

1 The χ^2 in the presence of correlations

Suppose to have an ensemble of n measurements having the following structure:

$$m_i \pm \sigma_{i,\text{stat}} \pm \sigma_{i,\text{unc}} \pm \sigma_{i,\text{corr}}^{(1)} \pm \cdots \pm \sigma_{i,\text{corr}}^{(k)}, \quad (1)$$

where m_i , with $i = 1, \dots, n$, is the central value of the i -th measurement, $\sigma_{i,\text{stat}}$ its (uncorrelated) statistical uncertainty, $\sigma_{i,\text{unc}}$ its uncorrelated systematic uncertainty¹, and $\sigma_{i,\text{corr}}^{(l)}$, with $l = 1, \dots, k$, its correlated systematic uncertainties. With this information at hand, one can construct the full covariance matrix V_{ij} as follows (see for example Ref. [3]):

$$V_{ij} = (\sigma_{i,\text{stat}}^2 + \sigma_{i,\text{unc}}^2) \delta_{ij} + \sum_{l=1}^k \sigma_{i,\text{corr}}^{(l)} \sigma_{j,\text{corr}}^{(l)}. \quad (2)$$

This is a clearly symmetric matrix. Given a set of predictions t_i corresponding to the n measurements of the ensemble, the χ^2 takes the form:

$$\chi^2 = \sum_{i,j=1}^n (m_i - t_i) V_{ij}^{-1} (m_j - t_j) = \mathbf{y}^T \cdot \mathbf{V}^{-1} \cdot \mathbf{y}, \quad (3)$$

where in the second equality we have used the matricial notation and defined $y_i = m_i - t_i$. A convenient way to compute the χ^2 relies on the Cholesky decomposition of the covariance matrix \mathbf{V} . In particular, it can be proven that any symmetric and positive definite matrix \mathbf{V} can be decomposed as:

$$\mathbf{V} = \mathbf{L} \cdot \mathbf{L}^T, \quad (4)$$

where \mathbf{L} is a lower triangular matrix whose entries are related recursively to those of \mathbf{V} as follows:

$$\begin{aligned} L_{kk} &= \sqrt{V_{kk} - \sum_{j=1}^{k-1} L_{kj}^2}, \\ L_{ik} &= \frac{1}{L_{kk}} \left(V_{ik} - \sum_{j=1}^{k-1} L_{ij} L_{kj} \right), \quad k < i, \\ L_{ik} &= 0, \quad k > i. \end{aligned} \quad (5)$$

It is then easy to see that the χ^2 can be written as:

$$\chi^2 = |\mathbf{L}^{-1} \cdot \mathbf{y}|^2. \quad (6)$$

But the vector $\mathbf{x} \equiv \mathbf{L}^{-1} \cdot \mathbf{y}$ is the solution of the linear system:

$$\mathbf{L} \cdot \mathbf{x} = \mathbf{y}, \quad (7)$$

that can be efficiently solved by forward substitution, so that:

$$\chi^2 = |\mathbf{x}|^2. \quad (8)$$

Following this procedure, one does not need to compute explicitly the inverse of the covariance matrix \mathbf{V} , simplifying significantly the computation of the χ^2 .

¹There could be more than one uncorrelated systematic uncertainty. In this case, $\sigma_{i,\text{unc}}$ is just the square root of the sum in quadrature of all the uncorrelated systematic uncertainties.

2 Additive and multiplicative uncertainties

The correlated systematic uncertainties $\sigma_{i,\text{corr}}^{(l)}$ may be either *additive* or *multiplicative*. The nature of the single uncertainties is typically provided by the experiments that release the measurements. A typical example of multiplicative uncertainty is the luminosity uncertainty but there can be others.

Now let us express all the correlated systematic uncertainties $\sigma_{i,\text{corr}}^{(l)}$ as relative to the associate central value m_i , so that we define²:

$$\sigma_{i,\text{corr}}^{(l)} \equiv \delta_{i,\text{corr}}^{(l)} m_i \quad (9)$$

and let us also define $s_i^2 \equiv \sigma_{i,\text{stat}}^2 + \sigma_{i,\text{unc}}^2$ so that Eq. (2) can be rewritten as:

$$V_{ij} = s_i^2 \delta_{ij} + \left(\sum_{l=1}^k \delta_{i,\text{corr}}^{(l)} \delta_{j,\text{corr}}^{(l)} \right) m_i m_j. \quad (10)$$

Now we split the correlated systematic uncertainties into k_a additive uncertainties and k_m multiplicative uncertainties, such that $k_a + k_m = k$. This way Eq. (10) takes the form:

$$V_{ij} = s_i^2 \delta_{ij} + \left(\sum_{l=1}^{k_a} \delta_{i,\text{add}}^{(l)} \delta_{j,\text{add}}^{(l)} + \sum_{l=1}^{k_m} \delta_{i,\text{mult}}^{(l)} \delta_{j,\text{mult}}^{(l)} \right) m_i m_j. \quad (11)$$

It is well known that this definition of the covariance matrix is problematic in that it results in the so-called D’Agostini bias of the multiplicative uncertainties [2]. A possible solution to this problem is the so-called t_0 -prescription [8], where the experimental central value m_i in the multiplicative term is replaced by a fixed theoretical predictions $t_i^{(0)}$, typically computed in a previous fit in which the “standard” definition of the covariance matrix in Eq. (2) (often referred to as *experimental* definition) is used. Applying the t_0 prescription, the covariance matrix takes the form:

$$V_{ij} = s_i^2 \delta_{ij} + \sum_{l=1}^{k_a} \delta_{i,\text{add}}^{(l)} \delta_{j,\text{add}}^{(l)} m_i m_j + \sum_{l=1}^{k_m} \delta_{i,\text{mult}}^{(l)} \delta_{j,\text{mult}}^{(l)} t_i^{(0)} t_j^{(0)}. \quad (12)$$

3 Artificial generation of correlated systematics

In order to implement the definition of the χ^2 discussed above, it is necessary to have the experimental information in terms of the correlated systematic uncertainties $\sigma_{i,\text{corr}}^{(l)}$. This is what the experimental collaborations usually release. However, in some cases this information is given in terms of a covariance matrix. Therefore, one needs to find a workaround to generate correlated systematic uncertainties out of a covariance matrix.

Given a $n \times n$ symmetric matrix \mathbf{C} , it will have n orthonormal eigenvectors $\mathbf{x}^{(i)}$, such that $\mathbf{x}^{(i)} \cdot \mathbf{x}^{(j)} = \delta_{ij}$, each of which will have a non-negative eigenvalue λ_i associated:

$$\mathbf{C} \cdot \mathbf{x}^{(i)} = \lambda_i \mathbf{x}^{(i)}, \quad i = 1, \dots, n. \quad (13)$$

²Note that this redefinition does not change the nature of the uncertainties, additive uncertainties remain additive as well as multiplicative uncertainties remain multiplicative.

If we define:

$$\sigma_{i,\text{corr}}^{(l)} = \sqrt{\lambda_l} x_i^{(l)}, \quad i, l = 1, \dots, n, \quad (14)$$

one can show that:

$$\sum_{l=1}^n \sigma_{i,\text{corr}}^{(l)} \sigma_{j,\text{corr}}^{(l)} = C_{ij}. \quad (15)$$

To prove this equality we start from the following matricial relation:

$$\mathbf{C} = \mathbf{Q} \cdot \mathbf{\Lambda} \cdot \mathbf{Q}^{-1}, \quad (16)$$

where $\mathbf{\Lambda}$ is a diagonal matrix with the eigenvalues λ_i on the diagonal ($\Lambda_{ij} = \lambda_i \delta_{ij}$), while \mathbf{Q} is a matrix whose columns are the eigenvectors $\mathbf{x}^{(i)}$ ($Q_{ij} = x_i^{(j)}$). In addition, since in this particular case $\mathbf{x}^{(i)} \cdot \mathbf{x}^{(j)} = \delta_{ij}$, this implies that:

$$\mathbf{Q}^T \cdot \mathbf{Q} = \mathbf{I} \quad \Rightarrow \quad \mathbf{Q}^{-1} = \mathbf{Q}^T, \quad (17)$$

so that:

$$\mathbf{C} = \mathbf{Q} \cdot \mathbf{\Lambda} \cdot \mathbf{Q}^T. \quad (18)$$

It follows that:

$$C_{ij} = \sum_{k,l=1}^n Q_{ik} \Lambda_{kl} Q_{jl} = \sum_{k,l=1}^n x_i^{(k)} \lambda_k \delta_{kl} x_j^{(l)} = \sum_{l=1}^n \lambda_l x_i^{(l)} x_j^{(l)} = \sum_{l=1}^n \sigma_{i,\text{corr}}^{(l)} \sigma_{j,\text{corr}}^{(l)}, \quad (19)$$

as required.

The matrix \mathbf{C} can be regarded as the correlated contribution to the full covariance matrix \mathbf{V} . In particular, considering Eqs. (2) and (10), one can write:

$$\mathbf{V} = \mathbf{U} + \mathbf{C}, \quad (20)$$

where \mathbf{U} is a diagonal matrix of uncorrelated uncertainties:

$$U_{ij} = s_i^2 \delta_{ij}. \quad (21)$$

This defines the matrix \mathbf{C} as:

$$\mathbf{C} = \mathbf{V} - \mathbf{U}, \quad (22)$$

such that, given a $n \times n$ covariance matrix \mathbf{V} along with its uncorrelated contribution \mathbf{U} , one can generate a set of n *artificial* correlated systematics according to Eq. (14), where \mathbf{C} is given in Eq. (22), for each of the n measurements. This allows us to implement Eq. (12) for the construction of the covariance matrix.

4 Determining the systematic shifts

In order to visualise the effect of systematic uncertainties, it is instructive to compute the *systematic shift* generated by the systematic uncertainties. To do so, we need to write the χ^2 in terms of the so-called “nuisance parameters” λ_α . One can show that the definition of the χ^2 in Eq. (3) is equivalent to [3]:

$$\chi^2 = \sum_{i=1}^n \frac{1}{s_i^2} \left(m_i - t_i - \sum_{\alpha=1}^k \lambda_\alpha \sigma_{i,\text{corr}}^{(\alpha)} \right)^2 + \sum_{\alpha=1}^k \lambda_\alpha^2. \quad (23)$$

The optimal value of the nuisance parameters can be computed by minimising the χ^2 with respect to them, that is imposing:

$$\frac{\partial \chi^2}{\partial \lambda_\beta} = 0. \quad (24)$$

This yields the system:

$$\sum_{\beta=1}^k A_{\alpha\beta} \lambda_\beta = \rho_\alpha, \quad (25)$$

with:

$$A_{\alpha\beta} = \delta_{\alpha\beta} + \sum_{i=1}^n \frac{\sigma_{i,\text{corr}}^{(\alpha)} \sigma_{i,\text{corr}}^{(\beta)}}{s_i^2} \quad \text{and} \quad \rho_\alpha = \sum_{i=1}^n \frac{m_i - t_i}{s_i^2} \sigma_{i,\text{corr}}^{(\alpha)}, \quad (26)$$

that determines the values of λ_β . The quantity:

$$d_i = \sum_{\alpha=1}^k \lambda_\alpha \sigma_{i,\text{corr}}^{(\alpha)} \quad (27)$$

in Eq. (23) can be interpreted as a shift caused by the correlated systematic uncertainties. Defining the shifted predictions as:

$$\bar{t}_i = t_i + d_i, \quad (28)$$

the χ^2 reads:

$$\chi^2 = \sum_{i=1}^n \left(\frac{m_i - \bar{t}_i}{s_i} \right)^2 + \sum_{\alpha=1}^k \lambda_\alpha^2. \quad (29)$$

Therefore, up to a penalty term given by the sum of the square of the nuisance parameters, the χ^2 takes the form of the uncorrelated definition. In order to achieve a visual assessment of the agreement between data and theory, it appears natural to compare the central experimental values m_i to the shifted theoretical predictions \bar{t}_i in units of the uncorrelated uncertainty s_i .

5 Effect of cuts on the χ^2

A relevant question to ask is what is the effect of possible cuts on the input dataset on the form of the χ^2 in the presence of correlations. In order to address this question, we consider a simple dataset made of two datapoints equipped with a single uncorrelated uncertainty and a single fully-correlated uncertainty:³

$$S : \quad \{m_i \pm \sigma_i^u \pm \sigma_i^c\}, \quad i = 1, 2. \quad (30)$$

The resulting covariance matrix reads:

$$\mathbf{V} = \begin{pmatrix} (\sigma_1^u)^2 + (\sigma_1^c)^2 & \sigma_1^c \sigma_2^c \\ \sigma_1^c \sigma_2^c & (\sigma_2^u)^2 + (\sigma_2^c)^2 \end{pmatrix}, \quad (31)$$

³Note that we need to include *also* an uncorrelated uncertainty because otherwise the covariance matrix would be singular. This is a symptom of the fact that if no uncorrelated uncertainties are present, the two points would be totally dependent on each other. This observation will be relevant in the following.

with inverse:

$$\mathbf{V}^{-1} = \frac{1}{(\sigma_1^u \sigma_2^u)^2 + (\sigma_1^c \sigma_2^u)^2 + (\sigma_1^u \sigma_2^c)^2} \begin{pmatrix} (\sigma_2^u)^2 + (\sigma_2^c)^2 & -\sigma_1^c \sigma_2^c \\ -\sigma_1^c \sigma_2^c & (\sigma_1^u)^2 + (\sigma_1^c)^2 \end{pmatrix}. \quad (32)$$

In addition, we have a set of two theoretical predictions $\{t_i\}$ that allows us to define a column vector of residuals $\mathbf{r}^T = (r_1, r_2)$ with $r_i = m_i - t_i$. The χ^2 then takes the following explicit expression:

$$\chi^2 = \mathbf{r}^T \cdot \mathbf{V}^{-1} \cdot \mathbf{r} = \frac{r_1^2 [(\sigma_2^u)^2 + (\sigma_2^c)^2] + r_2^2 [(\sigma_1^u)^2 + (\sigma_1^c)^2] - 2r_1 r_2 \sigma_1^c \sigma_2^c}{(\sigma_1^u \sigma_2^u)^2 + (\sigma_1^c \sigma_2^u)^2 + (\sigma_1^u \sigma_2^c)^2}. \quad (33)$$

Now, suppose we want to exclude the datapoint with $i = 2$ from the definition of the χ^2 . The possibly most natural way to proceed is to exclude the point, along with its uncertainties, directly from the set in Eq. (30). This straightforwardly leads to the following χ^2 :

$$\chi_{\text{cut } 1}^2 = \frac{r_1^2}{(\sigma_1^u)^2 + (\sigma_1^c)^2}. \quad (34)$$

However, this procedure might be questioned in that it artificially modifies the original dataset by effectively defining a *new* dataset where the point $i = 2$ is not present. In the absence of correlations this is a sound procedure because each single datapoint can be regarded as an independent subset of the original set. However, this is not the case when correlations are present.

Therefore, it is necessary to devise a method that avoids any modification of the dataset but yet allows one to prevent specific datapoints to contribute to the χ^2 . In general, the purpose of excluding datapoints from a fit is to avoid that the model used to compute the predictions is forced outside its application region. One can therefore *enforce* that the model exactly agrees with the datapoints to be excluded. This effectively amounts to set $t_2 = m_2$, or equivalently $r_2 = 0$, reducing Eq. (33) to:

$$\chi_{\text{cut } 2}^2 = \frac{r_1^2 [(\sigma_2^u)^2 + (\sigma_2^c)^2]}{(\sigma_1^u \sigma_2^u)^2 + (\sigma_1^c \sigma_2^u)^2 + (\sigma_1^u \sigma_2^c)^2}. \quad (35)$$

This has the clear advantage that the experimental information, including correlations, remains totally intact. In order to see that Eq. (35) has the right properties, we first take the limit for $\sigma_2^c \rightarrow 0$. This gives:

$$\lim_{\sigma_2^c \rightarrow 0} \chi_{\text{cut } 2}^2 = \frac{r_1^2}{(\sigma_1^u)^2 + (\sigma_1^c)^2}, \quad (36)$$

which reproduces the result of Eq. (34). This was to be expected because no reference to the point $i = 2$ can remain after the removal of its correlation to the point $i = 1$. In addition, it is also very instructive to take the limit $\sigma_2^u \rightarrow 0$ Eq. (35), which gives:

$$\lim_{\sigma_2^u \rightarrow 0} \chi_{\text{cut } 2}^2 = \frac{r_1^2}{(\sigma_1^u)^2}. \quad (37)$$

In this case, not only any reference to the point $i = 2$ drops out, but also the dependence of the χ^2 on the correlation uncertainty of the point $i = 1$ disappears. Despite at first sight this looks counterintuitive, it seems to be

the correct behaviour. To see this, we observe that, if the point $i = 2$ has no uncorrelated uncertainty, its fluctuations are totally driven by the fluctuations of the point $i = 1$. Therefore, the point $i = 2$ is completely dependent on the point $i = 1$ and thus cannot give any contribution to the χ^2 . In addition, the correlation uncertainty of the point $i = 1$ must also be irrelevant because, being $i = 2$ totally correlated to $i = 1$, any correlation of $i = 1$ to $i = 2$ translates into a correlation to itself and this cannot contribute to the χ^2 either. Importantly, if we take the limit $\sigma_1^u \rightarrow 0$, the χ^2 diverges no matter the value of σ_1^c . This is consistent with setting $\sigma_1^u = \sigma_2^u$ since the beginning in Eq. (30) which indeed gives a singular χ^2 because the covariance matrix is singular. Eq. (34) does not produce these features hanging in favour of Eq. (35).

6 Monte Carlo replica generation

In this section we consider the Monte Carlo generation of artificial replicas. The aim is to obtain a formula for the generation of replicas that gives rise to a figure of merit distributed correctly, *i.e.* that follows a χ^2 distribution with as many degrees of freedom as number of data points. Following Ref. [5], we require that the χ^2 , that in matricial form can be expressed as:

$$\chi^2 = \mathbf{y}^T \cdot \mathbf{V}^{-1} \cdot \mathbf{y}, \quad (38)$$

where \mathbf{V} is the $n \times n$ covariance matrix, being n the number of datapoints, and \mathbf{y} is a column vector with n entries that can be expressed as:

$$y_j = f_j - m_j, \quad (39)$$

with m_j a given value (to be thought as the central value) and f_j its fluctuated value, is distributed like:

$$\chi^2 \sim \sum_{i=1}^n z_i^2 = |\mathbf{z}|^2, \quad (40)$$

where z_1, \dots, z_n are independent, standard normal random variables, *i.e.* such that:

$$\langle z_i \rangle = 0 \quad \text{and} \quad \langle z_i z_j \rangle = \delta_{ij}, \quad (41)$$

where the symbol $\langle \dots \rangle$ indicates the average.⁴ This immediately tells us that:

$$\langle \chi^2 \rangle = n. \quad (44)$$

As already discussed above, Eq. (38) can alternatively be written as:

$$\chi^2 = |\mathbf{L}^{-1} \cdot \mathbf{y}|^2, \quad (45)$$

⁴The relevant identities are:

$$\langle z_i^{2n-1} \rangle = 0 \quad \text{and} \quad \langle z_i^{2n} \rangle = (2n-1)!! \quad n = 1, 2, 3, \dots, \quad (42)$$

where $(2n-1)!! = 1 \cdot 3 \cdot \dots \cdot (2n-1)$ is the odd semi-factorial. When considering combinations of two random variables the result is:

$$\langle z_i^n z_j^m \rangle = \begin{cases} \langle z_i^{n+m} \rangle & i = j \\ \langle z_i^n \rangle \langle z_j^m \rangle & i \neq j \end{cases} \quad (43)$$

where \mathbf{L} is the Cholesky decomposition of the matrix \mathbf{V} , such that:

$$\mathbf{V} = \mathbf{L} \cdot \mathbf{L}^T. \quad (46)$$

Comparing Eqs. (45) and (40), one can infer that the following relation should hold:

$$\mathbf{L}^{-1} \cdot \mathbf{y} \sim \mathbf{z}, \quad (47)$$

which immediately implies:

$$\mathbf{y} \sim \mathbf{L} \cdot \mathbf{z}, \quad (48)$$

or equivalently:

$$f_j \sim m_j + \sum_{i=1}^n L_{ji} z_i. \quad (49)$$

In conclusion, in order for the χ^2 to have the correct distribution, the fluctuations f_j around the experimental central values m_j of a given Monte Carlo replica have to be computed according to Eq. (49).

We now show that Eq. (49) reproduces the statistical properties of the original ensemble of data. To do so, we compute average and correlations of Eq. (49) using Eq. (41). The average gives:

$$\langle f_j \rangle = m_j + \sum_{i=1}^n L_{ji} \langle z_i \rangle = m_j, \quad (50)$$

as expected. The correlation instead gives:

$$\begin{aligned} \langle f_j f_k \rangle &= \left\langle \left(m_j + \sum_{i=1}^n L_{ji} z_i \right) \left(m_k + \sum_{i'=1}^n L_{ki'} z_{i'} \right) \right\rangle \\ &= m_j m_k + \sum_{i,i'=1}^n L_{ji'} L_{ki'} \langle z_i z_{i'} \rangle = \langle f_j \rangle \langle f_k \rangle + V_{jk}, \end{aligned} \quad (51)$$

again as expected. These are fundamental requirements of a sound Monte Carlo ensemble (see also Eq. (10) of Ref. [8]).

Eq. (49) should be compared to the more “standard” Monte-Carlo replica generation formulas such as those given in Refs. [6, 7]. We now specifically consider Eq. (20) of Ref. [7] that we report here in terms of relative uncertainties δ ’s matching our notation:

$$\tilde{f}_j \sim m_j \prod_{m=1}^{k_m} \left(1 + \delta_{\text{mult},j}^{(m)} x_m \right) \left(1 + \delta_{\text{unc},j} z_j + \sum_{a=1}^{k_a} \delta_{\text{add},j}^{(a)} y_a \right), \quad (52)$$

where z_j , y_a , and x_m are independent random variables. We now expand the product in the front up to second order in x

$$\prod_{m=1}^{k_m} \left(1 + \delta_{\text{mult},j}^{(m)} x_m \right) = 1 + \sum_{m=1}^{k_m} \delta_{\text{mult},j}^{(m)} x_m + \sum_{m=1}^{k_m} \sum_{\substack{m'=1 \\ m' \neq m}}^{k_m} \delta_{\text{mult},j}^{(m)} \delta_{\text{mult},j}^{(m')} x_m x_{m'} + \text{trilinear}, \quad (53)$$

where the “trilinear” corrections correspond to combinations of the kind $x_i x_j x_k$, with $i \neq j \neq k \neq i$, such that their average vanishes: $\langle x_i x_j x_k \rangle = \langle x_i \rangle \langle x_j \rangle \langle x_k \rangle = 0$. This allows us to expand Eq. (52) as:

$$\begin{aligned}
\tilde{f}_j &\sim m_j \left(1 + \delta_{\text{unc},j} z_j + \sum_{a=1}^{k_a} \delta_{\text{add},j}^{(a)} y_a + \sum_{m=1}^{k_m} \delta_{\text{mult},j}^{(m)} x_m \right. \\
&+ \delta_{\text{unc},j} \sum_{m=1}^{k_m} \delta_{\text{mult},j}^{(m)} z_j x_m + \sum_{a=1}^{k_a} \sum_{m=1}^{k_m} \delta_{\text{add},j}^{(a)} \delta_{\text{mult},j}^{(m)} y_a x_m \\
&\left. + \sum_{m=1}^{k_m} \sum_{\substack{m'=1 \\ m' \neq m}}^{k_m} \delta_{\text{mult},j}^{(m)} \delta_{\text{mult},j}^{(m')} x_m x_{m'} \right) + \text{trilinear} .
\end{aligned} \tag{54}$$

Since Eq. (54) involves only terms that have at most one power of the same random variable, the average of \tilde{f}_j automatically yields:

$$\langle \tilde{f}_j \rangle = m_j , \tag{55}$$

in agreement with the expectation. We now compute the correlations making use of the combination rules of random variables. This gives:

$$\begin{aligned}
\langle \tilde{f}_j \tilde{f}_k \rangle &\sim m_j m_k \left(1 + \delta_{\text{unc},j}^2 \delta_{jk} + \sum_{a=1}^{k_a} \delta_{\text{add},j}^{(a)} \delta_{\text{add},k}^{(a)} + \sum_{m=1}^{k_m} \delta_{\text{mult},j}^{(m)} \delta_{\text{mult},k}^{(m)} \right. \\
&+ \delta_{\text{unc},j}^2 \sum_{m=1}^{k_m} \delta_{\text{mult},j}^{(m)} \delta_{\text{mult},k}^{(m)} + \sum_{a=1}^{k_a} \sum_{m=1}^{k_m} \delta_{\text{add},j}^{(a)} \delta_{\text{add},k}^{(a)} \delta_{\text{mult},j}^{(m)} \delta_{\text{mult},k}^{(m)} \\
&\left. + \sum_{m=1}^{k_m} \sum_{\substack{m'=1 \\ m' \neq m}}^{k_m} \delta_{\text{mult},j}^{(m)} \delta_{\text{mult},k}^{(m)} \delta_{\text{mult},j}^{(m')} \delta_{\text{mult},k}^{(m')} + \mathcal{O}(\delta^6) \right) \\
&= \langle \tilde{f}_j \rangle \langle \tilde{f}_k \rangle + V_{jk} + \mathcal{O}(\sigma_{\text{mult}}^2 s^2, \sigma_{\text{mult}}^2 \sigma_{\text{add}}^2) .
\end{aligned} \tag{56}$$

In conclusion, Eq. (52) reproduces exactly the central values of the original ensemble (see Eq. (55)) while it reproduces correlations only up to corrections proportional to the fourth power of the typical uncertainty. Interestingly, these corrections are all proportional to σ_{mult}^2 . This means that, in the absence of multiplicative uncertainties, somewhat expectedly Eq. (52) *exactly* reproduces the correlations. On the other hand, if multiplicative uncertainties are large these correction become potentially sizeable preventing Eq. (52) from correctly reproducing the correlations of the original ensemble.

The case of Eq. (13) of Ref. [6] is more involved because of the presence of the square-root sign for the relative multiplicative uncertainties. This case can be handled again by expanding to second order the factor depending on multiplicative uncertainties with the only difference that, due to the expansion:

$$\sqrt{1+x} = 1 + \frac{1}{2}x - \frac{1}{8}x^2 + \mathcal{O}(x^3) , \tag{57}$$

the square root produces a factor $1/2$ in front of the linear terms and a factor $-1/8$ in front of the quadratic terms. This is expected to reduce the impact of the corrections terms in Eq. (56).

It is instructive to work out whether Eq. (49) can be directly related to the Monte-Carlo replicas generation formulas given in Refs. [6, 7]. The aim is to find a closed-form expression for the Cholesky-decomposition matrix \mathbf{L} of the covariance matrix \mathbf{V} in terms of the single uncertainties. Through Eq. (49), this will finally allow us to derive a formula of the same kind of those given in Refs. [6, 7]. The basic observation is that the covariance matrix can be written as:

$$\mathbf{V} = \mathbf{D} + \sum_{l=1}^k \mathbf{V}^{(l)} = \sum_{l=0}^k \mathbf{L}^{(l)} \cdot \mathbf{L}^{(l)T}, \quad (58)$$

with:

$$L_{ij}^{(l)} = \begin{cases} s_i \delta_{ij} & l = 0, \\ \sigma_i^{(l)} \delta_{j1} & 1 \leq l \leq k. \end{cases} \quad (59)$$

On the other hand, we also know that:

$$\mathbf{V} = \mathbf{L} \cdot \mathbf{L}^T, \quad (60)$$

therefore:

$$\mathbf{L} \cdot \mathbf{L}^T = \sum_{l=0}^k \mathbf{L}^{(l)} \cdot \mathbf{L}^{(l)T}. \quad (61)$$

The question is to express \mathbf{L} in terms of all the $\mathbf{L}^{(l)}$. Unfortunately though this equation has no general solution. However, a step forward can be done by exploiting the fact that the matrices $\mathbf{V}^{(l)}$ are positive semidefinite (*i.e.* they have zero determinant) and thus their Cholesky decomposition is not unique. In particular, provided that $k \leq n$, one can choose:

$$\tilde{L}_{ij}^{(l)} = \sigma_i^{(l)} \delta_{jl}, \quad 1 \leq l \leq k \leq n, \quad (62)$$

such that:

$$\left(\sum_{l=1}^k \tilde{\mathbf{L}}^{(l)} \right) \cdot \left(\sum_{l'=1}^k \tilde{\mathbf{L}}^{(l')} \right) = \sum_{l=1}^k \mathbf{V}^{(l)}. \quad (63)$$

However, choosing:

$$\tilde{\mathbf{L}} = \mathbf{L}^{(0)} + \sum_{l=1}^k \tilde{\mathbf{L}}^{(l)}, \quad (64)$$

and multiplying it for its transpose would give:

$$\tilde{\mathbf{L}} \cdot \tilde{\mathbf{L}}^T = \mathbf{V} + \sum_{l=1}^k \left(\mathbf{L}^{(0)} \cdot \tilde{\mathbf{L}}^{(l)T} + \tilde{\mathbf{L}}^{(l)} \cdot \mathbf{L}^{(0)T} \right), \quad (65)$$

that gives back the covariance matrix plus a spurious term. Writing the expression above in terms of single entries, one finds:

$$\sum_{m=1}^n \tilde{L}_{im} \tilde{L}_{jm} = V_{ij} + s_i \sigma_j^{(i)} + s_j \sigma_i^{(j)}. \quad (66)$$

The peculiarity here is that the upper index of $\sigma_j^{(i)}$ and $\sigma_i^{(j)}$, since $k \leq n$ (and in fact usually $k \ll n$, *i.e.* the number of correlated uncertainties is typically much smaller than the number of data points), exceeds the allowed range. To fix this problem we just assume $\sigma_j^{(i)} = 0$ for $i > k$. This finally gives:

$$\sum_{m=1}^n \tilde{L}_{im} \tilde{L}_{jm} = V_{ij}, \quad \text{for } i, j > k, \quad (67)$$

that means that, by adopting the Cholesky decomposition in Eq. (62), part of the covariance matrix (usually the largest part) is exactly recovered. Therefore, $\tilde{\mathbf{L}}$ Eq. (64) seems to be a generally good approximator to the real Cholesky decomposition of the covariance matrix \mathbf{V} . Plugging this equation into Eq. (49) immediately gives:

$$f_j \sim m_j + z_j s_j + \sum_{l=1}^k z_l \sigma_{j,\text{corr}}^{(l)}. \quad (68)$$

It is interesting to focus on similarities and differences with the formulas present in the literature. We will specifically consider Eq. (13) of Ref. [6] and Eq. (20) of Ref. [7].

We start by considering the case in which there are only uncorrelated and correlated additive uncertainties. In this case, Refs. [6, 7] are in mutual agreement and only in partial agreement with our Eq. (68), the difference being that the random numbers used to generate the fluctuation associated to the k correlated uncertainties are equal to the first k of the n used for the uncorrelated uncertainties (remember that we required $k \leq n$). Of course, if $k \ll n$, as it often happens, the agreement between Eq. (68) and Refs. [6, 7] is asymptotically good.

When considering also correlated multiplicative uncertainties there are differences across formulas of Refs. [6, 7] and Eq. (68). While Eq. (68) treats multiplicative uncertainties on the same footing as additive uncertainties (consistently with the construction of the covariance matrix in Eq. (2)), Refs. [6, 7] introduce a multiplicative factor. In addition, Ref. [6] distinguishes between relative and absolute multiplicative uncertainties.

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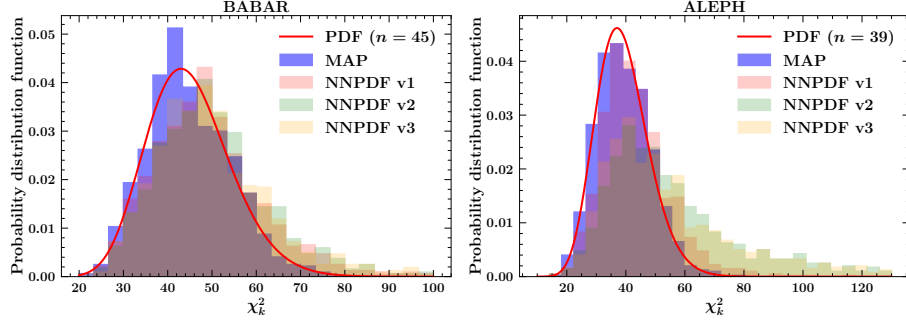


Figure 1: Distribution of the χ^2 's of the Monte Carlo replica w.r.t. to the corresponding central values for BABAR (left) and ALEPH (right) experiments compared to the probability distribution function with to the appropriate number of degrees of freedom given by the number of data points. The blue histograms are produced using Eq. (49), the red histogram is produced using Eq. (13) of Ref. [6] where multiplicative uncertainties are treated as relative, while the green histogram is produced using Eq. (20) of Ref. [7].

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