Fitting unintegrated PDFs for LHC

EINN2009, Milos, Greece

Albert Knutsson, DESY

Collaborators: A. Bacchetta, H. Jung, K. Kutak

Outline

- Introduction unintegrated PDFs and Monte Carlo
- Determining the unintegrated PDFs
 - the fitting procedure
 - the x dependence
 - the k_t-dependence
- Summary







J. Collins, H. Jung hep-ph/0508280









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Need k_t-dependence. Kinematics correctly treated by using unintegrated PDFs. Gives significant transverse momentum of in the final state.







• Take derivative from PDF:

$$f(x,k_{\perp}^2) \sim \frac{d\Delta x g(x,k_{\perp}^2)_{\rm DGLAP}}{d\log k_{\perp}^2}$$

• The KMR approach. Use normal PDFs. Let the last emission generate transverse momentum via the Sudakov form factor.

→ Can be used in the DGLAP approach, where a strong ordering in k_t is assumed.

In this talk:

•CCFM (Catani Ciafaloni Fiorani Marchesini) approach.

Parton evolution with angular ordering instead of strong ordering in k_t.

Use a unintegrated PDF with parameters determined by fits to data.





• uPDF starting disitribution (example):

$$xA_0(x, k_T, \bar{q}_0) = N \cdot x^{-B} \cdot (1-x)^C \cdot exp(-\frac{(k_T - \mu)^2}{2\sigma^2})$$

 Defined at some starting scale and evolved to higher scales by emissions of gluons according to the *CCFM* evolution scheme.
 Angular ordering of emitted gluon (Color coherence). No explicit k_t-ordering.



- CCFM is usually referred to as the bridge between DGLAP and BFKL.
- CCFM and uPDFs are fully implemented in the general purpose ep/pp MC generator Cascade (H. Jung, Comput.Phys.Commun.143:100-111,2002).





Determining the uPDF parameters



ty

Work done together with A. Bacchetta, H. Jung, K. Kutak

$$xA_0(x, k_T, \bar{q}_0) = N \cdot x^{-B} \cdot (1-x)^C \cdot exp(-\frac{(k_T - \mu)^2}{2\sigma^2})$$

N: Normalization B: Small x behaviour C=4: Large x behaviour (Roughly fixed by momentum sum rules.) μ , σ : Determines the shape of the intrinsic k_T of the gluon

The parameters N,B,C, μ , $\!\sigma,\,$ are not theoretically calculable.

We need to fit the uPDF to experimental data.





- Calculate cross-section using Monte Carlo for a given set of parameter values
 Compare to data, calculate Chi2 and feed it to MINUIT
- 3. MINUIT (e.g. the MIGRAD method) estimates new parameter values
- 4. Iterate 1. 3. until Chi2 is minimized





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This means that if MINUIT needs 100 iterations to minimize Chi2, the generator is run 100 times, not simultaneously:

If one MC generator run takes 1 hour (understatement), the minimization takes 100hours.

One may need exclusive measurements

A lot of MC statistics. Minimization >> 100h.

Also delicate: Fitting several "event types" simultaneously, e.g. Charm production and inclusive jet production Above method makes separated event generation difficult.





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New Approach: Describe parameter dependence before parameter fitting, by using a grid in parameter space.







Simplest possible example 1 parameter, 1 data cross-section

1. Build up the grid









Simplest possible example 1 parameter, 1 data cross-section

2. Determine polynomial using SVD









Simplest possible example 1 parameter, 1 data cross-section

3. Minimize Chi2 to data







If you have a CPU farm (or use the *GRID*) this ultimately takes the time of running the MC generator once.





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2. Fit polynomials to the Monte Carlo grid.

$$\sigma_{\text{poly}} = A + \sum_{1}^{N} B_i \cdot p_i + \sum_{1}^{N} C_i \cdot p_i^2 + \sum_{i=1}^{N} \sum_{j=i+1}^{N} D_{ij} \cdot p_i p_j + H.O.$$

A, B, C and D are determined by fitting the polynomial to the parameter grid by Singular Value Decomposition.





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$$A, B, C \text{ and } D \text{ are determined}$$
by fitting the polynomial to the parameter grid by Singular Value Decomposition. Takes care of correlation between parameters





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$$\begin{split} \sigma_{\text{poly}} &= A + \sum_{1}^{N} B_i \cdot p_i + \sum_{1}^{N} C_i \cdot p_i^2 + \sum_{i=1}^{N} \sum_{j=i+1}^{N} D_{ij} \cdot p_i p_j + H.O. \\ A, B, C \text{ and } D \text{ are determined} \\ \text{by fitting the polynomial to the} \\ \text{parameter grid by SVD.} \end{split}$$

Step 1. and 2. are done for each data point in the measurement.





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Step 1. and 2. are done for each data point in the measurement.

3. Determine PDF parameters, p_i , by fitting all the polynomials to data simultaneously Also this takes only a few seconds.

Step 3. is done by Chi2-minimization using MINUIT.





$$xA_0(x, k_T, \bar{q}_0) = N \cdot x^{-B} \cdot (1-x)^C \cdot exp(-\frac{(k_T - \mu)^2}{2\sigma^2})$$

Used in the CASCADE MC generator: Evolved according to the CCFM equation – parton showers – (hadronization)

•First goal determine the x-dependence: Normalization (N) and the small x behaviour (B)





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•Use the proton structure function, F2. The data covers a large range in x

3 • 10⁻⁵ < x < 2 • 10⁻¹ 1 < Q² < 150 GeV²

....but should be fairly insensitive to the kt-dependent part of the gluon. Inclusive measurement with no restrictions on the hadronic final state.







$$xA_0(x, k_T, \bar{q}_0) = N \cdot x^{-B} \cdot (1-x)^C \cdot exp(-\frac{(k_T - \mu)^2}{2\sigma^2})$$

Fitting F2 in the range x < 0.005, Q2>4.5, gives:

<u>Minimum</u> N = 0.807 +/- 0.016 B = 0.029 +/- 0.004 Chi2/ndf=1.2 This is a good fit which reconstructs the parameter values in an already existing PDF tuned to F2 within the same kinematic range... (Good validation of the method.)



Fitting the x dependence to the proton structure - F2

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= not fitted

Bad description of data outside the fitted range.



DESY

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Bad description of data outside the fitted range.







However if we open up the phase space and fit all F2 data points, we obtain the minimum:

<u>Minimum</u> N = 0.767 +/- 0.001 B = 0.028 +/- 0.000 Chi2/ndf = 5.4

Essentially the same minimum

Data at high and low x are still not described.

Suggests that we need more freedom in the fit needed.





Inspired by the CTEQ group we added an extra factor in the PDF parameterization:

$$xA_0(x,k_T,\bar{q}_0) = \mathbf{N} \cdot x^{-\mathbf{B}} \cdot (1-x)^C \underbrace{(1-\mathbf{D}x)}_G(k_T)$$

...and performed a 3 dimensional fit of N,B and D.

Fitting the x dependence to the proton structure - F2

$$xA_0(x,k_T,\bar{q}_0) = N \cdot x^{-B} \cdot (1-x)^C \cdot (1-Dx) \cdot G(k_T)$$

Fitting F2 over the full range in x gives a slightly different gluon then before. uPDF allowed to be more pronounced at low and high x:

 $xA(x,k_{t}^{2},\mu^{2})$ -0.03 ^{-0.08} (1+5.1x) 3.5 **Minimum** N = 0.487 + - 0.0073 B = 0.097 +/- 0.003 2.5 D = -5.10 + - 0.35Chi2/ndf = 2.8 2 (Before: Chi2/ndf = 5.4) k_t²=1 GeV² 1.5 -3 -2 -1 log(x) χ^2 /ndf χ^2 /ndf χ^2 /ndf 30 50 25 20 20 10 0.02 0.16 6 **0[°]′7** № -15 0.04 -15 -10

PHYSICS AT THE TERA SCALE Helmholtz Alliance

Fitting the x dependence to the proton structure - F2





<u>Minimum</u> N = 0.487 +/- 0.007 B = 0.097 +/- 0.003 D = -5.10 +/- 0.35 Chi2/ndf = 2.8

The extra factor (1-Dx) gives a significant improvement of the data description at low and high x.

Blue dashed line is the new fit







Fitting the k_t dependence of the uPDF







Integrated PDF: DGLAP

- LO: Gluon collinear with proton $k_{t,gluon} = 0$ $\Delta E_{T,jets} = 0$ in HCM
- Higher orders: $k_{t,\text{gluon}} \neq 0$ $\Delta E_{T,\text{jets}} \neq 0$

Unintegrated PDF: CCFM or BFKL

 $k_{t,\text{gluon}} \neq 0$ $\Delta E_{t,\text{jets}} \neq 0$ already at LO







Fit unintegrated gluon density to HERA di-jet data

Target hard di-jets. Dominated by BGF, sensitivity to gluon.



Data from H1 Collab., A. Aktas et al., Eur. Phys. J. C33 (2004) 477 Inclusive Dijet Production at Low x_Bj in DIS





 $\langle Q^2 \rangle = 13 \text{ GeV}^2$

 $\langle \mathbf{x} \rangle = 1.2 \times 10^{-3}$

 $\langle Q^2 \rangle = 17 \text{ GeV}$

 $\langle \mathbf{x} \rangle = 1.4 \times 10^{-3}$

 $\langle O^2 \rangle = 25 \text{ GeV}$

 $\langle Q^2 \rangle = 39 \text{ GeV}$

 $\langle \mathbf{x} \rangle = 2.2 \times 10^{-3}$

 $\langle \mathbf{x} \rangle = 1.7 \times 10^{-3}$

 $\langle \mathbf{x} \rangle = 7.2 \times 10^{-4}$

 $\langle \mathbf{x} \rangle = 7.2 \times 10^{-4}$

 $\langle x \rangle = 7.8 \times 10^{-4}$

0









determined from fit to F2-data (H.Jung, Comp.Phys.Com. 143:100-111,2002) The new fit to the dijet data suggest stronger rising x and a shifted gaussian for k_t.

• Sensitivity to the low k_t-region.

...but still we are in the perturbative region - require hard jets in the final state. When measuring F2 at low Q2, there is risk that we leave the perturbative range.

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Saturation effects mimicked?
 The suppression at low k_t is also
 seen when using saturated PDFs.

Saturation of parton density due to recombination of partons.

Jung, Kutak. arXiv 0812.4082

•So far only unintegrated gluons (i.e. indirectly also sea quarks) in CASCADE.

•Valence quarks expected to be relevant for LHC. For example high Pt production:

Two scale process. With relevant physics for both $x \rightarrow 0$ and $x \rightarrow 1$. High sensitivity to parton dynamics.

Currently the k_t dependent quark PDF is taken from derivated CTEQ5.1.

Deak, Hautmann, Jung, Kutak. Published in JHEP within short. arXiv:0908.0538

- Different data sensitive to different parts of the gluon.
- The x-dependent part of the uPDF is under control. Adding the extra factor (1-Dx) in the parameterization improves the description of F2 significantly.
- New interesting knowledge can be obtained from di-jets, suggests a suppressed gluon at low k_t. Similar effect as if using saturation approach.
- We successfully use a fast approach for fitting the uPDF. The method is based on grid interpolation in the uPDF parameter space.

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Chi2 as a function of Bg for different N values

Dijets using Et and Eta cross-sections

(data: DESY 03-160)

Scans

 The statistical error of the MC predictions are propagated to the coefficients of the polynomial fitted to the MC grid. Only the statistical errors of the MC enters, and thus a simple Chi2 calculation can be used:

$$\chi^2 = \Sigma \frac{(X_{MC} - X_{Polynomial})^2}{(\text{Error}_{MC})^2}$$

•In the fit of the MC parameters to the experimental data the uncorrelated errors and the different correlated errors are treated separately according to:

$$\chi^2 = \Sigma \frac{(X_{Data} - X_{Polynomial})^2}{\alpha^2} - \Sigma_j \Sigma_{j'} B_j (A^{-1})_{jj'} B_{j'}$$

 $lpha^2=$ Sum of uncorrelated errors (data and polynomial)

 $\sum_{j} \sum_{j'} B_j (A^{-1})_{jj'} B_{j'} = \text{Term related to the correlated systematic errors} (\text{vector } B\text{), and their correlations (matrix } A\text{)}$

(From the CTEQ group, hep/ph/0101051)

Fitting F2 in x<0.005, Q2>4.5 GeV. Using the usual PDF parameterization (N x^(-B))

N = 0.807 +/- 0.016 B = 0.029 +/- 0.004 Chi2/ndf = 1.2

Fitting the full F2, 1<Q2<150 GeV using the usual parameterization (N x^(-B))

N = 0.767 +/- 0.001 B = 0.028 +/- 0.000 Chi2/ndf = 5.4

Fitting F2 in x<0.005, Q2>4.5 GeV using (N x^(-B) (1-Dx))

N = 0.550 +/- 0.043 B = 0.082 +/- 0.011 D = -5.38 +/- 1.2 Chi2/ndf = 1.1

Fitting the full F2, 1<Q2<150 GeV using (N x^(-B) (1-Dx)) N = 0.487 +/- 0.007 B = 0.097 +/- 0.003 D = -5.10 +/- 0.35 Chi2/ndf = 2.8

The unintegrated gluon density

The uPDF starting distribution:

$$xA_0(x, k_T, \bar{q}_0) = N \cdot x^{-B} \cdot (1-x)^C \cdot exp(-\frac{(k_T - \mu)^2}{2\sigma^2})$$

 $\begin{array}{l} \textit{N: Normalization (fitted)} \\ \textit{B: Small x behaviour (fitted)} \\ \textit{C=4: Large x behaviour (kept fixed)} \\ \mu, \ \sigma \ \textit{Determines the shape of the intrinsic } k_T \ \textit{of the gluon below } k_T = 1.2 \ \textit{GeV} (\ \mu \ \textit{fitted}) \end{array}$

Calculated at some starting scale (\overline{q}_0). The uPDF is calculated for higher scales by emissions of gluons according to the *CCFM* evolution scheme.

The parameters N,B,C, μ , $\!\sigma,\,$ are not theoretically calculable.

We need to fit the uPDF to experimental data. Albert Knutsson - EINN 2009 - Greece

The new approach was developed for tuning Monte Carlo models

Suggested already 12 years ago...

"Tuning and test of fragmentation models based on identified particles and precision event shape data."

Z.Phys.C73:11-60,1996

Also work on Tuning MC in Lund.

We try to carry out the same method for fitting uPDFs.

DESY

Number of Monte Carlo grid points > Coefficients

Overdetermined system

$$\sigma_{\text{poly}}(p_1, p_2) = A + B_1 p_1 + B_2 p_2 + C_1 p_1^2 + C_2 p_2^2 + C_3 p_1 p_2 + H.O.$$

i.e.

$$\begin{split} P_{n,m}X_m &= \sigma_{n,poly} \quad \text{where} \quad X_n = (A, B_1, B_2, C_1, C_2, C_3, \ldots) \\ P_n &= (1, p_1, p_2, p_1^2, p_1 p_2, p_2^2, \ldots) \\ n &= \textit{Grid point} \end{split}$$

Approach based on SVD algorithm:

To obtain solution we minimize $|PX-\sigma|$ by χ^2 -minimization

Could also use MINUIT, but it is sensitive on starting values.

•The method is implemented into a program – PROFFIT check for updates on www.hepforge.org/PROFFIT.

•A lot of data available for tuning in hztool

("HZTool is a library of routines which will allow you to reproduce an experimental result using the four-vector final state from Monte Carlo generators."

In the future replaced by RIVET)

Comparison to existing uPDF

Coefficients in 4 th order				
polynomial determined from:	SVD	MINUIT	MINUIT ba	d starting values
Chi2 [Polynomial-MC1/ndf·	-25404.9082 765676.064 357293.297 3091.77111 114841.02 166905.85 -31421.3098 -927589.152 -60480.3599 2524.9688 -1064150.37 5799612.85 12981.5342 2592311.1 313456.597 -26463.8328 429940.571 318899.245 -23885.6361 1446.24668 855372.625 -3554618. 97974.8469 -5838295.7 -214807.392 -9020.6301 10567702.9 -437402.716	-25315.4624 762720.857 358067.861 2347.15353 140499.037 157433.813 -32900.9618 -927572.803 -61180.0691 4162.80871 -1135039.64 5804476.94 16228.0397 2623536.1 315635.922 -26091.529 419854.565 320294.755 -23727.0438 525.83837 918985.733 -3552083.15 95580.1082 -5848044.35 -216430.586 -9326.67473 10534823.1 -439016.175	358.56777 694969.672 -52198.7414 3582.44715 52826.8157 1929633.41 -61072.7505 98180.9293 -12538.2387 1618.16049 -465510.961 1334921.4 104971.842 1662889.02 284934.203 -22961.2676 553178.07 -72213.1127 4989.53369 308.886517 -135820.813 354725.482 203313.617 536985.25 -978685.349 -22832.1865 2166665.32 427727.795	For example here, large difference between Coefficients. Resulting in that MINUIT gets stuck in local minimum
	1.0			-