

Automatic Synchronization of Drift-Time Spectra and Maximum Drift-Time Measurement of an MDT

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Abstract

A method for the reliable automatic synchronization of the drift-time spectra of ATLAS monitored drift-tubes and for the determination of the maximum drift-time is presented. The Fermi function for the measurement of the maximum drift time has been adapted to the particular shape of the drift-time spectrum characteristic for Ar:CO₂(93:7) at 3 *bar* which is used in the ATLAS muon chambers.

1 Introduction

The monitored drift tubes used in the ATLAS muon spectrometer measure the time interval τ_d between the passage of a muon through the tube and the instance of time when the electrons which are the ionization product of muon-atom collisions in the gas arrive at the anode wire of the tube. This drift time τ_d is a strictly monotone function of the distance r of the muon track from the wire. In practice, τ_d is not measured directly, but the time interval

$$\tau_m := t_\mu + \tau_d + \tau_{\mu,electronics} - t_{trigg}, \quad (1)$$

where t_μ is the time when the muon hits the tube, t_{trigg} the time when the trigger gives the start pulse of the TDC connected to the tube, and $\tau_{\mu,electronics}$ is the propagation time of the MDT signal. The term $t_\mu + \tau_{\mu,electronics} - t_{trigg}$ is the same for all events. It can, however, slightly differ from tube to tube. It defines the position of the drift-time spectrum of the tube with respect to the event time.

If the temperature, the pressure, and the gas mixture are the same in all drift tubes of a muon chamber, they have the same r-t relation after their drift time

spectra have been synchronized. In the present article, we describe an algorithm which allows a *reliable automatic* synchronization of the spectra. We will use the leading edge of the drift time spectrum to match the spectra. Our algorithm also enables us to measure the length of the drift time spectra. The length of the spectrum is a very useful quantity, as differences in the lengths from tube to tube indicate a different operational behavior of tubes.

2 Phenomenological Parameterization of the Rising and Falling Edge of the Drift-Time Spectrum

Figure 1 shows the typical drift-time spectrum of a uniformly illuminated MDT of 3 *cm* diameter filled with Ar:CO₂(93:7) at 3 *bar* absolute pressure and 3080 *V* operating voltage. The data presented in figure 1 were obtained with the MPI BOS prototype chamber in the CERN H8 beam in September 1999. At zero drift

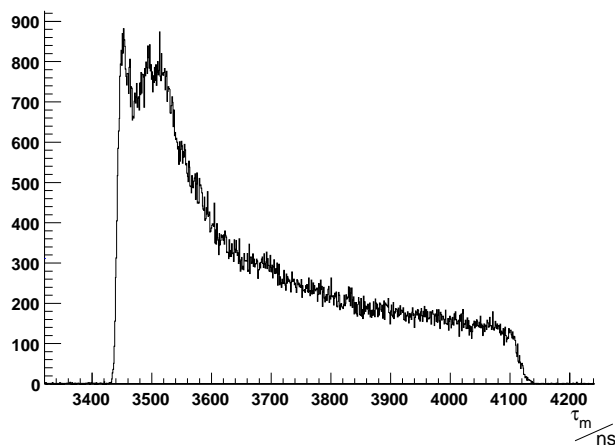


Figure 1: Drift-time spectrum of a drift tube filled with Ar:CO₂(93:7) at 3 bar absolute pressure.

time we see a sharp rise and, at 700 *ns* drift time, a fall-off, which is less steep than the rise because of the inefficiency of the tube at its wall.

We will use the step at the beginning of the drift time spectrum to synchronize the drift time spectra in a chamber. For a tube with perfect resolution the rising edge of the spectrum would have infinite slope and the initial part of the spectrum would be a step function. The resolution of the drift tube makes this rise finite. This smeared out step can quite well be described by a Fermi function F ,

$$F(t) := \frac{A_0}{1 + e^{-\frac{t-t_0}{T_0}}}; \quad (2)$$

herein A_0 is the maximum value F achieves for $t \rightarrow \infty$, $F(t_0) = \frac{A_0}{2}$, and the "temperature" T_0 corresponds to the slope of the edge. The slope at $t = t_0$ is $\frac{A_0}{4T_0}$.

In practice, due to accidental hits, one can find entries for $t \ll t_0$ which the Fermi function does not describe. Therefore, on the left hand side of t_0 , it looks as if the well-known drift-time spectrum sits on a plateau of height p_0 . In this region, the Fermi function F has to be replaced by the modified function

$$G(t) := p_0 + F(t) = p_0 + \frac{A_0}{1 + e^{-\frac{t-t_0}{T_0}}}. \quad (3)$$

In order to synchronize the drift-time spectra of the tubes in a muon chamber one has to fit this function G to the leading of the spectrum of each individual tube and shift it by $-t_0$.

On first sight, one would expect the function G to fit the falling edge of the drift-time spectrum as well with different values for p_0 , A_0 , and T_0 . In case of Ar:CO₂, however, the drift-time distribution is not constant on the left hand side of the falling edge, but approximately linearly decreasing. Hence, a more appropriate parameterization of the falling edge is

$$H(t) := p_m + \frac{\alpha_m t + A_m}{1 + e^{\frac{t-t_m}{T_m}}}. \quad (4)$$

This function approximates for $t \ll t_m$ a line with the intercept A_m and the slope α_m . Figure 2 shows a characteristic fit of G and H to the drift-time spectrum.

3 Algorithm

As the functions G and H only describe the rising and the falling edge of the spectrum, one cannot fit these functions to the complete spectrum. In principal there are three methods which circumvent this problem.

The first method is to describe the area in between the edges by a polynomial as described in [1]. The major drawback of this method is that it introduces extra parameters, which are of no interest, into the fit. As these parameters are correlated to the values of t_0 and t_m it can affect these values.

In the second approach, one selects the area of the spectrum where G and H describe its shape correctly. However, our experience is that even after a proper choice of these areas the fit can fail.

The method which is described here is to select an area for each of the parameters p_0 , A_0 , p_m , α_m , and A_m , and calculate them analytically. The selection is performed by an automatic pattern recognition. This pattern recognition also returns the areas for the fits. In the fit the parameters mentioned above are fixed to the calculated values. This guarantees a fast and reliable convergence of the fit.

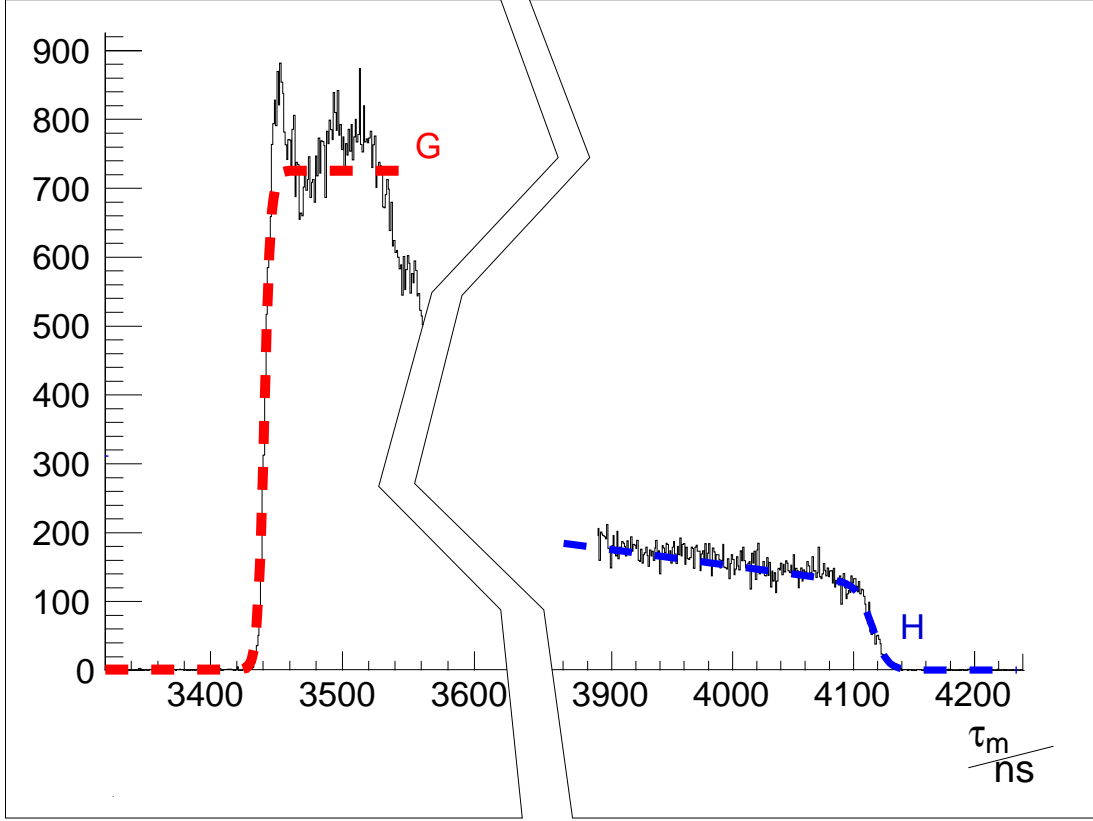


Figure 2: Typical fits of the leading and falling edge of the drift-time spectrum

3.1 Determination of the Fit Area and Calculation of the fixed Parameters

The areas which have to be selected are shown in figure 3. The fitted functions are approximately flat in the p_0 , A_0 , and p_m region and the function values are approximately the corresponding parameters. Therefore the mean of the bin contents in the corresponding regions is the best value for the parameters. In the region used for the determination of the parameters A_m and α_m the fitted function approximates a line with these two parameters as its intersection and slope. To determine these parameters one puts a line to the spectrum in this region.

The fits are performed inside the areas marked with “ t_0 fit” and “ t_{max} fit”.

3.1.1 The Variable Bin Width Histogram (VBH)

The pattern recognition is based on the concept of bins with equal content [3]. We will refer to that by the term “variable bin width histogram(VBH)”. This histogram has a fixed number of entries (bin content) in each histogram bin. The

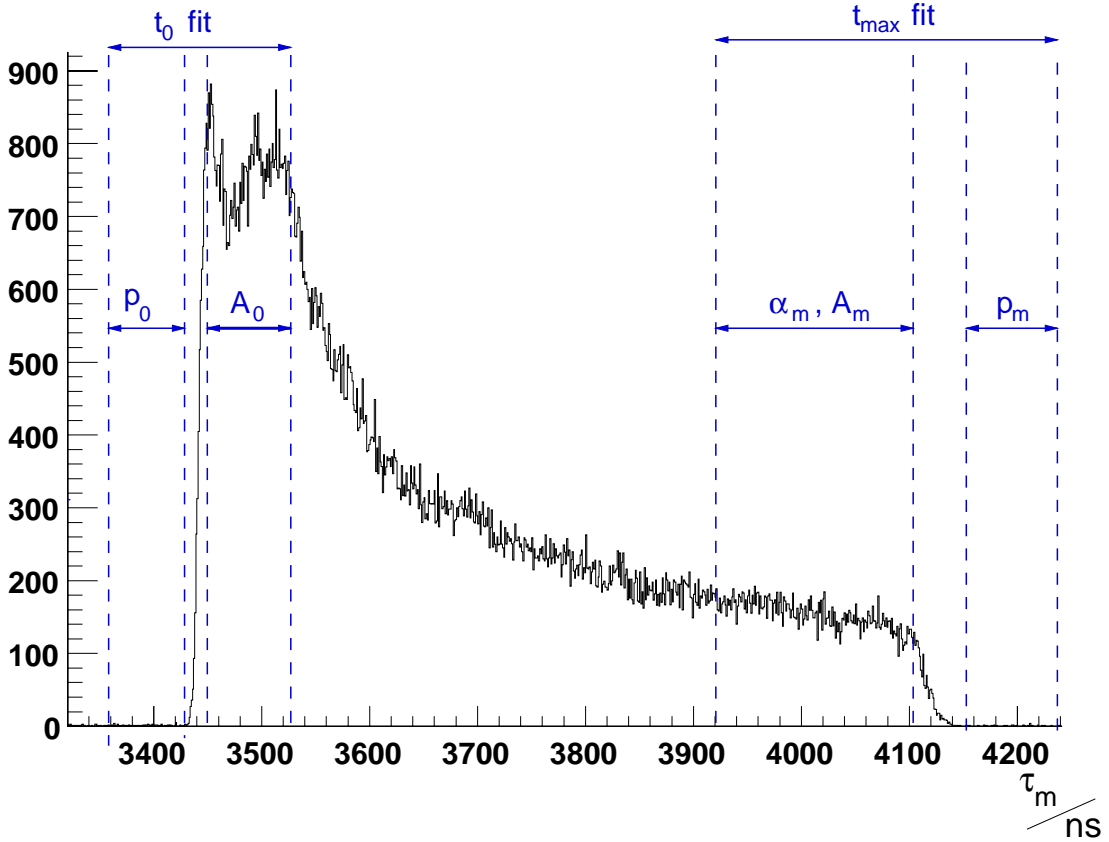


Figure 3: Relevant areas of the spectrum

density of the data is represented by the width of the bin. Wide bins correspond to a low density and small bins to a high density.

It is known, that in a conventional histogram the binning should be chosen such that each bin has a certain minimum number of entries per histogram bin. A binning that is too fine, leads to a high statistical fluctuation for the number of entries in each bin, and a binning that is too coarse results in a bad resolution. The VBH performs the task of finding the ideal binwidth automatically, and for each region of the histogram separately.

As the TDC-data comes already binned due to the digitization, and as one wants a unique assignment of a hit to a VBH-bin, the bin content has to be set higher than the biggest number of entries in a TDC-bin. The pattern recognition relies on the bin content being two times the content of the highest TDC-bin.

As the background is hopefully low, the first bin tends to cover not only the complete p_0 area, but also part of the rising edge of the spectrum. This would prevent the pattern recognition from finding the p_0 area. The same goes for the p_m area. Therefore we impose a maximum bin width. This setting is one of the values which has to be tuned to the properties of the detector gas, the tube

diameter, and the TDC-digitization. In our case we use a maximum bin width of 20 ns for the t_0 fit and 40 ns for the t_{max} fit.

For our one-dimensional problem the creation of this histogram is quite simple. The TDC entries are sorted by the time value. The start of the lowest VBH bin is the lowest TDC value. One walks through the sorted hits. The algorithm for creating the VBH is shown in figure 4.

In figure 5 the areas which are important for the pattern recognition are shown magnified. In addition to the spectrum in TDC-binning the bin width of the VBH histogram is shown.

3.1.2 Selection of the Areas

In the upper left magnification in figure 5 one can see that the start of the slope lies in the last VBH bin which has the maximum bin width. The left border of this bin is the right border of the p_0 area.

In the plateau region, the bin width of the VBH fluctuates between two values. This is due to the fact that the TDC-data are binned. As we have set the bin content to two times the highest TDC bin, the bin width is either two times or three times the TDC bin width. The principle of the A_0 area selection is to select the smallest region which contains all bins which have the smallest bin width. This can lead to problems if the spectrum has spikes. A small modification of this concept makes our algorithm more robust. We sort the VBH bins by size, starting with the smallest. We memorize the width of the fifth of these sorted VBH bin. The A_0 area is the smallest area which contains all VBH bins of this size. If the spectrum has no unusual spikes, the memorized width is the smallest bin width, otherwise the spike leads to smaller bins.

In the lower magnification in figure 5 one can see that, after the falling edge, all VBH bins have the maximum bin width. We search for the first VBH-bin which comes after the rising slope of the spectrum, which has the maximum bin width and whose right neighbour also has the maximum bin width.

The right border of this bin is the start of the p_m area. As the background in the left of the spectrum is flat, the p_m region extends to the end of the TDC range.

For the α_m - p_m area we need the part of the spectrum which decreases approximately linearly. As this area is quite large, the selection can be made quite coarse. The right border is simply set to a fixed distance from the left boarder of the p_m area (usually 40 ns). Its width is also set to a fixed length (400 ns is a good value for Ar:C0₂).

3.1.3 Calculation of the Parameters

The parameters p_0 , A_0 , and p_m are the mean of the TDC-bin-content over the corresponding areas. The parameters α_m and A_m are the result of a linear regression

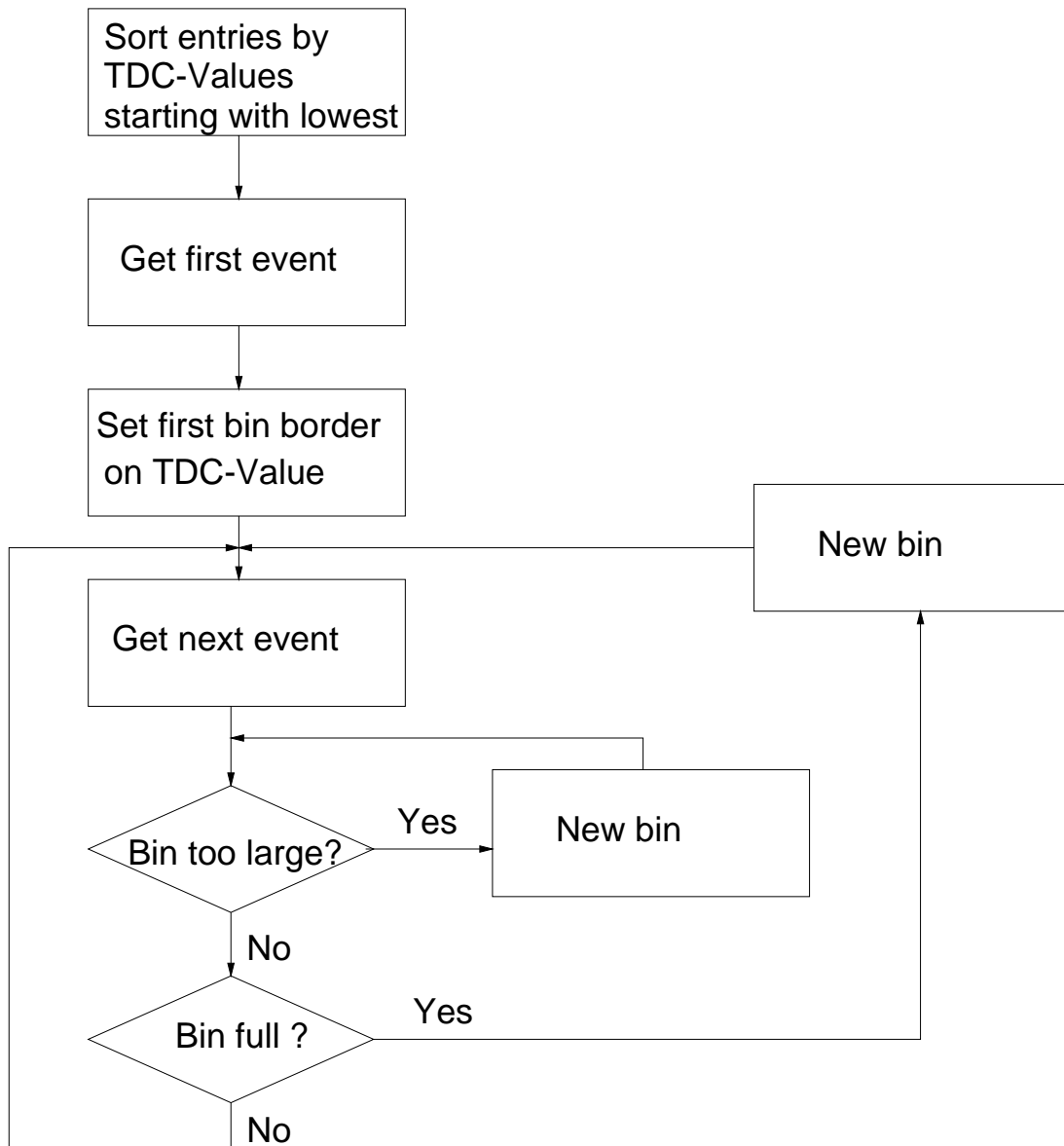


Figure 4: Filling the VBH

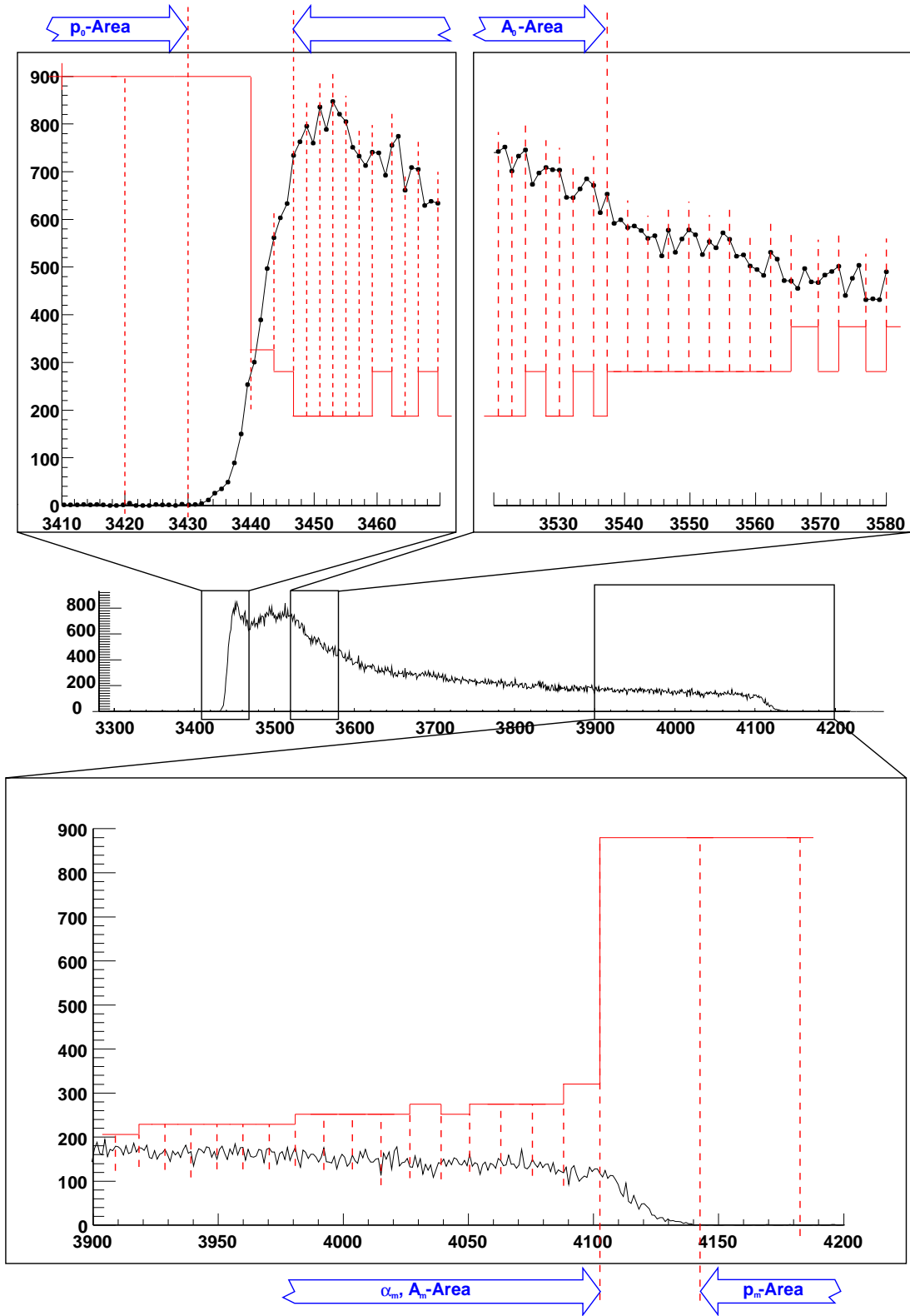


Figure 5: The important areas of the spectrum and the binning of the VBH

over the selected region.

3.2 Fitting

The area for the t_0 fit consists of the p_0 area, the A_0 area, and the intermediate region. The combination of the α_m - A_m area, the p_m area and the region in between these is the t_{max} fit area.

As a start value for t_0 and t_m the centers of the regions between the p_0/p_m and the $A_0/\alpha_m, A_m$ regions are used. The analytically calculated parameters are fixed. The fit is performed with a fitting package like MINUIT.

For the fit one does not use the VBH histogram, but the original TDC data. This has two reasons. First the maximum bin width does not represent the true density of hits in regions with a low density and second the histogram in TDC binning contains the maximum amount of information.

4 Test of the Pattern Recognition and Fitting Procedure

We tested the fit in about 250 cases from Monte Carlo and 50 cases from the MPI BOS prototype chamber acquired at the 1999 test beam in H8. These fits never failed unless the presence of the spectrum could not be seen by eye.

For a more quantitative study we created a big Monte-Carlo sample with 2080000 hits in one tube. We split these hits in two ways. First we made 52 subsamples with 40000 hits each and second we made 13 subsamples each containing 160000 hits. We let the algorithm run over all these subsamples.

The results for the quantity t_0 and the quantity $t_0 - 2T_0$ are shown in figure 6. Here one can also see that the fit never failed as all results were in a $\pm 3\sigma$ region around the mean of the distribution.

Looking at the RMS of the results one notices that, for 40000 events the fluctuation of $t_0 - 2T_0$ is a little bit and for 160000 events per fit considerably smaller than t_0 . If one looks at the behavior of these quantities for an increasing statistics one gets the following:

$$\begin{aligned} \frac{\sigma_{t_0, 40000 \text{ events}}}{\sigma_{t_0, 160000 \text{ events}}} &= \frac{0.44}{0.32} = 1.4, \\ \frac{\sigma_{t_0 - 2T_0, 40000 \text{ events}}}{\sigma_{t_0 - 2T_0, 160000 \text{ events}}} &= \frac{0.42}{0.20} = 2.1 \end{aligned} \tag{5}$$

So $t_0 - 2T_0$ follows the expected $\frac{1}{\sqrt{N}}$ statistical scaling rule, but not t_0 .

The reason for this seems to be an uncertainty in the determination of A_0 . As the corresponding region is not completely flat, and the spectrum drops at both sides of the region, the value strongly depends on the selection of the region. This selection has a low precision.

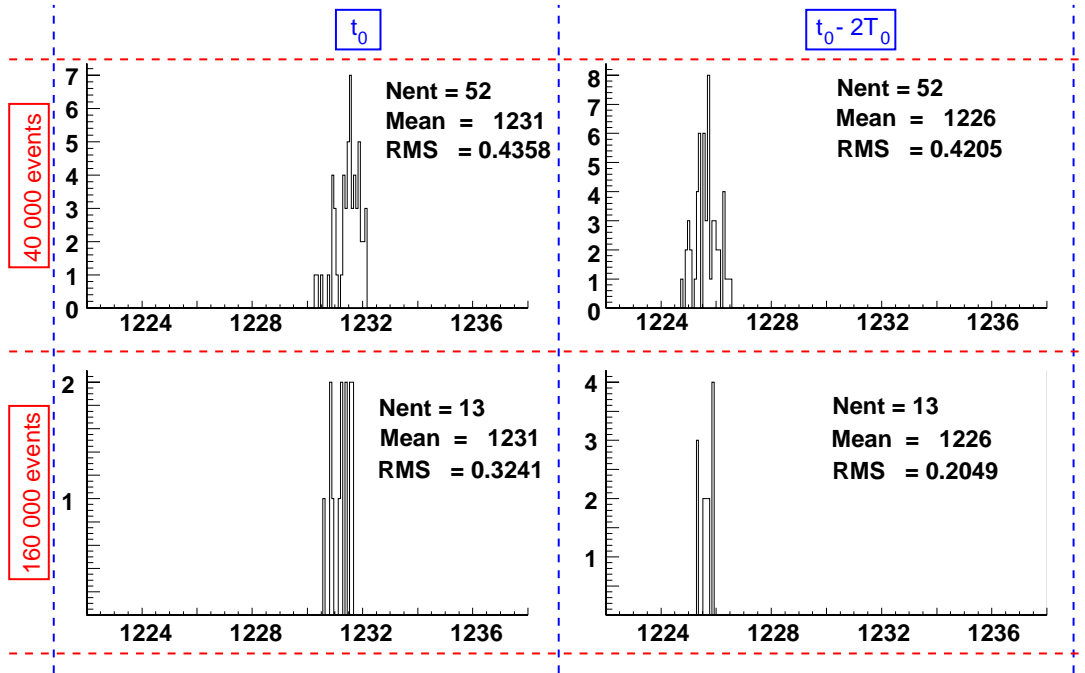


Figure 6: Test results for the t_0 fit

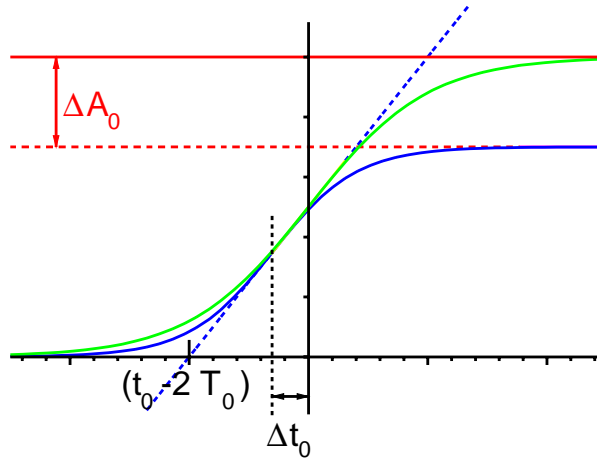


Figure 7: Correlations in the Fermi function

In figure 7 one sees how A_0 influences t_0 . This figure shows the function G with two different sets of parameters. They both approximate the same rising edge, but A_0 is shifted. As t_0 is the value, where

$$G(t = t_0) - p_0 = \frac{1}{2}A_0, \quad (6)$$

t_0 will be shifted. But the tangent of G at $t = t_0$ hardly changes. The point where this line crosses the background level is at $t = t_0 - 2T_0$. Therefore t_0 is strongly correlated to A_0 , while $t_0 - 2T_0$ is not. As A_0 is an unstable parameter, $t_0 - 2T_0$ is more stable than t_0 for high statistics.

As one has less hits with a high drift time, the values for t_m have a higher RMS. The value corresponding to $t_0 - 2T_0$ in the t_{max} fit is $t_m + 2T_m$. As the slope at the end of the spectrum is not well defined due to the low statistic it is not useful to use this value for the synchronisation of drift-time spectra.

5 Summary

The use of a pattern recognition using variable bin width histograms not only makes an automatic selection of the fit regions, but enables us to calculate some of the fit parameters analytically. This results in a stable and fully automatic method of synchronizing the spectra and measuring the maximum drift time.

The parameters of the pattern recognition which have to be set, namely the maximum bin width and the widths of some areas, may depend on the TDC properties and the background noise, but are the same for all tubes. So these parameters have to be set only once for a given setup.

For high statistics (more than 40000 hits) the value $t_0 - 2T_0$ is more stable than t_0 and therefore more useful for the synchronisation of the spectra.

Acknowledgement

We would like to thank A. Staude for the fruitful discussion we had with him on our algorithm.

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