound-proton
nuclear modifications $R_i^{p/A}$



Dijets at 5.02 TeV *new!*

data from: [CMS Collaboration, Phys.Rev.Lett. 121 (2018) 062002]



Results in line with
the reweighting study

[Eskola, PP & Paukkunen,
Eur.Phys.J.C 79 (2019) 511]

Still finding it difficult to fit
the forwardmost data points

→ currently excluded
from the fit

Excellent fit!

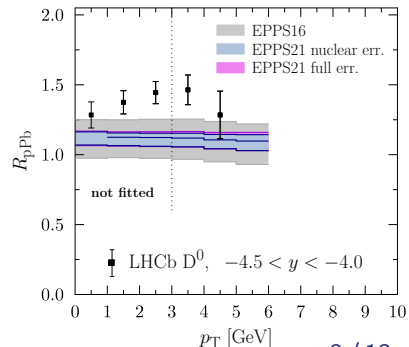
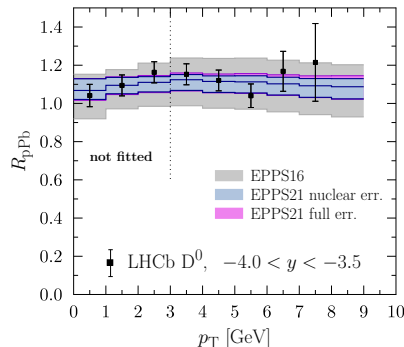
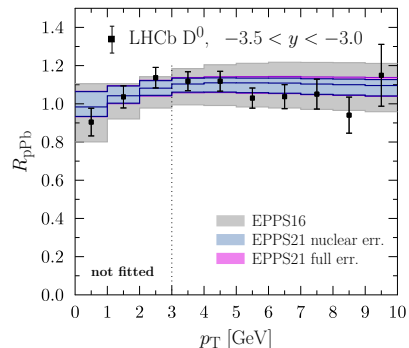
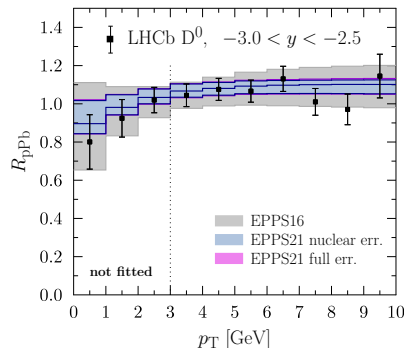
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[Eskola, Helenius, PP & Paukkunen,
JHEP 05 (2020) 037]

Using the NLO pQCD
S-ACOT- m_T GM-VFNS

[Helenius & Paukkunen,
JHEP 05 (2018) 196]

Using a $p_T > 3$ GeV cut
to reduce theoretical
uncertainties



Excellent fit!

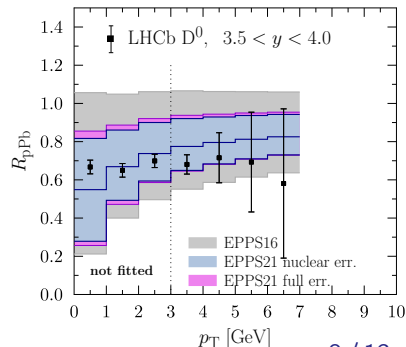
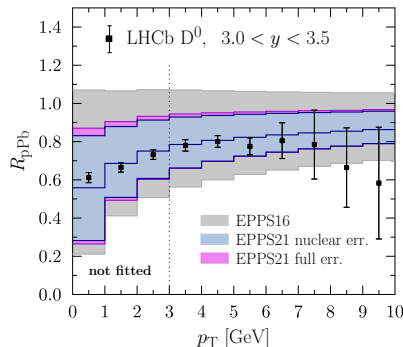
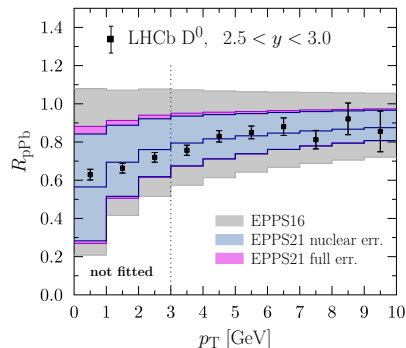
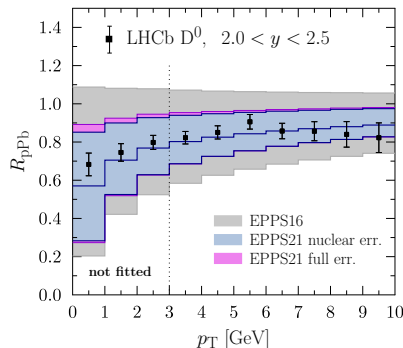
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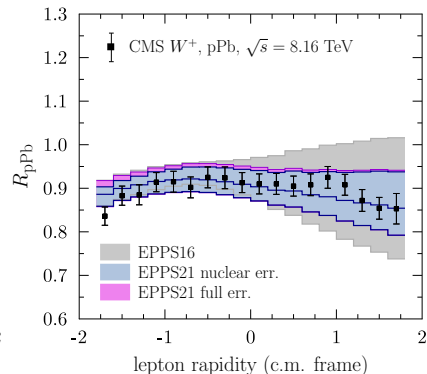
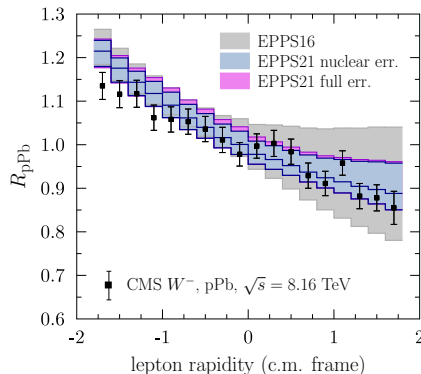


Excellent fit!

Using the mixed-energy
nuclear modification ratio

$$R_{\text{pPb}} = \frac{d\sigma_{8.16 \text{ TeV}}^{\text{pPb}}/d\eta_\mu}{d\sigma_{8.0 \text{ TeV}}^{\text{pp}}/d\eta_\mu}$$

to cancel the free-proton
PDF uncertainty



Fully consistent with the dijets and D^0 s

- Important check on the nPDF universality & factorization

These data do not appear to give additional flavour-separation constraints
on top of those we had already in EPPS16

- Looking forward to increased precision at LHC Run 3

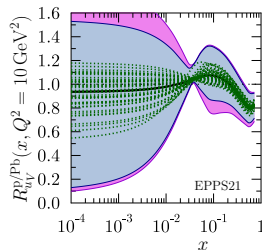
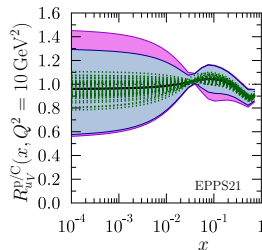
Fit results – valence

Bound-proton modifications *Prelim.*

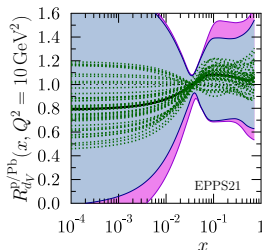
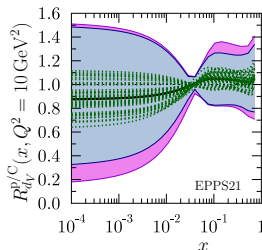
Carbon

Lead

u_V



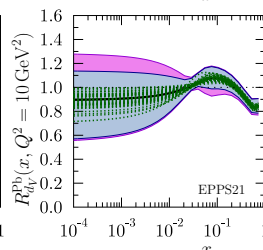
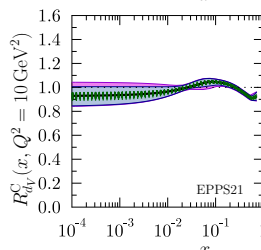
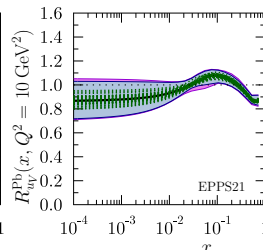
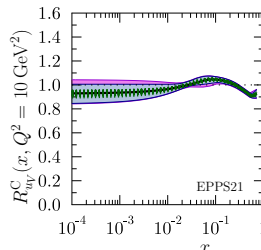
d_V



Full-nucleus modifications *Prelim.*

Carbon

Lead



$$R_i^{p/A} = \frac{f_i^{p/A}}{f_i^p}$$

$$R_i^A = \frac{Z f_i^{p/A} + N f_i^{n/A}}{Z f_i^p + N f_i^n}$$

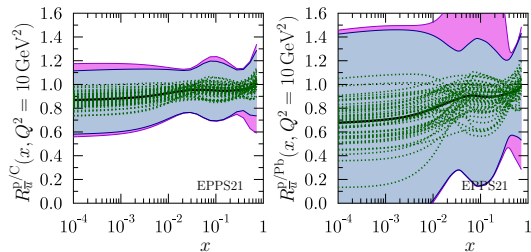
Fit results – sea

Bound-proton modifications *Prelim.*

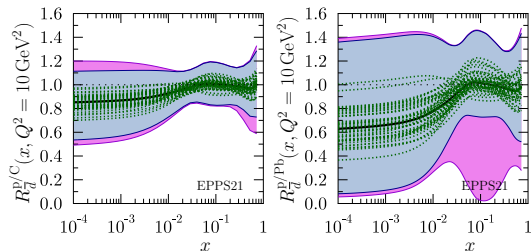
Carbon

Lead

$\bar{u}$



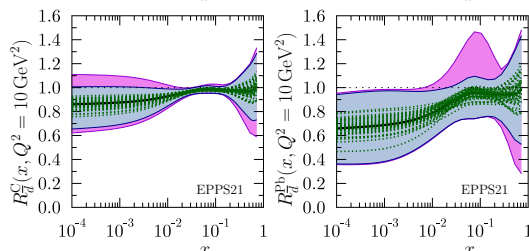
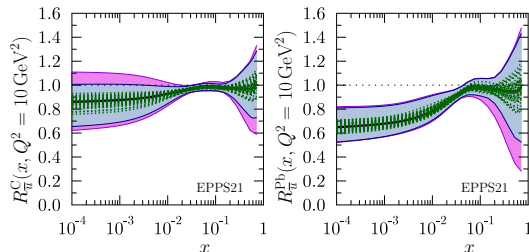
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Full-nucleus modifications *Prelim.*

Carbon

Lead



$$R_i^{p/A} = \frac{f_i^{p/A}}{f_i^p}$$

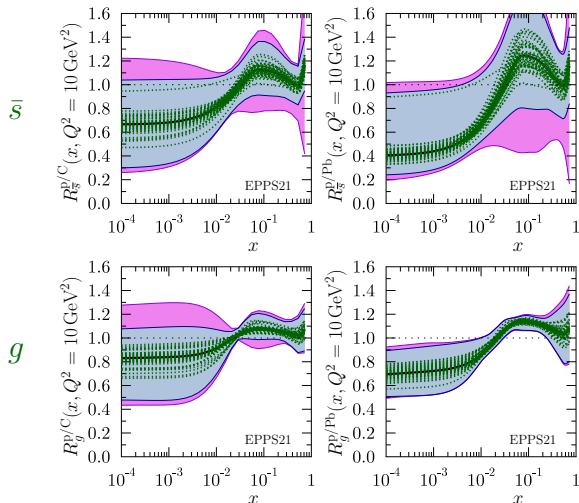
$$R_i^A = \frac{Z f_i^{p/A} + N f_i^{n/A}}{Z f_i^p + N f_i^n}$$

Fit results – strange and glue

Bound-proton modifications *Prelim.*

Carbon

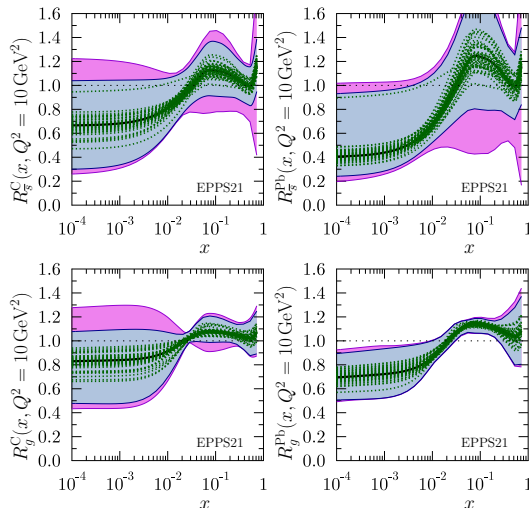
Lead



Full-nucleus modifications *Prelim.*

Carbon

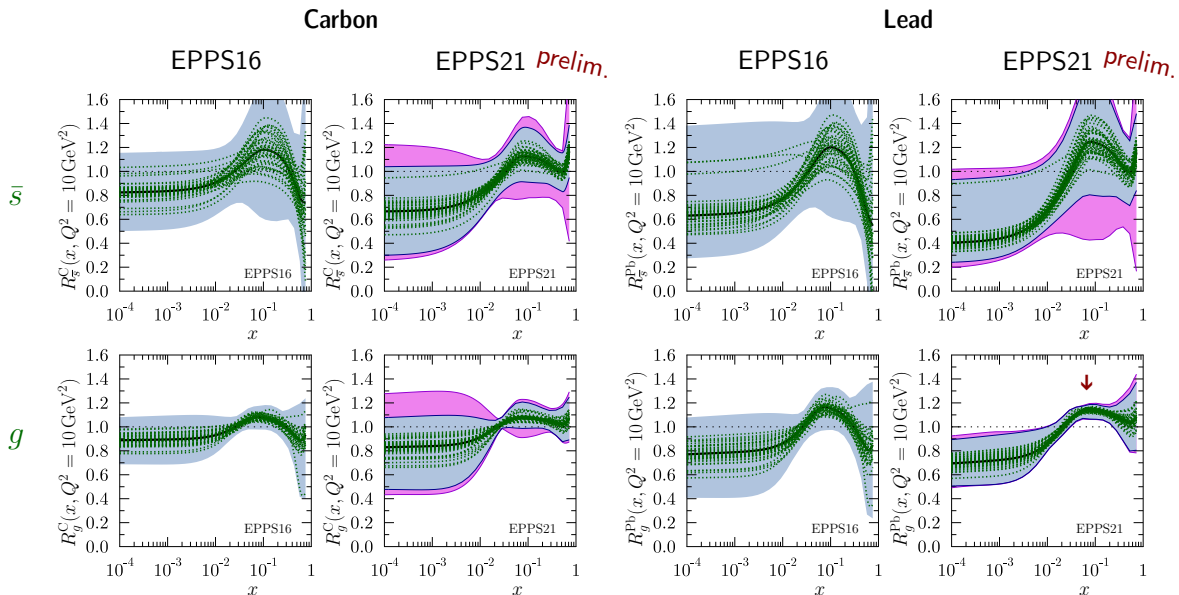
Lead



$$R_i^{p/A} = \frac{f_i^{p/A}}{f_i^p}$$

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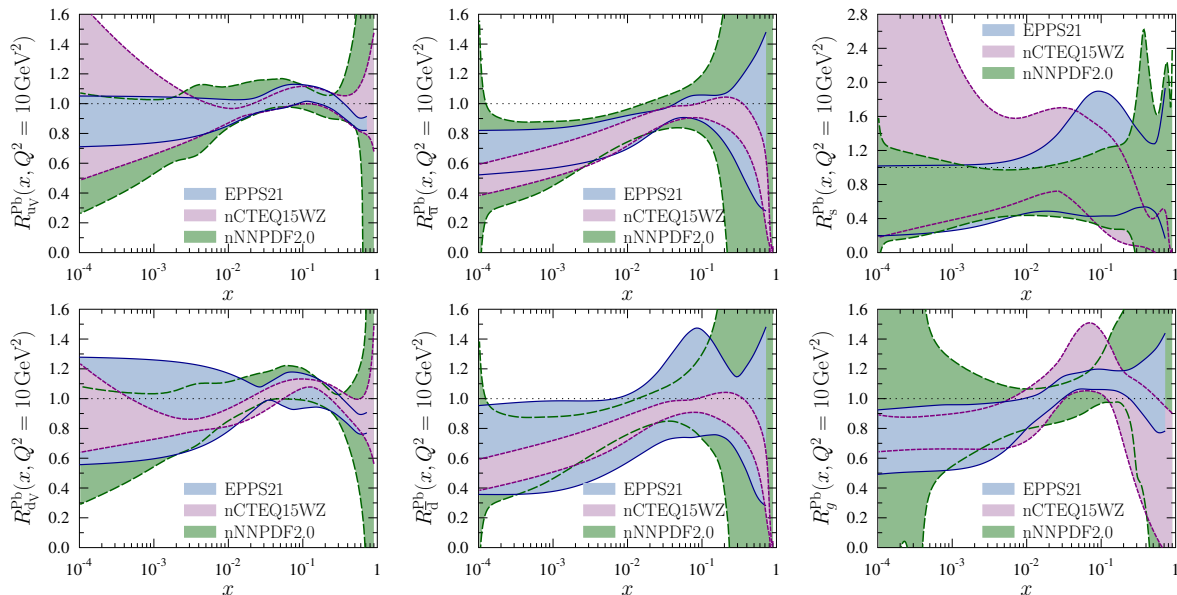
Comparison with EPPS16



- Better control over gluon (anti)shadowing → sub-10% level uncertainties at mid- x !
- Flavour separation (esp. strangeness) remains a difficult beast to tame

Comparison with nNNPDF2.0 and nCTEQ15WZ

prelim.



- All three consistent within uncertainties, but significant differences in the uncertainty estimates
- Best constrained gluons in the (prelim.) EPPS21 fit!

Summary

A new EPPS nPDF fit on its way...

- Obtained a good fit for multiple new nuclear data from LHC (W s, D^0 s and dijets) and JLab
- Strong new constraints on gluon modifications in lead $\rightarrow$ sub-10% level mid- x uncertainties!
- Flavour separation uncertainties remain large, esp. for the sea quarks

Study the baseline free-proton PDF sensitivity / error

- We use data as nuclear modification ratios to decorrelate nuclear-modification and free-proton d.o.f.s to best possible extent
- Residual baseline free-proton PDF uncertainty still relevant where direct data constraints are scarce (sea-quark flavour separation, light-nuclei gluons)

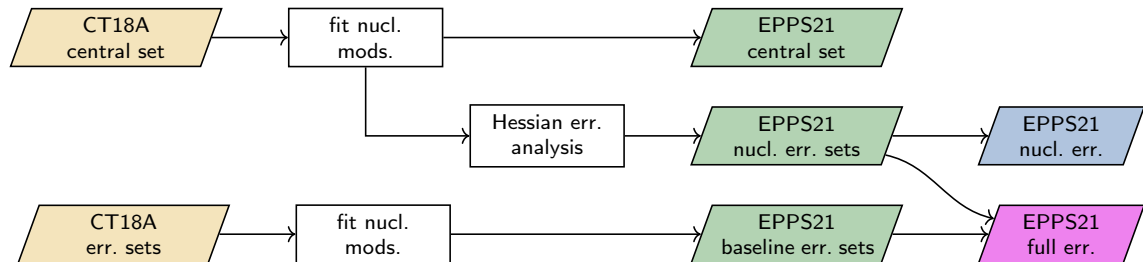
TODO:

- Still charting the possible parametrization choices

Backup

Charting the baseline proton-PDF uncertainty *new!*

We study baseline-PDF sensitivity by fitting nuclear modifications separately for each CT18A error set



The full error is calculated with:

$$\begin{aligned}
 (\delta X_{\text{full}}^{\pm})^2 &= (\delta X_{\text{nuc.}}^{\pm})^2 + \underbrace{(\delta X_{\text{baseline}}^{\pm})^2}_{N_{\text{baseline param.}}} \\
 &= \sum_k \left(\max_{\min} \left\{ \underbrace{X[S_k^+] - X[S_0]}_{\text{calculate with } f_i^{p/A}[S_k^+] = \underbrace{R_i^{p/A}[S_k^+]}_{\text{fitted using the } f_i^p[S_k^+] \text{ proton PDFs}}}, X[S_k^-] - X[S_0], 0 \right\} \right)^2
 \end{aligned}$$

Comparison with EPPS16 – valence

Carbon

Lead

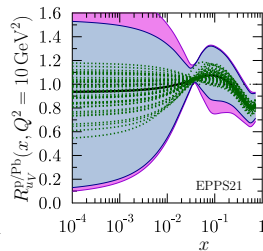
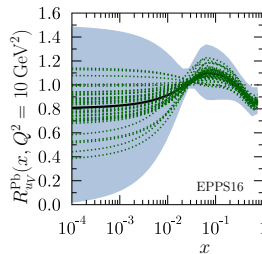
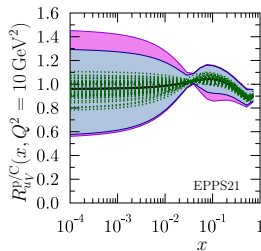
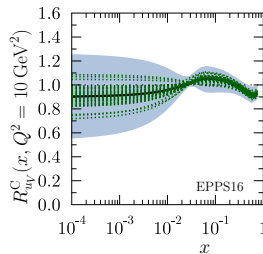
EPPS16

EPPS21 *Prelim.*

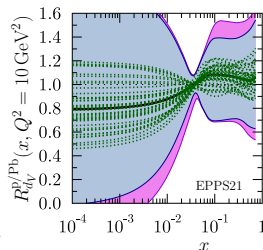
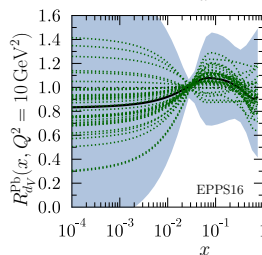
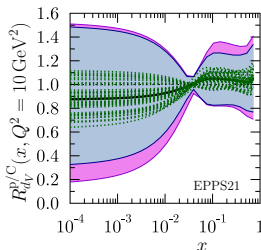
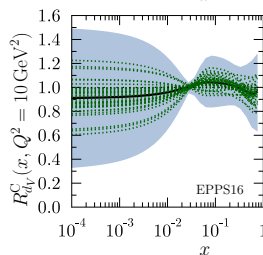
EPPS16

EPPS21 *Prelim.*

u_V



d_V



Comparison with EPPS16 – sea

Carbon

Lead

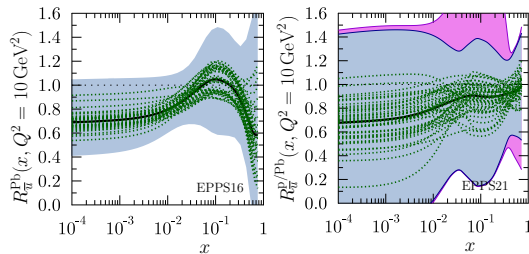
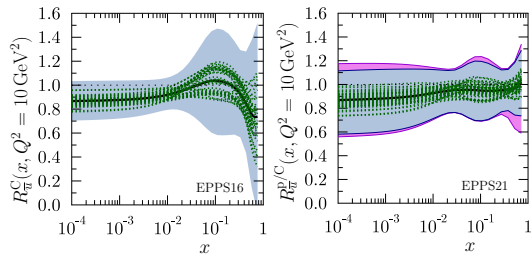
EPPS16

EPPS21 *Prelim.*

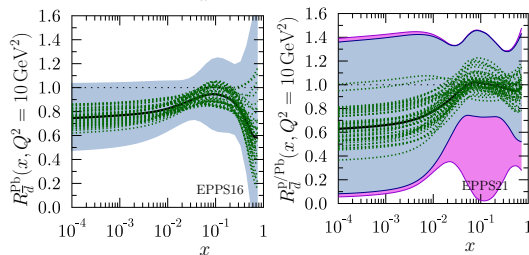
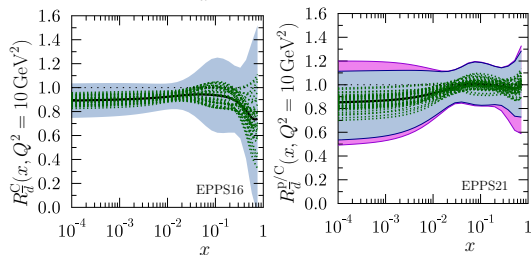
EPPS16

EPPS21 *Prelim.*

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Comparison with EPPS16 – strange and glue

Carbon

Lead

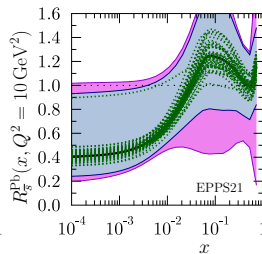
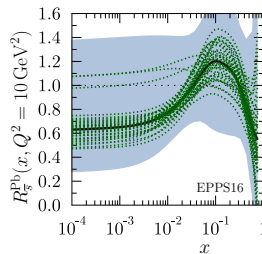
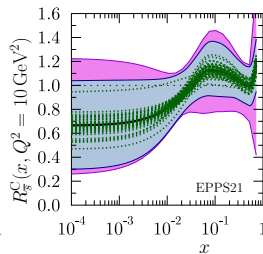
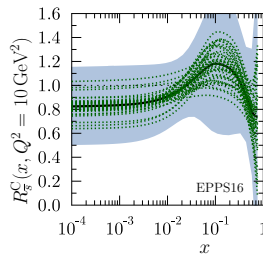
EPPS16

EPPS21 *Prelim.*

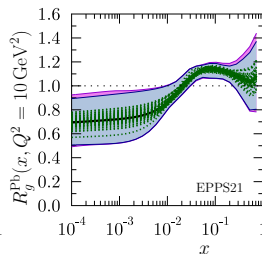
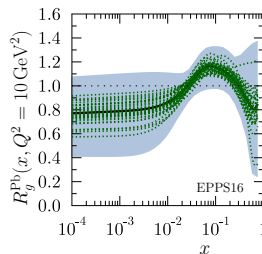
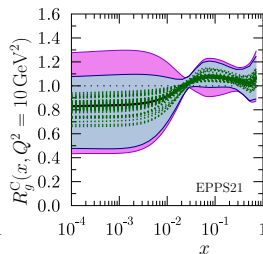
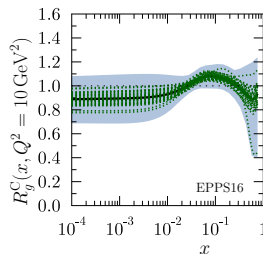
EPPS16

EPPS21 *Prelim.*

$\bar{s}$



g



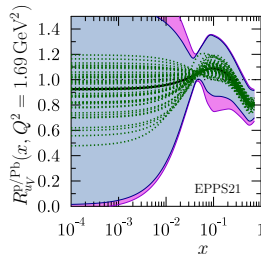
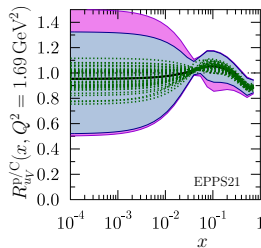
Fit results at parametrization scale – valence

Bound-proton modifications *Prelim.*

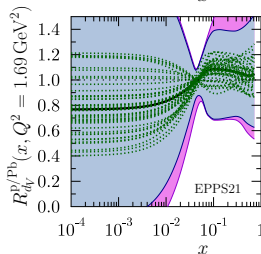
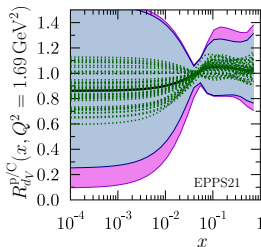
Carbon

Lead

u_V



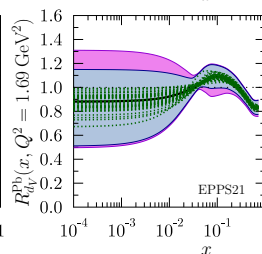
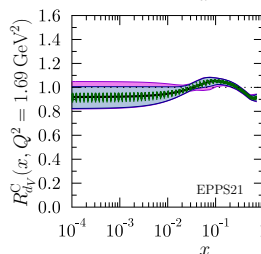
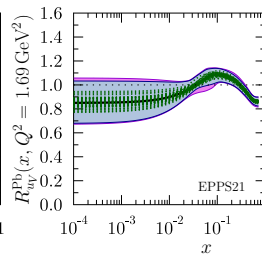
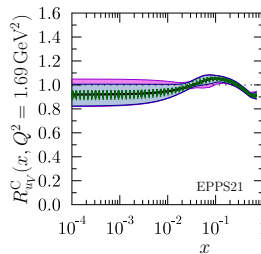
d_V



Full-nucleus modifications *Prelim.*

Carbon

Lead



$$R_i^{p/A} = \frac{f_i^{p/A}}{f_i^p}$$

$$R_i^A = \frac{Z f_i^{p/A} + N f_i^{n/A}}{Z f_i^p + N f_i^n}$$

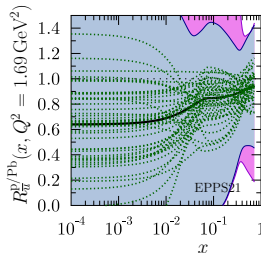
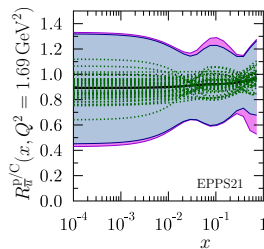
Fit results at parametrization scale – sea

Bound-proton modifications *Prelim.*

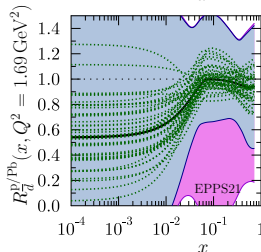
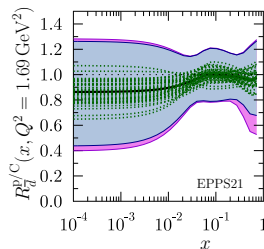
Carbon

Lead

$\bar{u}$



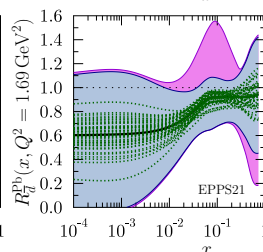
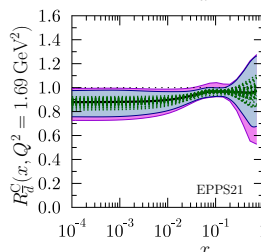
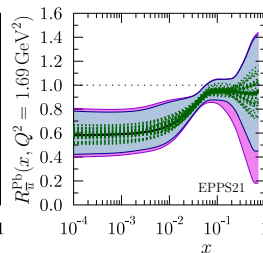
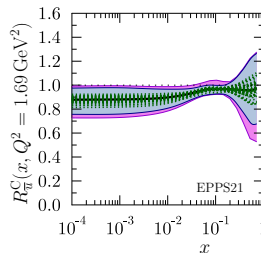
$\bar{d}$



Full-nucleus modifications *Prelim.*

Carbon

Lead



$$R_i^{p/A} = \frac{f_i^{p/A}}{f_i^p}$$

$$R_i^A = \frac{Z f_i^{p/A} + N f_i^{n/A}}{Z f_i^p + N f_i^n}$$

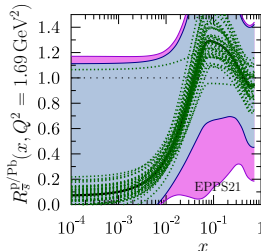
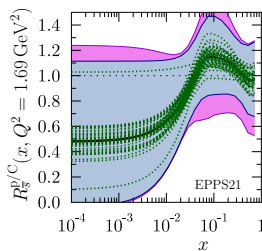
Fit results at parametrization scale – strange and glue

Bound-proton modifications *Prelim.*

Carbon

Lead

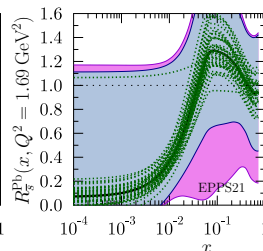
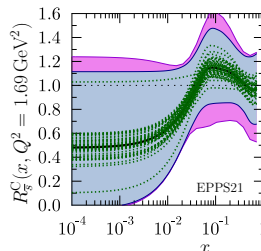
$\bar{s}$



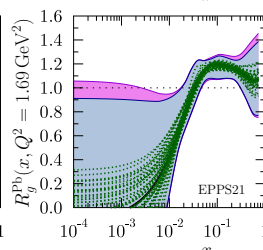
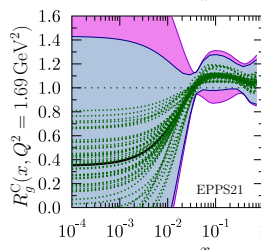
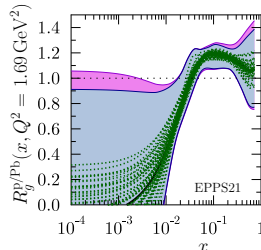
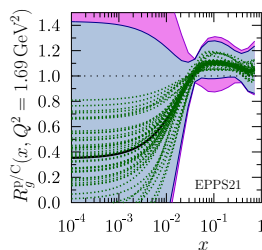
Full-nucleus modifications *Prelim.*

Carbon

Lead



g



$$R_i^{p/A} = \frac{f_i^{p/A}}{f_i^p}$$

$$R_i^A = \frac{Z f_i^{p/A} + N f_i^{n/A}}{Z f_i^p + N f_i^n}$$