

Step by step TWIM Calibration

7. April 2022

0.1 Overview

For this calibration I used the data of Experiment S455 in March 2021, Run 273, subruns 1-48.

1. Alignment of the energy per Anode for each Section
2. Alignment of the energy per Section
3. Drift time Calibration

Info: I always count from 0 to 15.

0.2 Alignment of the energy per Anode for each Section

For the Energy, you should first align all the gain per anode by plotting for each section:

$E_{raw}[anode\ i]$ vs $E_{raw}[anode\ ref]$.

anode ref = the 5th anode

I plot these 2D histos only for events where the 16 anodes per section have seen an ion. (no specific tpat selection needed).

0.2.1 Computing

First run the program "small_script_hist.C" for all subruns. Then use "hadd" to add up the .root files. The combined .root file can then be used for the scrip called "retrieve_fits_hist.C ". This one makes nice canvases for the plots $anode[i]$ vs $anode_ref$ and stores the fit parameters under `parameters_twim_anodes.csv`.

In this directory you find the `parameters_twim_anodes.csv` I retrieved from this first calibration step. The offset (=gain) should be near to 1.

Design of `parameters_twim_anodes.csv`

It stores `section(s)`, `anodenr(i)`, `slope`, `offset`

$E_sum_ref * slope[i][s] + offset[i][s]$

To calibrate you have to do it the other way round:

$E_cal_anode[i][s] = E_anode[i][s] / slope[i][s] - offset[i][s] / slope[i][s]$

0.3 Alignment of the energy per Section

- you should select event where the ions loss their energy in one section only
- then you should select a limited range in ToF (100 ps range)
- then you should calculate for each section $E_{sum}[s]$, the sum of the 16 anodes (using $E_cal_anode[i][s]$ from the previous step).
- you should see a small shift between the four sections
- correct from this shift by pol1
- as result you get: $E_{al_step2final}[s][a] = OffsetPerSection[s] + GainPerSection[s] * E_cal_anode[i][s]$

0.3.1 Computing

Run the macro "twim_sum_energy.C" using all subruns as input parameter. As output you get a .root file with 1D histos with the summed TWIM energy for each section. Use this output .root file as input file for the macro "e_sum_cal.C". This uses TSpectra etc. to (linearly) calibrate the E_sum energy for all sections. The according fit parameters are stored in the parameter file "sum_anodes_parameters.csv".

Now you can use the macro "twim_final_cal.C" (as input parameter the name of the subruns). This macro uses both parameter files "parameters_twim_anodes.csv" (anode fit) and "sum_anodes_parameters.csv" (summed energy fit). Now E_sum is calibrated.

0.4 Drift time Calibration

After this calibration you can extract the angle before GLAD from TWIM Music. From the signal of Music we do not get directly position information, but timing. Using the information from MW1 and MW2 the position on each anode can be extrapolated (2X and 2Y positions are needed in MW1 and MW2. For the precise reconstruction see macro `rft_analysis_tref.C` and `mw_position.C`):

$$X_{al} = FF_slope * anode_pos + FF_offset$$

This has to be done for each fission fragment (FF).

For this calibration only selected FF combinations are used (no tpat selection:

- subcase 2X2Y-D: one left down and one right down
- subcase 2X2Y-U: one left up and one right up
- subsubcase 2X2Y-LD-RU: one left down and one right up
- subsubcase 2X2Y-LU-RD: one left up and one right down
- subsubcase 2X2Y-L: one left down and one left up
- subsubcase 2X2Y-R: one right down and one right up

Important to mention, I only select events where each anode has exactly one value assigned, not more. All sections have a reference anode. The time for the anode is calculated:

$$t_drift[s][a] = time[s][a] - time_ref[s]$$

Plotting X versus drift time, the drift velocity on each anode can be extracted. X is the position of the ion in front of each anode, extrapolated from X data obtained with MW1 and MW2.

From this you retrieve X_{cal} for each anode of each section:

$$X_{al}[s][a] = X_OffsetPerAnodePerSection[s][a] + DV_PerAnodePerSection[s][a] * DT_{raw}[s][a] \quad (\text{where } DV_PerAnodePerSection \text{ is the drift velocity, the theoretical value should be } 0.056 \text{ mm/ns})$$

0.4.1 Δx correction

As next step the aligned positions from all 16 anodes ($X_{al}[s][a]$) for one section can be fitted for each event. For the fit I used anode 1-10 (see figure 1).

From this linear fit you get $X_{fit}[s][a]$.

Then you plot $\Delta X[s][a] = X_{fit}[s][a] - X_{al}[s][a]$ versus $X_{al}[s][a]$.

Here you will see that it is not flat. This is normal because there are non-linearity due the cathode effect and attachment. I fit this using a TSpline.

$$X_{cal} = X_{al} + TSpline(X_{al})$$

After correction with this TSpline check that $\Delta X = (X_{fit} - X_{cal})$ versus X_{cal} is horizontal. The width of the projection of this 2D-plot on the Y axis will give you the position resolution per anode. I have something around 20 micrometer sigma.

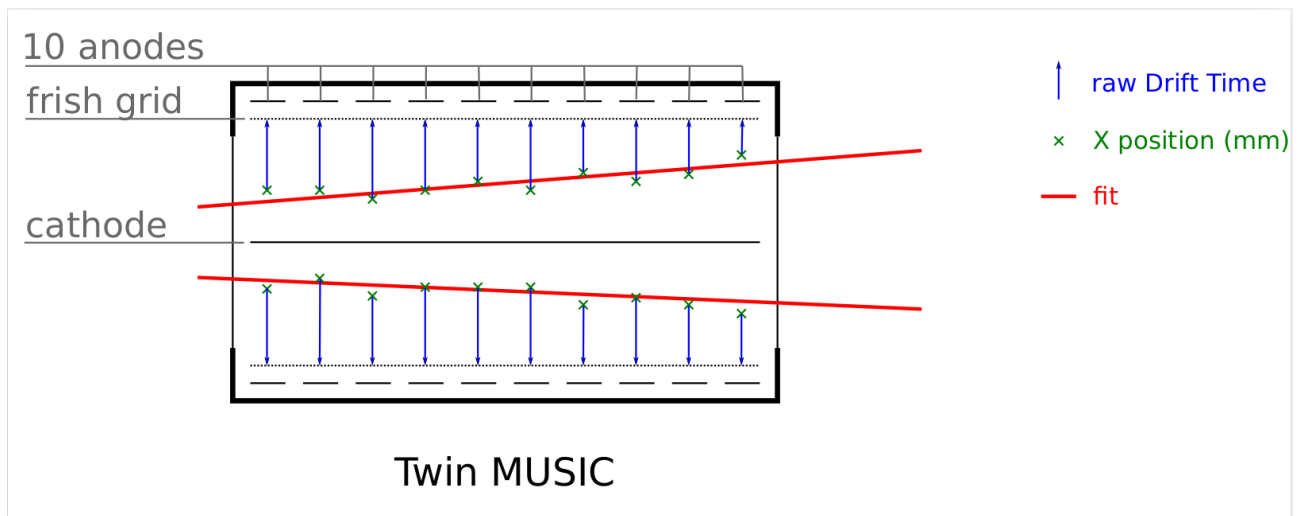


Abbildung 1: Calibrated positions (green) vs. fit (red). The fit is done using anode 1-10.

0.4.2 Computing

Use the macro "drift_analysis_tref.C". From this you get the .root files. Use again "hadd" to combine all subruns. Then use retrieve_drift_fits.C. As input parameter use the comined .root file from the drift_analysis_tref.C output. The macro retrieve_drift_fits.C gives the parameters for the drift time calibration.

Design of the parameter file: section,anode,drift time, offset.

For the DeltaX vs Xal computation the macro "delta_x_final.C" is used (I don't remember why I have also a macro called "delta_x_drift.C", it does not consider the multiwires).

To get the final Tspline Fit and the arrangement I use the macro "retrieve_tspline_dt.C". As input parameter I use again the combined .root output from the macro "delta_x_final.C".

To get the position resolution of the TWIM Music anodes I use as input the output from the "retrieve_tspline_dt.C" macro. The resolution (sigma) is stored in the parameter file "xcal_std_dev_params.csv"