

On statistical methods of structure function extraction

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Several methods of statistical analysis are proposed and analyzed in application for a specific task – extraction of the structure functions from the cross sections of deep inelastic interactions of any type. We formulate the method based on the orthogonal weight] functions and on an optimization procedure of errors minimization as well as methods underlying common χ^2 minimization. Effectiveness of these methods usage is estimated by the statistical parameters such as bias, extraction variance etc., for sample deep inelastic scattering data set.

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Introduction

Precise extraction and analysis of the nucleon's structure functions from the deep inelastic scattering experiments plays great role in current understanding of the particles structure and development of quantum field theory. Deep inelastic scattering (DIS) processes or probing internal structure of nucleon by a pointlike particle at small distances and high energies provide such information about nucleon structure. As soon as experimental facilities to investigate polarized particle interactions have appeared, interest moved mainly to the study of the polarized DIS. Phenomenologically obtained cross sections and asymmetries contain such information as polarized structure functions g_1 and g_2 (or sometimes called g_5 or g_6) or partonic composition of nucleon's spin. To extract these data precisely it's necessary to count on radiative corrections and background effects for obtaining pure cross sections (so-called unfolding of radiative smearing, see e.g.) for further extraction of the structure functions. This paper is devoted to some statistical aspects in application to extraction of the structure functions from the Born interaction parameters.

The cross sections for neutral and charged current deep inelastic scattering both on unpolarized and polarized nucleon targets can be written in the following “separable” form (see e.g. [1]):

$$d^2\sigma/dxdy \equiv \sigma_{xy} = \sigma_0 \sum_k Y_k(y, x, Q^2) W_k(x, Q^2),$$

where $Y_k(y, x, Q^2)$ are known functions, which in the Bjorken limit depend on y only and they are y -polynomial. $W_k(x, Q^2)$ are hadronic structure functions. For massless leptons the cross section is parameterized by a set of three unpolarized structure functions

$$F_1(x, Q^2), F_2(x, Q^2), F_3(x, Q^2)$$

as well as five polarized functions

$$g_1(x, Q^2), g_2(x, Q^2), g_3(x, Q^2), g_3(x, Q^2), g_3(x, Q^2), g_3(x, Q^2).$$

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One can derive desired (un)polarized structure functions operating directly with asymmetries, nevertheless, as asymmetries and corresponding cross sections are being related with each other, we propose here methods operating only with the cross sections. Methods presented below can be easily generalized for more specific cross section expressions, e.g. extended phenomenological models with larger number of the structure functions or more complicated form of the Y_k functions.

The structure of the given article is the following: after initial experimental (statistical) data being introduced we sketch orthogonal weight function method and its optimization technique for the simplest case of the cross section that includes two structure functions $f(x)$ and $g(x)$ for methodological purpose (in the case of polarized particles DIS one can get such expressions by subtracting corresponding cross sections with opposite spin directions). Then we offer χ^2 minimization methods adapted to a given task. Last section compares mentioned methods by means of numerical estimation.

Initial experimental data

Consider normalized on a σ_0 the cross section of the simplest form

$$\sigma_{xy} = Y_+(y)f(x) + Y_-(y)g(x), \quad (1)$$

where Y_{\pm} reads

$$Y_{\pm} = 1 \pm (1 - y)^2. \quad (2)$$

Let's suppose that we know values of cross sections σ_{xy} (in other words – a counted number of events in the bin of a histogram) and its errors $\Delta\sigma_{xy}$ at some x, y -lattice (grid) of experimental kinematical points. Defined grid can be either regular or irregular, so we will distinguish the following cases:

1. Regular (rectangular) x, y -grid

$$M = \mathbf{X} \otimes \mathbf{Y}, \quad \mathbf{X} = x_1, \dots, x_m; \quad \mathbf{Y} = y_1, \dots, y_n.$$

2. Regular (rectangular) x, Q^2 -grid

$$\tilde{M} = \mathbf{X} \otimes \mathbf{Q}^2, \quad \mathbf{X} = x_1, \dots, x_m; \quad \mathbf{Q}^2 = Q_1^2, \dots, Q_n^2.$$

3. Non-regular (irregular) x, y - and x, Q^2 -grid – arbitrary set of points (bins).

Although, it's not crucial here for the lattice to be regular (besides experimenters mostly gather events on irregular set of bins), but at first we implement methods for regular grids and then propose schemes to generalize and apply them for an irregular set of points (bins) using simple interpolation considerations.

Orthogonal weight function method.

Lets begin with the rectangular x, y -grid M and take the parameters a_{\pm} of some preliminary chosen weight function $\omega(y; a)$ in such a way that the following orthogonality condition fulfils:

$$\sum_y \omega(y; a_{\pm}) Y_{\pm} = 0,$$

whereas the following requirement holds:

$$\sum_y \omega(y; a_{\mp}) Y_{\pm} \neq 0.$$

In another way, preceding expressions mean that ω acts as a projection operator. Then one can extract structure function by projection in the following form:

$$\begin{aligned} f(x) \pm \Delta f(x) &= \frac{1}{\sum_y \omega(y; a_-) Y_+(y)} \sum_y \omega(y; a_-) [\sigma_{xy} \pm \Delta \sigma_{xy}], \\ g(x) \pm \Delta g(x) &= \frac{1}{\sum_y \omega(y; a_+) Y_-(y)} \sum_y \omega(y; a_+) [\sigma_{xy} \pm \Delta \sigma_{xy}]. \end{aligned} \quad (3)$$

For example, one may choose the simplest weight functions as a $\omega(y; a) = 1 + ay$, then a_{\pm} parameters take the following values:

$$a_{\pm} = -\frac{\sum_y Y_{\pm}(y)}{\sum_y y Y_{\pm}(y)}. \quad (4)$$

Note that in this case a_{\pm} depend only on chosen data grid.

As presented above the procedure of structure function extraction implies rough estimation of the standard deviation (error) values in eq. (3). To find correctly uncertainty in the fitted structure functions one should adhere to the standard procedure of the variance (noted here as D) calculation of random variables function, which assumes here for a given x -bin and a set of n y -bins (the same holds for g -function)

$$\begin{aligned} D[f(\sigma(x, y_1), \dots, \sigma(x, y_n))] &= \frac{1}{\left[\sum_{i=1}^n \omega(y_i; a_-) Y_+(y_i) \right]^2} \times \left\{ \sum_{i=1}^n \omega(y_i; a_-)^2 D[\sigma(x, y_i)] + \right. \\ &\quad \left. + \sum_{i \neq j} \omega(y_i; a_-) \omega(y_j; a_-) \rho[\sigma(x, y_i), \sigma(x, y_j)] \sqrt{D[\sigma(x, y_i)] D[\sigma(x, y_j)]} \right\}. \end{aligned} \quad (5)$$

In the case of the uncorrelated data with diagonal correlation matrix $\rho[\sigma(x, y_i), \sigma(x, y_j)] = \delta_{ij}$ we specify correct estimation for deviation of the structure function values

$$\Delta f(\sigma(x, y_1), \dots, \sigma(x, y_n)) = \frac{1}{\left| \sum_{i=1}^n \omega(y_i; a_-) Y_+(y_i) \right|} \sqrt{\sum_{i=1}^n \omega(y_i; a_-)^2 D(\sigma(x, y_i))}, \quad (6)$$

that can be larger or smaller than

$$\frac{1}{\sum_{i=1}^n \omega(y_i; a_-) Y_+(y_i)} \sum_{i=1}^n \omega(y_i; a_-) \Delta \sigma(x, y_i), \quad (7)$$

dependently of the $\omega(y; a)$ sign. In these formulas we assume that the lattice has no uncertainty, i.e. $D[a_{\pm}] = 0$ as given initially.

Optimization procedure

To minimize errors of the structure functions that are extracted from experimental data $\sigma_{xy} \pm \Delta \sigma_{xy}$ one can apply an optimization procedure of the following type. Let's introduce auxiliary functions

$$A(a) = \sum_y \omega(y; a) Y_+(y), \quad B(a) = \sum_y \omega(y; a) Y_-(y),$$

$$S(x; a) = \sum_y \omega(y; a) \sigma_{xy}, \quad \Delta S(x; a) = \sum_y \omega(y; a) \Delta \sigma_{xy}.$$

Two systems of equations for $f(x)$ and $g(x)$ and errors $\Delta f(x)$ and $\Delta g(x)$ are

$$\begin{cases} A(a)f(x) + B(a)g(x) = S(x; a), \\ A(b)f(x) + B(b)g(x) = S(x; b); \end{cases}$$

$$\begin{cases} A(a)\Delta f(x) + B(a)\Delta g(x) = \Delta S(x; a), \\ A(b)\Delta f(x) + B(b)\Delta g(x) = \Delta S(x; b). \end{cases}$$

To find their solutions define the following determinants:

$$\Delta(a, b) = \begin{vmatrix} A(a) & B(a) \\ A(b) & B(b) \end{vmatrix}, \quad \Delta_1(x; a, b) = \begin{vmatrix} S(x; a) & B(a) \\ S(x; b) & B(b) \end{vmatrix}, \quad \Delta_2(x; a, b) = \begin{vmatrix} A(a) & S(x; a) \\ A(b) & S(x; b) \end{vmatrix}$$

and

$$\delta \Delta_1(x; a, b) = \begin{vmatrix} \Delta S(x; a) & B(a) \\ \Delta S(x; b) & B(b) \end{vmatrix}, \quad \delta \Delta_2(x; a, b) = \begin{vmatrix} A(a) & \Delta S(x; a) \\ A(b) & \Delta S(x; b) \end{vmatrix}.$$

Then we get the solution

$$f(x) = \frac{\Delta_1(x; a, b)}{\Delta(a, b)}, \quad \Delta f(x) = \frac{\delta \Delta_1(x; a, b)}{\Delta(a, b)}, \quad g(x) = \frac{\Delta_2(x; a, b)}{\Delta(a, b)}, \quad \Delta g(x) = \frac{\delta \Delta_2(x; a, b)}{\Delta(a, b)}.$$

The optimal values of the parameters a and b can be found from the condition of the errors minimization

$$\min_{\{a, b\}} [w_f |\Delta f(x)| + w_g |\Delta g(x)|], \quad (8)$$

where w_f and w_g – optional weight factors.

So, optimization procedure implies determination of the optimal parameters a_k and b_k for each experimental point x_k in order to minimize errors in this point. As a result we have the estimation for the mean values of the structure functions extracted at given experimental points and rough estimations for corresponding errors (deviations):

$$f(x) = \frac{\Delta_1(x; a, b)}{\Delta(a, b)}, \quad \Delta f(x) = \frac{\delta \Delta_1(x; a, b)}{\Delta(a, b)}, \quad g(x) = \frac{\Delta_2(x; a, b)}{\Delta(a, b)}, \quad \Delta g(x) = \frac{\delta \Delta_2(x; a, b)}{\Delta(a, b)}, \quad (9)$$

for each $k = 1, 2, \dots, n$.

It should be noted that by such solution we get only approximate values for errors (likewise mentioned above argument about correct deviation values), nevertheless one can easily obtain correct values by finding minimum solution (8) analytically and repeating formulas (5), (6).

A difficulty arises from the fact that the method implies experimental data on a (rectangular) x, y -lattice which is rarely used, furthermore as a rule experimental bins chosen for analysis and fitting are not uniformly distributed (e.g. see kinematics in experimental reports [3, 4]). First of all we propose to apply the same scheme to regular (rectangular) x, Q^2 -lattice. The difference between x, y - and x, Q^2 -data consist in principle only in redefinition of the structure functions. Let's modify Y expressions (2):

$$Y_{\pm}(Q^2) = 1 \pm \left(1 - \frac{Q^2}{sx}\right)^2, \quad \tilde{Y}_+ = Q^4, \quad \tilde{Y}_- = 2Q^2,$$

$$\sigma_{x, Q^2} \sim 2 \frac{f(x)}{x} + \frac{\tilde{f}(x)}{x} \frac{1}{s^2 x^2} \tilde{Y}_+ - \frac{\tilde{f}(x)}{x} \frac{1}{sx} \tilde{Y}_-$$

and use the similar test weight function $\omega(Q^2; a) = 1 + aQ^2$. The same orthogonality relations take the form of

$$\sum_{Q^2} \omega(Q^2; a_{\pm}) \tilde{Y}_{\pm} = 0, \quad \sum_{Q^2} \omega(Q^2; a_{\pm}) \neq 0,$$

$$a_{\pm} = -\frac{\sum_{Q^2} \tilde{Y}_{\pm}}{\sum_{Q^2} Q^2 \tilde{Y}_{\pm}},$$

As a result one can obtain

$$\sum_{Q^2} \sigma_{x,Q^2} \omega(Q^2; a_{\pm}) = 2 \frac{f(x)}{x} \sum_{Q^2} \omega(Q^2; a_{\pm}) + \frac{\tilde{f}(x)}{x} \frac{1}{s^2 x^2} \sum_{Q^2} \omega(Q^2; a_{\pm}) \tilde{Y}_+ - \frac{\tilde{f}(x)}{x} \frac{1}{s x} \sum_{Q^2} \omega(Q^2; a_{\pm}) \tilde{Y}_-,$$

$$\sum_{Q^2} \sigma_{x,Q^2} \omega(Q^2; a_+) = 2 \frac{f(x)}{x} \sum_{Q^2} \omega(Q^2; a_+) - \frac{\tilde{f}(x)}{s x^2} \sum_{Q^2} \omega(Q^2; a_+) \tilde{Y}_-,$$

$$\sum_{Q^2} \sigma_{x,Q^2} \omega(Q^2; a_-) = 2 \frac{f(x)}{x} \sum_{Q^2} \omega(Q^2; a_-) + \frac{\tilde{f}(x)}{s^2 x^3} \sum_{Q^2} \omega(Q^2; a_-) \tilde{Y}_+,$$

where $\tilde{f}(x) = f(x) - g(x)$. Hence

$$f(x) = \frac{x}{2} \frac{\left[\sum_{Q^2} \sigma_{x,Q^2} \omega(Q^2; a_+) \sum_{Q^2} \omega(Q^2; a_-) \tilde{Y}_+ + x s \sum_{Q^2} \sigma_{x,Q^2} \omega(Q^2; a_-) \sum_{Q^2} \omega(Q^2; a_+) \tilde{Y}_- \right]}{\left[\sum_{Q^2} \omega(Q^2; a_+) \sum_{Q^2} \omega(Q^2; a_-) \tilde{Y}_+ + x s \sum_{Q^2} \omega(Q^2; a_-) \sum_{Q^2} \omega(Q^2; a_+) \tilde{Y}_- \right]},$$

$$g(x) = f(x) - x \frac{s^2 x^2 \left[\sum_{Q^2} \sigma_{x,Q^2} \omega(Q^2; a_-) \sum_{Q^2} \omega(Q^2; a_+) - \sum_{Q^2} \sigma_{x,Q^2} \omega(Q^2; a_+) \sum_{Q^2} \omega(Q^2; a_-) \right]}{\left[\sum_{Q^2} \omega(Q^2; a_+) \sum_{Q^2} \omega(Q^2; a_-) \tilde{Y}_+ + x s \sum_{Q^2} \omega(Q^2; a_-) \sum_{Q^2} \omega(Q^2; a_+) \tilde{Y}_- \right]}.$$

To obtain values of the absolute error one should substitute $\Delta \sigma_{x,Q^2}$ instead of σ_{x,Q^2} .

One of the advantages of x, Q^2 -grid consists in possibility of appropriate usage of the interpolation through Q^2 range, contrary to the y variable range, which combines different $Q^2 = sxy$ values in an unhandy way. Special optimization procedure similar to one described above by formulas (9) is also applicable over here. Used scheme will be reliable if one obtains data for rectangular x, Q^2 -region with narrow Q^2 range to neglect existing Q^2 -dependency of structure functions (but still with at least two different Q^2 values). It should be noted that for non-uniform grids the summing range \sum_{Q^2} may depend on selected x -bin and consequently $a_{\pm} \rightarrow a_{\pm}(x)$ becomes a function of the given x value. Nevertheless, above mentioned scheme works in the same way.

If one has x, Q^2 -lattice of the data, which cannot be grouped in the certain x -bins, but experimental points are distributed in the vicinity of certain x -values, one can eliminate these difficulties by interpolation methods, as briefly mentioned below.

Interpolation

Here we describe simplest ways to manage with data on irregular grids. One can use the following Taylor formula for the cross section value near given x_i value

$$\sigma_i^{int}(x, Q^2) = \sigma^{ex}(x_i, Q^2) + \frac{\sigma^{ex}(x_i, Q^2)}{\sigma^{fit}(x_i, Q^2)} (x - x_i) \partial_x \sigma^{fit}(x_i, Q^2) + \dots$$

$$\sigma^{int}(x, Q^2) = [\text{weighted average}] = \frac{\sum_i \sigma_i^{int}(x, Q^2) w(x - x_i)}{\sum_i w(x - x_i)}.$$

Here the additional fraction is included with the purpose of normalization - but it may be omitted. These formulas match up the case when one has data with a small dispersion in x -values. One can employ some external parameterization, denoted here as σ^{fit} , to fix missing experimental x points by simple interpolation. But total uncertainty of the extracted $f(x)$ and $g(x)$ values increases by the theoretical uncertainty of these parameterizations.

In case of significant Q^2 range Q^2 -dependence cannot be omitted and one can treat $f(x, Q^2)$ approximately using similar Taylor series of f for Q^2 near given Q_0^2 value

$$f(x) \rightarrow f^{int}(x, Q^2) = f(x, Q_0^2) + \frac{f(x, Q_0^2)}{f^{fit}(x_i, Q_0^2)} (Q^2 - Q_0^2) \partial_{Q^2} f^{fit} + \dots,$$

or introducing some fixing factor $\delta(x, Q^2)$ of a pre-given form, e.g. $f(x)\delta(x, Q^2)$, but these ideas require $f(x)$ and $g(x)$ as well as Y_{\pm} and \tilde{Y}_{\pm} to be redesigned.

χ^2 -minimization procedure for the parameters a_{\pm}

Below common χ^2 methods are applied to structure functions extraction. First we preserve usage of the a_{\pm} parameters (thus implicit method, requiring introducing of the orthogonal function ω) and modify common procedure. Let's construct the χ^2 -function of random cross-sections and parameters a_{-}, a_{+} as follows

$$\chi^2(\sigma(x, y_1), \dots, \sigma(x, y_n); a_{-}, a_{+}) = \sum_{i,j=1}^n [\sigma(x, y_i) - m_i(a_{-}, a_{+})] D^{-1}_{ij} [\sigma(x, y_j) - m_j(a_{-}, a_{+})],$$

where $D_{ij} = \rho[\sigma(x, y_i), \sigma(x, y_j)] \sqrt{D[\sigma(x, y_i)] D[\sigma(x, y_j)]}$ is the covariance matrix and $m(a_{-}, a_{+})$ are the expectation values:

$$m_i(a_{-}, a_{+}) = Y_{+}(y_i) \frac{\sum_{j=1}^n \omega(y_j; a_{-}) \sigma(x, y_j)}{\sum_{j=1}^n \omega(y_j; a_{-}) Y_{+}(y_j)} + Y_{-}(y_i) \frac{\sum_{j=1}^n \omega(y_j; a_{+}) \sigma(x, y_j)}{\sum_{j=1}^n \omega(y_j; a_{+}) Y_{-}(y_j)}.$$

As a_{\pm} are supposed to be estimated parameters only, we may neglect dependence of these mean values on the random cross-section values $\sigma(x, y_j)$ regarding them as a set of initial exact numbers.

Next step is to minimize the χ^2 function. In the case of the uncorrelated data with $\rho[\sigma(x, y_i), \sigma(x, y_j)] = \delta_{ij}$ we get the following estimator equations:

$$\begin{aligned} - \sum_{i=1}^n \frac{\partial m_i}{\partial \hat{a}_{-}} D_{ii}^{-1} [\sigma(x, y_i) - m_i(\hat{a}_{-}, \hat{a}_{+})] &= 0, \\ - \sum_{i=1}^n \frac{\partial m_i}{\partial \hat{a}_{+}} D_{ii}^{-1} [\sigma(x, y_i) - m_i(\hat{a}_{-}, \hat{a}_{+})] &= 0. \end{aligned}$$

Consistent solution of the system gives χ^2 -estimators \hat{a}_{\pm} . One can check consistency calculating the derivatives, which should not vanish

$$- \sum_{i=1}^n \frac{\partial^2 m_i}{\partial \hat{a}_{\pm}^2} D_{ii}^{-1} [\sigma(x, y_i) - m_i(\hat{a}_{-}, \hat{a}_{+})] + \sum_{i=1}^n \left(\frac{\partial m_i}{\partial \hat{a}_{\pm}} \right)^2 D_{ii}^{-1} \neq 0,$$

$$-\sum_{i=1}^n \frac{\partial^2 m_i}{\partial \hat{a}_+ \partial \hat{a}_-} D_{ii}^{-1} [\sigma(x, y_i) - m_i(\hat{a}_-, \hat{a}_+)] + \sum_{i=1}^n \frac{\partial m_i}{\partial \hat{a}_-} D_{ii}^{-1} \frac{\partial m_i}{\partial \hat{a}_+} \neq 0.$$

After that one should substitute obtained \hat{a}_\pm into the following expressions to get $f(x)$ and $g(x)$:

$$f(x) = \frac{\sum_{j=1}^n \omega(y_j; \hat{a}_-) \sigma(x, y_j)}{\sum_{j=1}^n \omega(y_j; \hat{a}_-) Y_+(y_j)}, \quad g(x) = \frac{\sum_{j=1}^n \omega(y_j; \hat{a}_+) \sigma(x, y_j)}{\sum_{j=1}^n \omega(y_j; \hat{a}_+) Y_-(y_j)}.$$

Apart from these values, it's necessary to get estimations for the statistical parameters such as bias, deviation etc. This analysis is discussed below.

χ^2 -minimization procedure for the functions of $f(x)$ and $g(x)$

Here we omit a_\pm parameter and estimate functions $f(x)$ and $g(x)$ explicitly (as before we refer to x as to fixed bin), thus changing χ^2 arguments:

$$\chi^2(\sigma(x, y_1), \dots, \sigma(x, y_n); f(x), g(x)) = \sum_{i,j=1}^n [\sigma(x, y_i) - m_i(f(x), g(x))] D_{ij}^{-1} [\sigma(x, y_j) - m_j(f(x), g(x))],$$

where $D_{ij} = \rho[\sigma(x, y_i), \sigma(x, y_j)] \sqrt{D[\sigma(x, y_i)] D[\sigma(x, y_j)]}$ is the covariance matrix and $m(f(x), g(x))$ are the expectation values:

$$m_i(f(x), g(x)) = Y_+(y_i) f(x) + Y_-(y_i) g(x).$$

Next step is to minimize the χ^2 function. In the case of the uncorrelated data with diagonal correlation matrix $\rho[\sigma(x, y_i), \sigma(x, y_j)] = \delta_{ij}$ we get the following estimator equations:

$$\begin{aligned} -\sum_{i=1}^n Y_+(y_i) D_{ii}^{-1} [\sigma(x, y_i) - Y_+(y_i) f(x) - Y_-(y_i) g(x)] &= 0, \\ -\sum_{i=1}^n Y_-(y_i) D_{ii}^{-1} [\sigma(x, y_i) - Y_+(y_i) f(x) - Y_-(y_i) g(x)] &= 0. \end{aligned}$$

Solving this equations we get the following consistent χ^2 -estimates for $f(x)$ and $g(x)$

$$\begin{aligned} \Delta &= \begin{vmatrix} \sum_{i=1}^n Y_+^2(y_i) D_{ii}^{-1} & \sum_{i=1}^n Y_+(y_i) Y_-(y_i) D_{ii}^{-1} \\ \sum_{i=1}^n Y_+(y_i) Y_-(y_i) D_{ii}^{-1} & \sum_{i=1}^n Y_-^2(y_i) D_{ii}^{-1} \end{vmatrix} \neq 0, \\ f(x) &= \frac{1}{\Delta} \begin{vmatrix} \sum_{i=1}^n Y_+(y_i) D_{ii}^{-1} \sigma(x, y_i) & \sum_{i=1}^n Y_+(y_i) Y_-(y_i) D_{ii}^{-1} \\ \sum_{i=1}^n Y_-(y_i) D_{ii}^{-1} \sigma(x, y_i) & \sum_{i=1}^n Y_-^2(y_i) D_{ii}^{-1} \end{vmatrix}, \\ g(x) &= \frac{1}{\Delta} \begin{vmatrix} \sum_{i=1}^n Y_+^2(y_i) D_{ii}^{-1} & \sum_{i=1}^n Y_+(y_i) D_{ii}^{-1} \sigma(x, y_i) \\ \sum_{i=1}^n Y_+(y_i) Y_-(y_i) D_{ii}^{-1} & \sum_{i=1}^n Y_-(y_i) D_{ii}^{-1} \sigma(x, y_i) \end{vmatrix}. \end{aligned} \tag{10}$$

Last method resembles previous optimization procedure (even in the forms of the matrices), but crucial point here is the absence of a parameter (it looks like Y is a weight function itself without any auxiliary parameter). For the case of correlated data one can refer to the same formulas using simple substitutions, e.g. in the matrix form:

$$\sum_{i=1}^n Y_\pm^2(y_i) D_{ii}^{-1} \Rightarrow \sum_{i,j=1}^n Y_\pm(y_i) D_{ij}^{-1} Y_\pm(y_j) = \mathbf{Y}_\pm^T D^{-1} \mathbf{Y}_\pm.$$

Considered case corresponds to the linear χ^2 approach, so using standard statistics one can get the following formulas, e.g. [2]:

$$\begin{aligned} D(f) &= \frac{1}{\Delta^2} F D F^T, & F &= (\mathbf{Y}_-^T D^{-1} \mathbf{Y}_-) \mathbf{Y}_+^T D^{-1} - (\mathbf{Y}_+^T D^{-1} \mathbf{Y}_-) \mathbf{Y}_-^T D^{-1}, \\ D(g) &= \frac{1}{\Delta^2} G D G^T, & G &= (\mathbf{Y}_+^T D^{-1} \mathbf{Y}_+) \mathbf{Y}_-^T D^{-1} - (\mathbf{Y}_-^T D^{-1} \mathbf{Y}_+) \mathbf{Y}_+^T D^{-1}. \end{aligned} \quad (11)$$

Main advantage of the last χ^2 approach (least squares method in the linear case) is that it provides consistent unbiased estimator for f and g with the smallest in its type estimator variance, according to the Gauss-Markov theorem.

Discussion and numerical results

For the purpose of numerical analysis let's take the model parameterization for the structure functions, in the following simplest form:

$$f(x) = C_f x^{\alpha_f} (1-x)^{\beta_f}, \quad g(x) = C_g x^{\alpha_g} (1-x)^{\beta_g}. \quad (12)$$

Then one can construct cross section σ_{xy} and its error $\Delta\sigma_{xy}$ at some points x_1, \dots, x_n and y_1, \dots, y_m , referring to these quantities as expectation value and deviation of some preselected probability distribution function. Also one may include additional random bias if desired. Using such initial assumptions it's easy to compare numerically above listed methods. As expected by the Gauss-Markov theorem, numerical analysis gives both the lowest values of the $\chi^2(f, g)$ function and the correct minimal variances $D(f)$, $D(g)$ for the last χ^2 method. At the same time minimum value $\chi^2(a_+, a_-)$ for χ^2 - a procedure equals to the minimum $\chi^2(f, g)$ value, although presence of the random data in the mean values $m(a_-, a_+)$ increases the errors $D[f(a)]$, $D[g(a)]$. Both orthogonal method and optimized orthogonal method have the same larger χ^2 value for different optimal sets of a and b and for nonoptimized a_{\pm} parameters, though they give different $D(f)$, $D(g)$ values. These estimations include bias to be calculated analytically, contrary to the linear χ^2 - f, g procedure without it. This causes additional peculiarity – one can get in this case either larger $D(f)$ and smaller $D(g)$ (or opposite) in comparison with the χ^2 methods. Detailed analysis can be carried out analytically using standard statistics methods after weight function definition. It should be noted as well the orthogonal weight function method gives the expectation values $f(a_{\pm})$ and $g(a_{\pm})$ (defined by (3)) equal to optimized expectation values $f(a_{\pm\text{opt}})$ and $g(a_{\pm\text{opt}})$ (defined by(9)), but they differs from f and g values obtained using formulas (10). The orthogonal weight functions approaches with formulas (3) and (9) can be used as a approximate estimations. The last mentioned least squares procedure gives reasonable unbiased estimation, and can be used for Born cross-section analysis. It should be noted that discussed methods can be easily generalized for more common case and for various specific purposes.

References

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