

# HPGe DETECTOR APPENDIX

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The following document is our appendix to our lab books that contain all of the figures that are being produced by our various codes, and the codes themselves, that have been produced over the four weeks of the experiment. Here, detailed plots are produced for all of the individual fits that have been done on all of the peaks, with all of the various fitting functions. Conclusions can be drawn from these figures, and so they are included here.

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## A1 FITTED PEAKS

For all of the known sources, we had to fit all of the peaks with various fitting functions to extract the central channel number for that peak. From this, it can then be fed into our tools that calibrates the detector properly, as we can produce a further fit that is of channel numbers as a function of peak energy.

To properly test the fits, we had to fit the data with three different functions to assess the quality of the fits, as well as how the mean and standard deviations of the fit varies with the fitting function. Then for each fit, we first ran a ‘rough’ fit, that is where we have manually gone through the plots, and isolated each peak extracting the approximate lower-bound, mean, and upper-bound of each peak. This then gave us a rough estimation of the location & mean of each peak, which was then passed into the fitting tool. From there, the new mean and standard deviation was extracted. From these new values, we then produced a ‘zoomed’ fit, by limiting the data to  $x = \mu \pm 2\sigma$  for each peak, which then constrained the peak correctly, producing a more accurate fit.

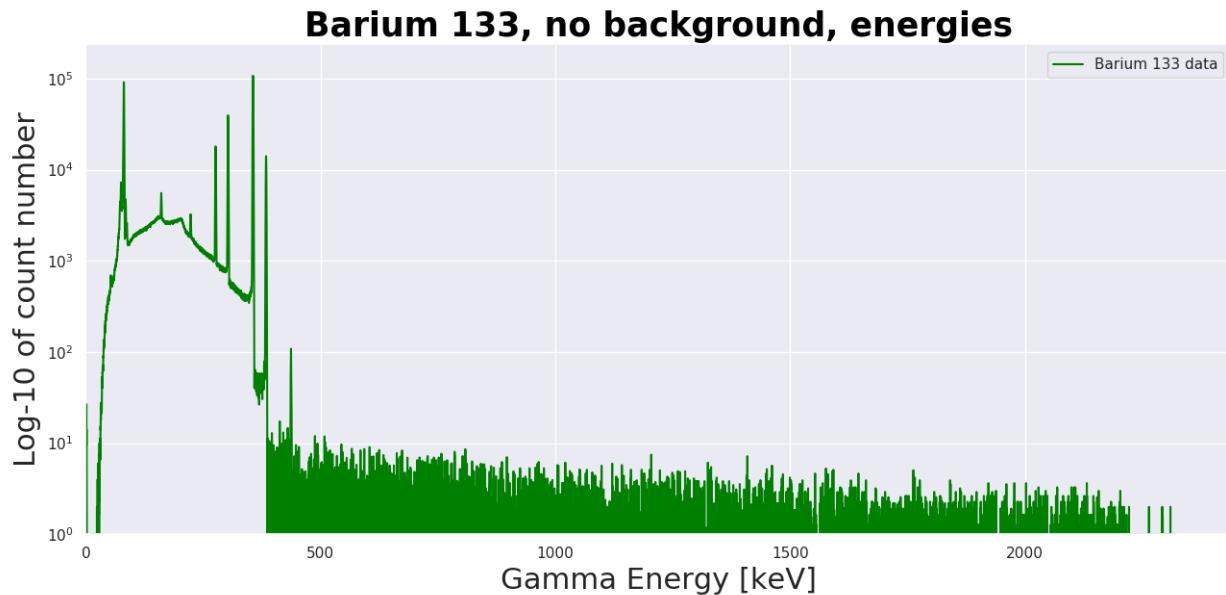
The reason behind this double-fitting process is that for some peaks, there are significant noise and background effects that causes the count numbers to behave significantly non-Gaussian, which can affect the location of the mean for the peak. By containing the fit to only the central maxima, we can reduce these effects significantly – and so the mean that we extract from the fit is much more likely to be that for the actual data, rather than any effects introducing any shifts.

All of the following plots are for the 24 hour data obtained, which is reflected by the high count numbers. By running each source for longer, we should be able to reduce the effects of any statistical noise fluctuations that may induce a slight shift in the data.

The following plots are broken down first by source, in decreasing atomic number, and then further broken down into subsections by the fit type used to produce each plot. For each fit, there are the the fit statistics and the mean & standard deviation for the peak.

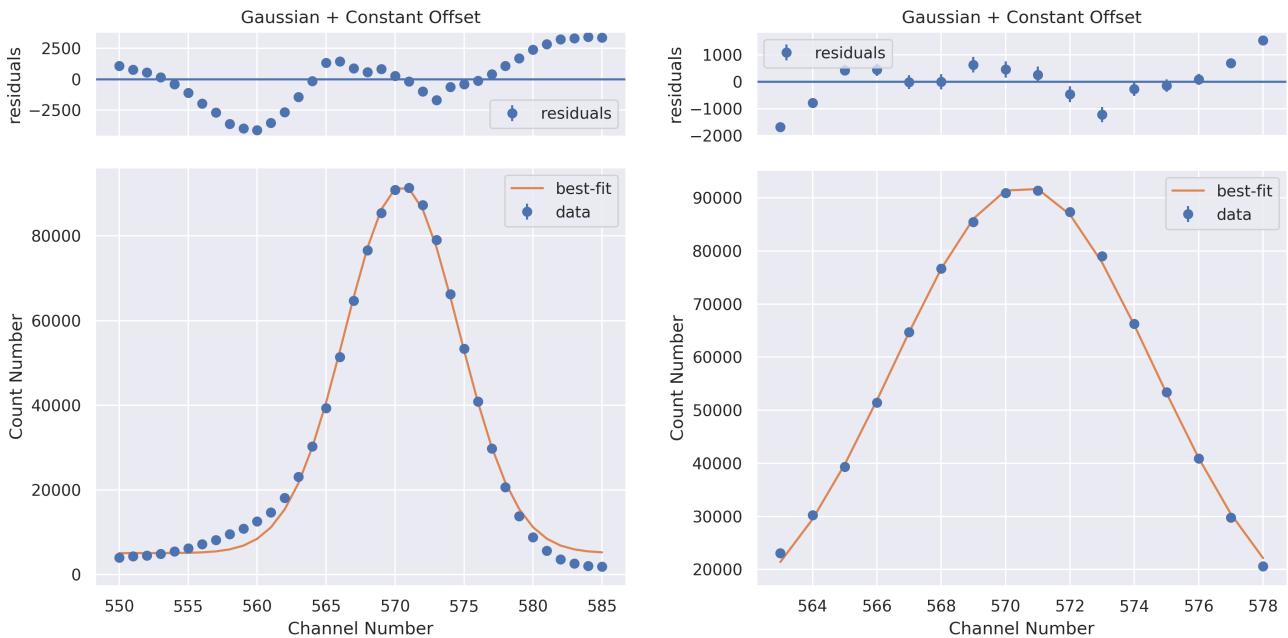
### A1.1 BARIUM-133

The  $^{133}\text{Ba}$  source had the most amount of identifiable peaks, totalling seven. This gave us a large number of peaks to fit, and so improved the accuracy of our energy calibration. All of these peaks were considered low energy, as they all occur below 400 keV. This is shown in the figure below



**Figure A.1:** The final spectrum obtained for the  $^{133}\text{Ba}$  source run for 24 hours. This has been converted to energies by using the quadratic energy calibration.

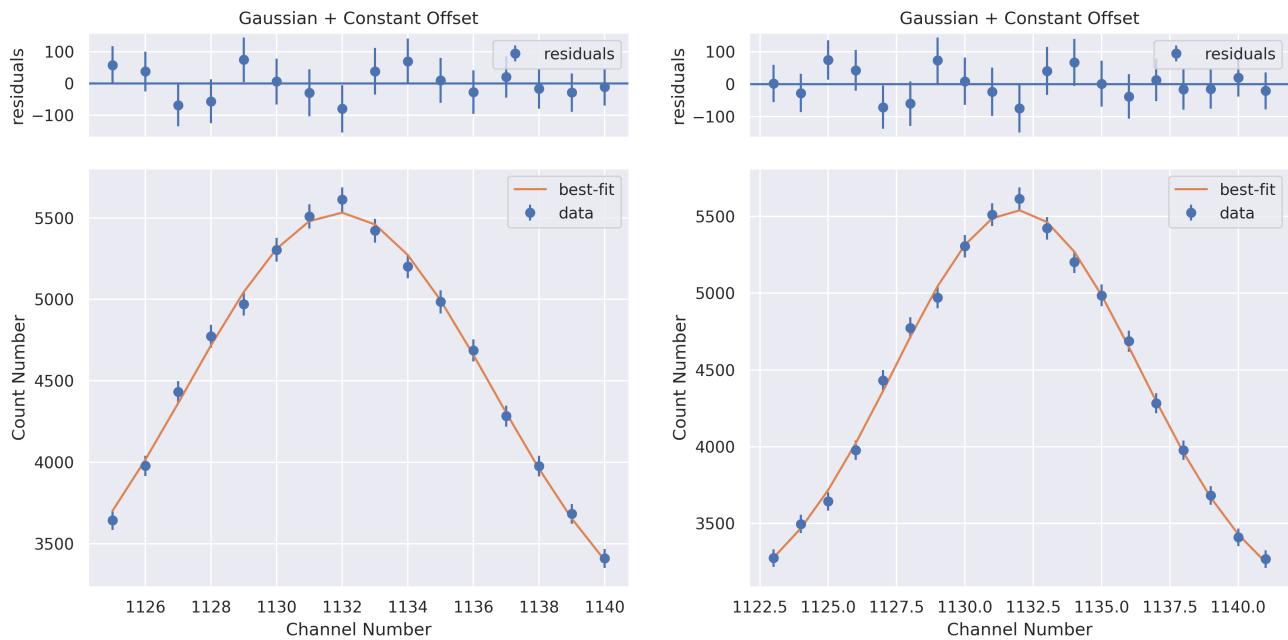
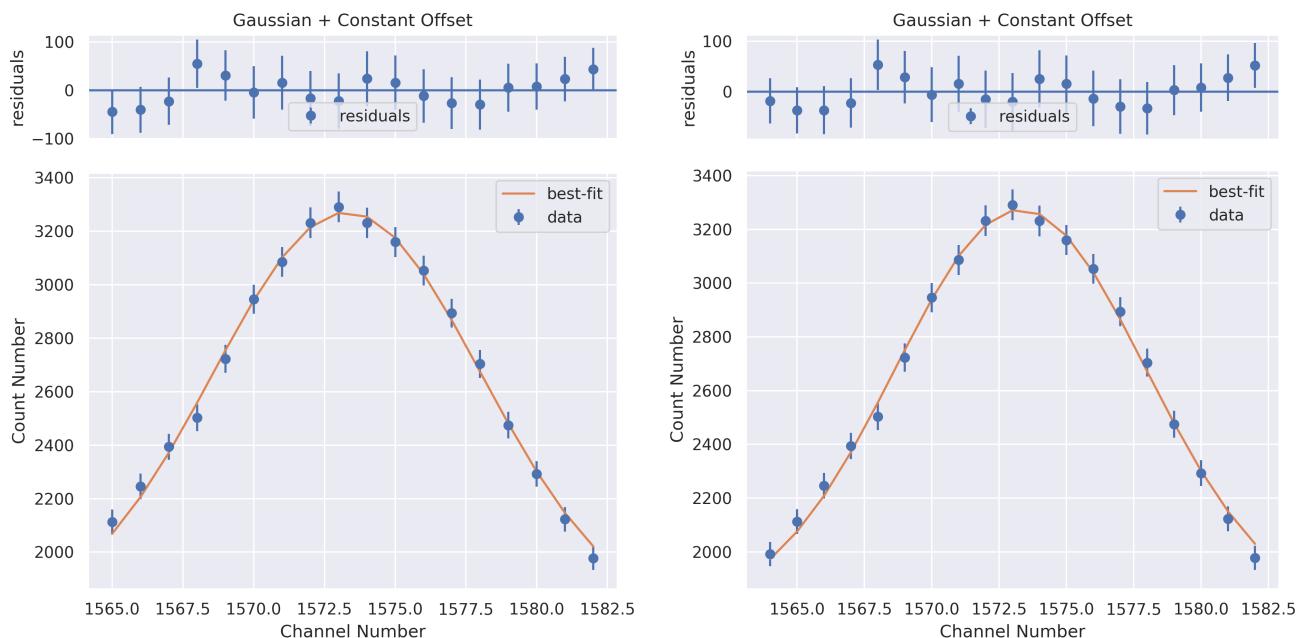
### A1.1.1 GAUSSIAN + OFFSET FIT

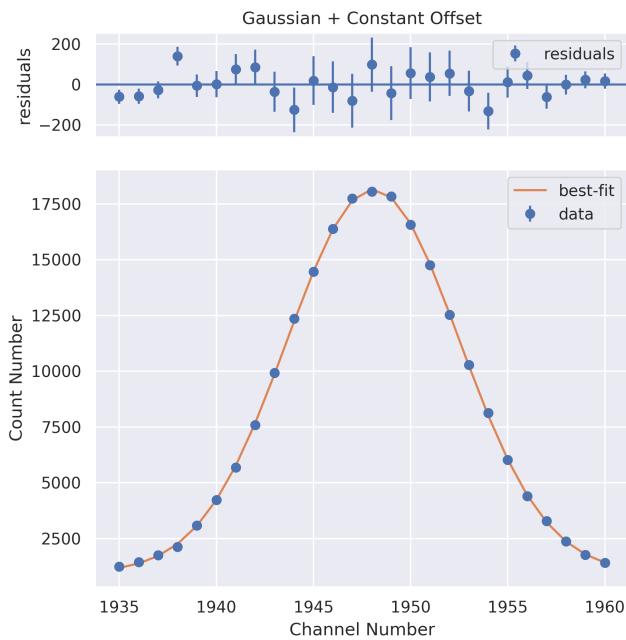


(a) Full peak with fit.  $\chi^2 = 29044$ ,  $\chi^2_\nu = 854$ , Prob = 0%,  $\mu = 570.5$ ,  $\sigma = 4.12$

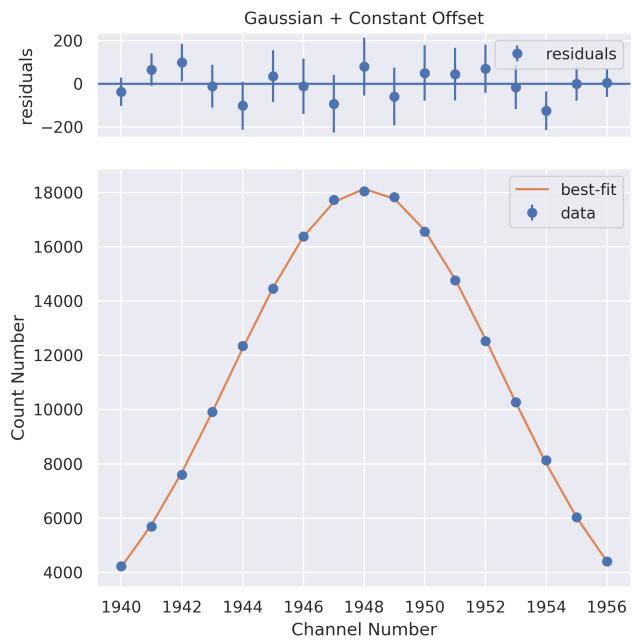
(b) Zoomed in peak with fit.  $\chi^2 = 310.6$ ,  $\chi^2_\nu = 22.2$ , Prob = 0%,  $\mu = 570.6$ ,  $\sigma = 4.00$

**Figure A.2:** Fit of full & zoomed in peak of  $^{133}\text{Ba}$  81 keV peak

**Figure A.3:** Fit of full & zoomed in peak of  $^{133}\text{Ba}$  161 keV peak**Figure A.4:** Fit of full & zoomed in peak of  $^{133}\text{Ba}$  223 keV peak

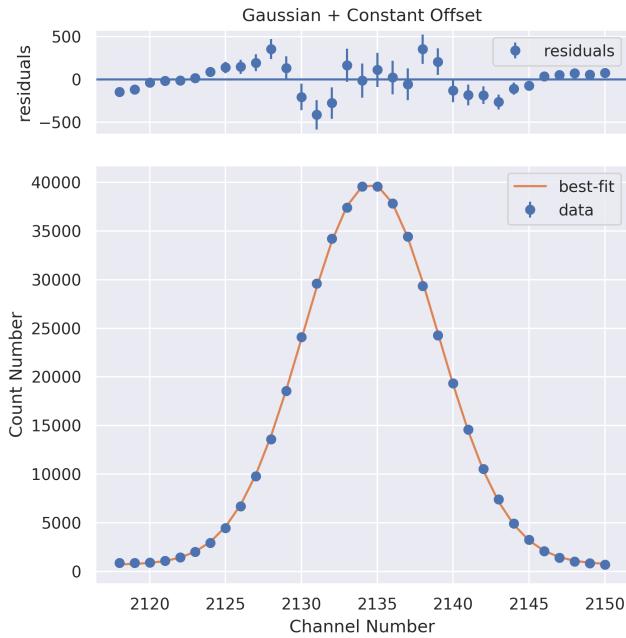


(a) Full peak with fit.  $\chi^2 = 24.2$ ,  $\chi^2_\nu = 1.01$ ,  
Prob = 44.9%,  $\mu = 1948.1$ ,  $\sigma = 4.43$

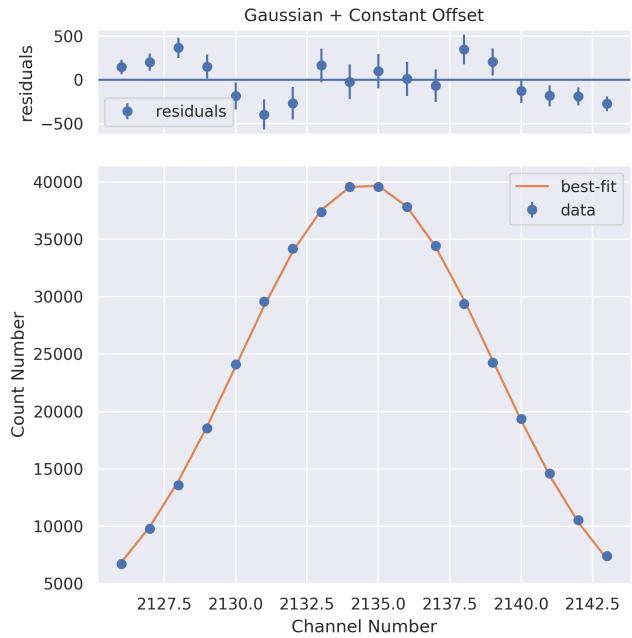


(b) Zoomed in peak with fit.  $\chi^2 = 6.92$ ,  $\chi^2_\nu = 0.46$ ,  
Prob = 96.0%,  $\mu = 1948.1$ ,  $\sigma = 4.48$

**Figure A.5:** Fit of full & zoomed in peak of  $^{133}\text{Ba}$  276 keV peak

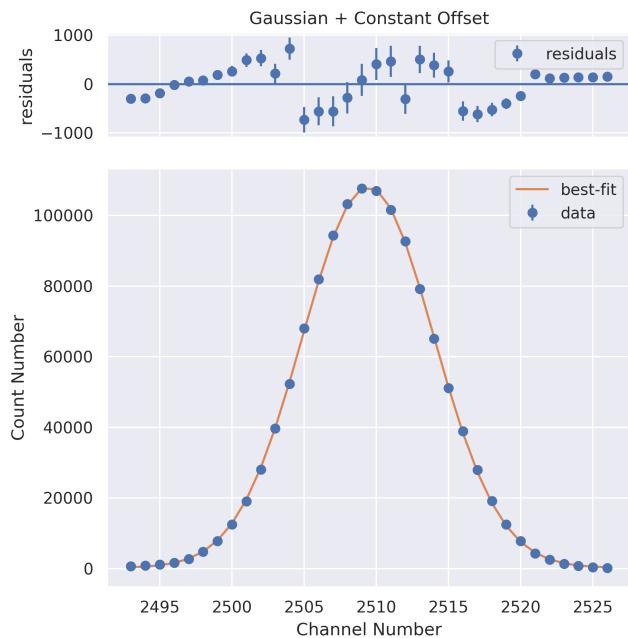


(a) Full peak with fit.  $\chi^2 = 121$ ,  $\chi^2_\nu = 4$ ,  
Prob = 0%,  $\mu = 2134.55$ ,  $\sigma = 4.45$

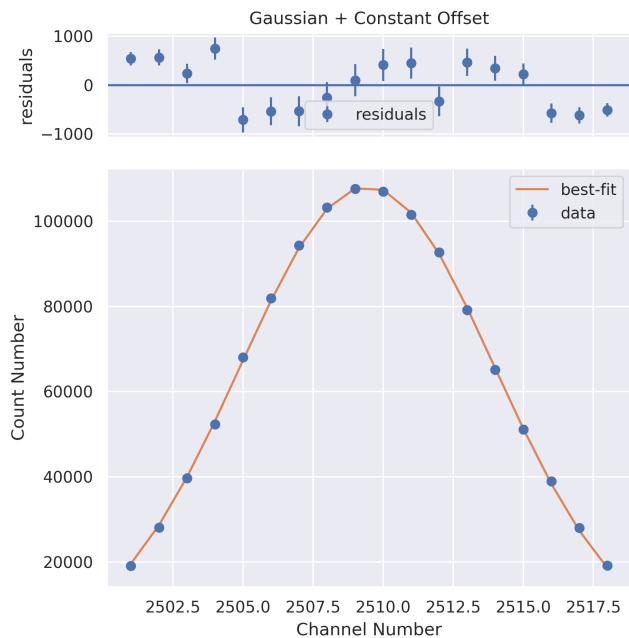


(b) Zoomed in peak with fit.  $\chi^2 = 51.0$ ,  $\chi^2_\nu = 3.2$ ,  
Prob = 0.02%,  $\mu = 2134.55$ ,  $\sigma = 4.56$

**Figure A.6:** Fit of full & zoomed in peak of  $^{133}\text{Ba}$  303 keV peak

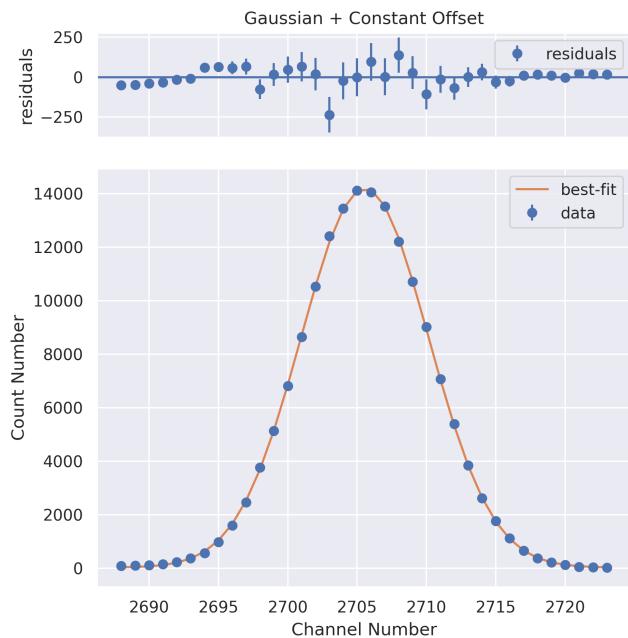


(a) Full peak with fit.  $\chi^2 = 608$ ,  $\chi^2_\nu = 19$ ,  
Prob = 0%,  $\mu = 2509.4$ ,  $\sigma = 4.5$

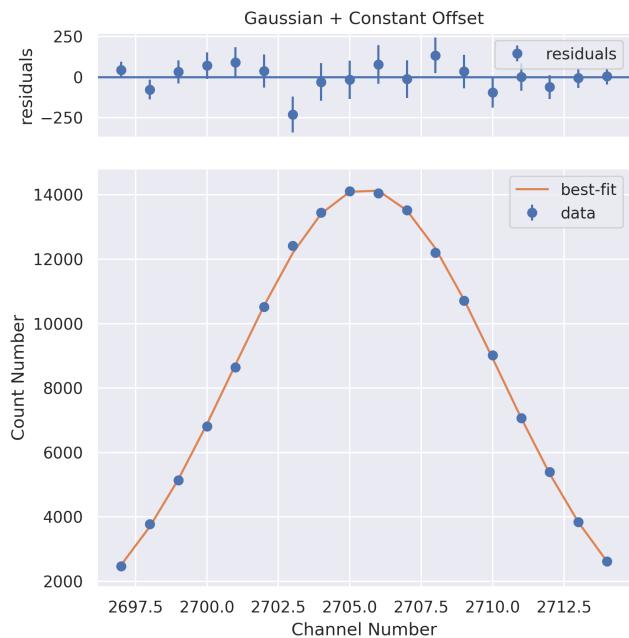


(b) Zoomed in peak with fit.  $\chi^2 = 100.1$ ,  $\chi^2_\nu = 6.3$ ,  
Prob = 0%,  $\mu = 2509.4$ ,  $\sigma = 4.5$

**Figure A.7:** Fit of full & zoomed in peak of  $^{133}\text{Ba}$  356 keV peak



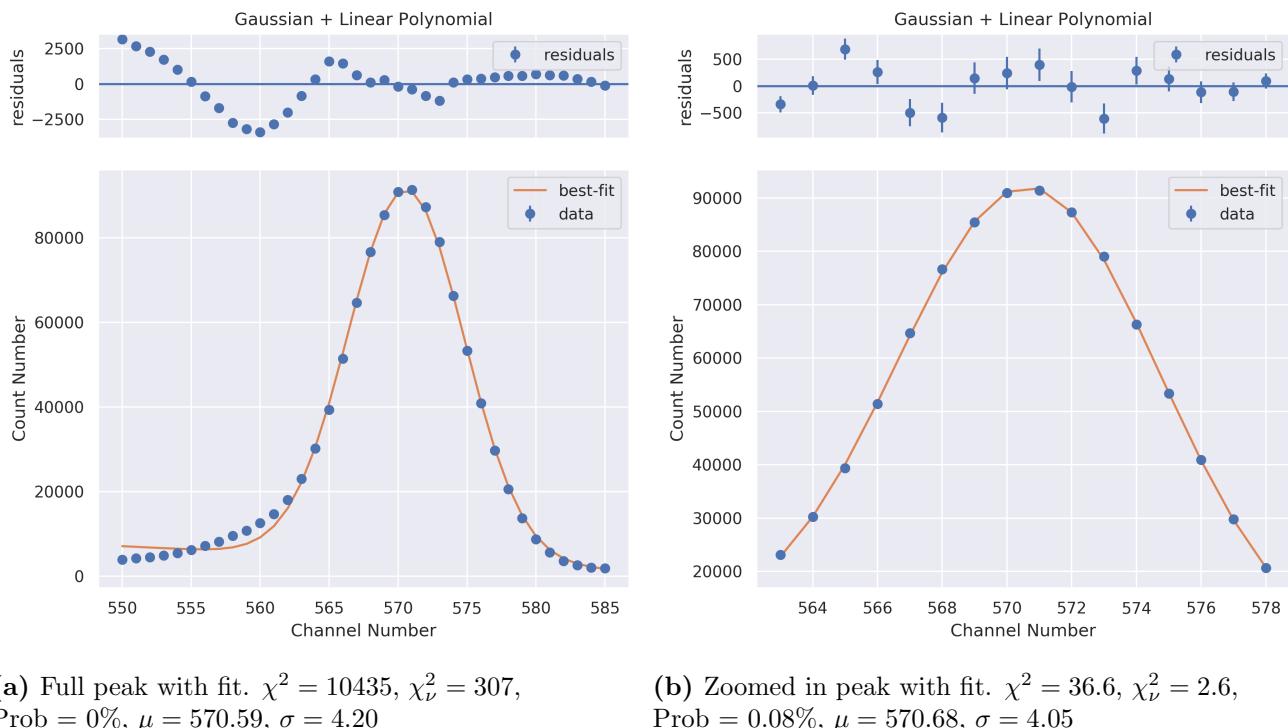
(a) Full peak with fit.  $\chi^2 = 127$ ,  $\chi^2_\nu = 4$ ,  
Prob = 0%,  $\mu = 2705.56$ ,  $\sigma = 4.59$



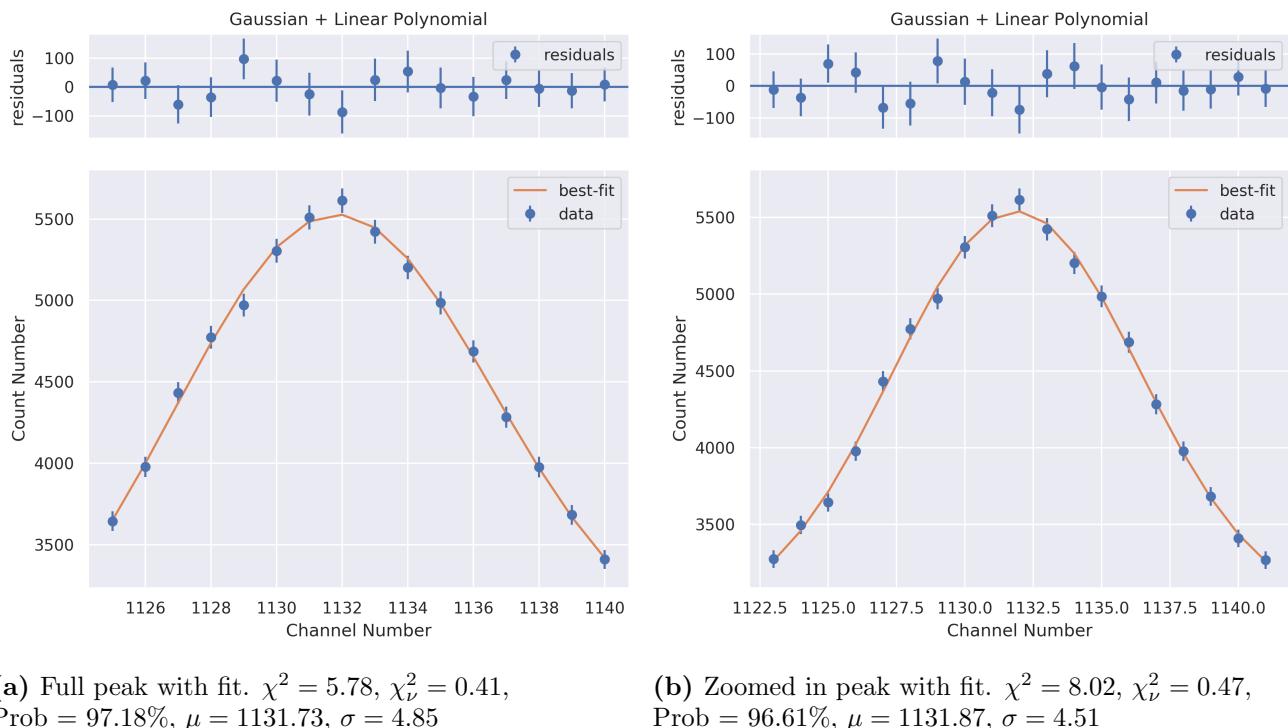
(b) Zoomed in peak with fit.  $\chi^2 = 12.5$ ,  $\chi^2_\nu = 0.8$ ,  
Prob = 70.9%,  $\mu = 2705.55$ ,  $\sigma = 4.64$

**Figure A.8:** Fit of full & zoomed in peak of  $^{133}\text{Ba}$  384 keV peak

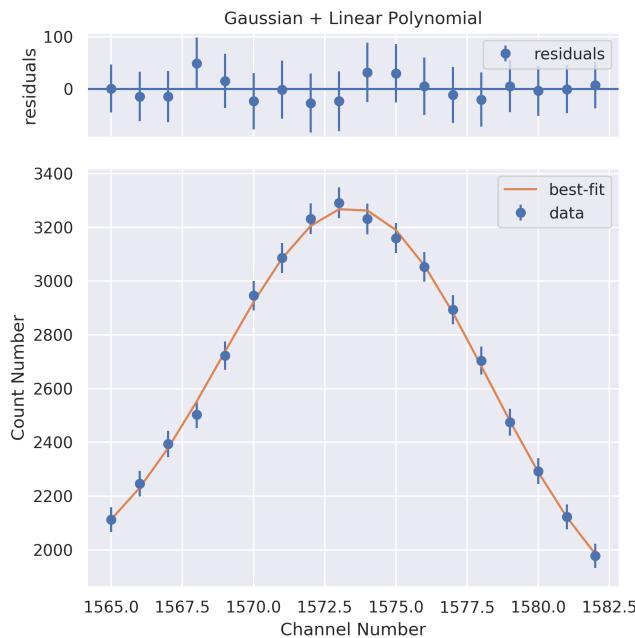
### A1.1.2 GAUSSIAN + LINEAR FIT



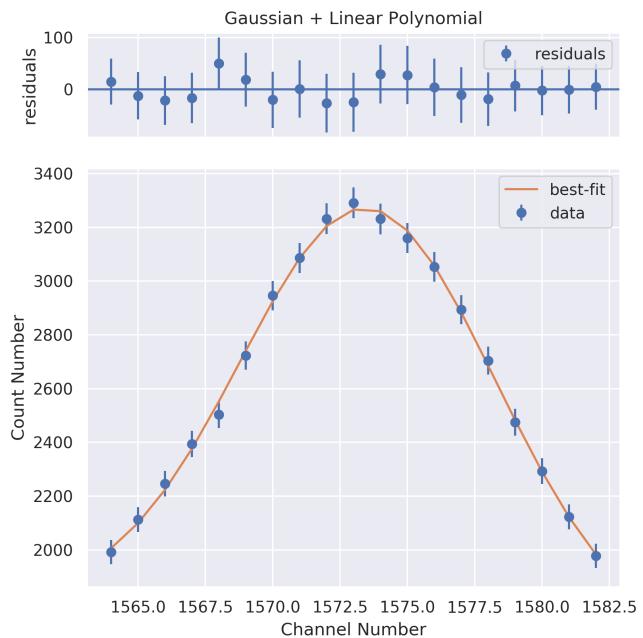
**Figure A.9:** Fit of full & zoomed in peak of  $^{133}\text{Ba}$  81 keV peak



**Figure A.10:** Fit of full & zoomed in peak of  $^{133}\text{Ba}$  161 keV peak

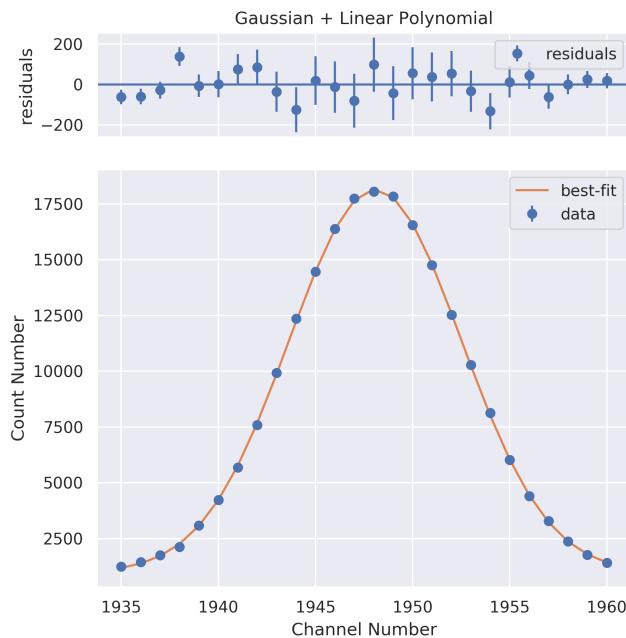


(a) Full peak with fit.  $\chi^2 = 2.65$ ,  $\chi_\nu^2 = 0.17$ ,  
Prob = 99.99%,  $\mu = 1573.55$ ,  $\sigma = 4.66$

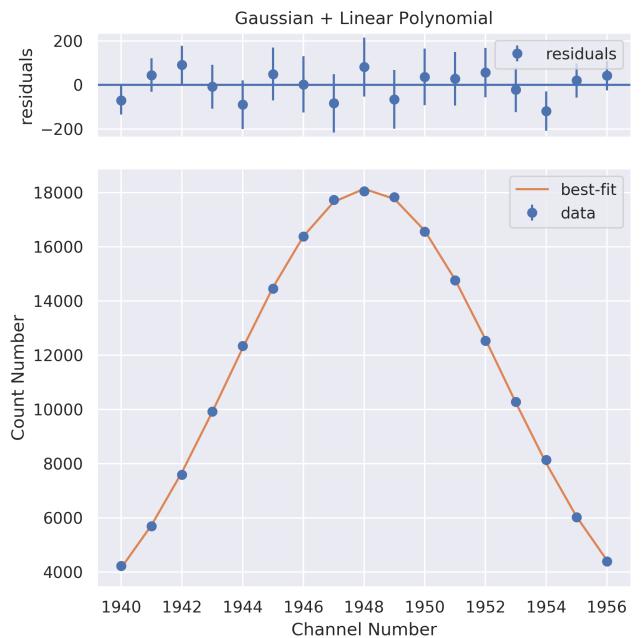


(b) Zoomed in peak with fit.  $\chi^2 = 2.93$ ,  $\chi_\nu^2 = 0.17$ ,  
Prob = 99.99%,  $\mu = 1573.52$ ,  $\sigma = 4.74$

**Figure A.11:** Fit of full & zoomed in peak of  $^{133}\text{Ba}$  223 keV peak

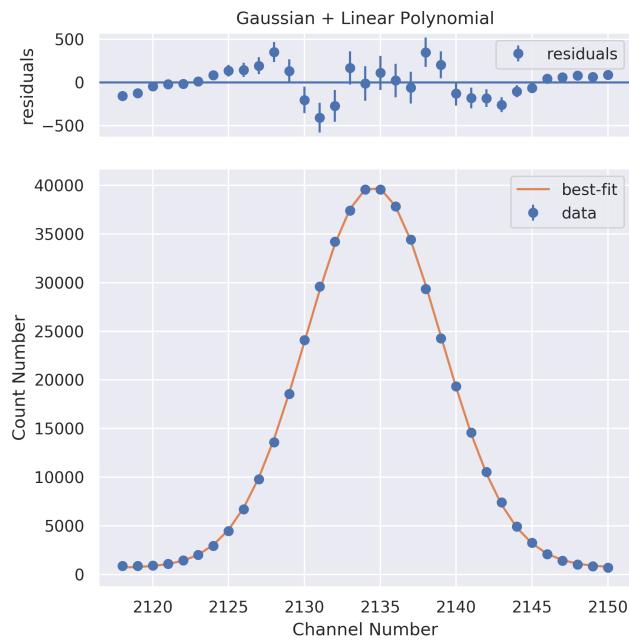


(a) Full peak with fit.  $\chi^2 = 24.39$ ,  $\chi_\nu^2 = 1.02$ ,  
Prob = 43.93%,  $\mu = 1948.08$ ,  $\sigma = 4.43$

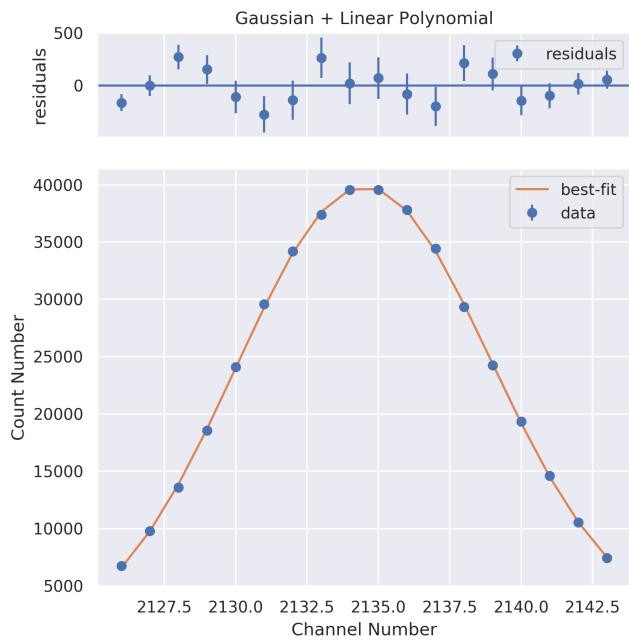


(b) Zoomed in peak with fit.  $\chi^2 = 7.08$ ,  $\chi_\nu^2 = 0.47$ ,  
Prob = 95.54%,  $\mu = 1948.06$ ,  $\sigma = 4.47$

**Figure A.12:** Fit of full & zoomed in peak of  $^{133}\text{Ba}$  276 keV peak

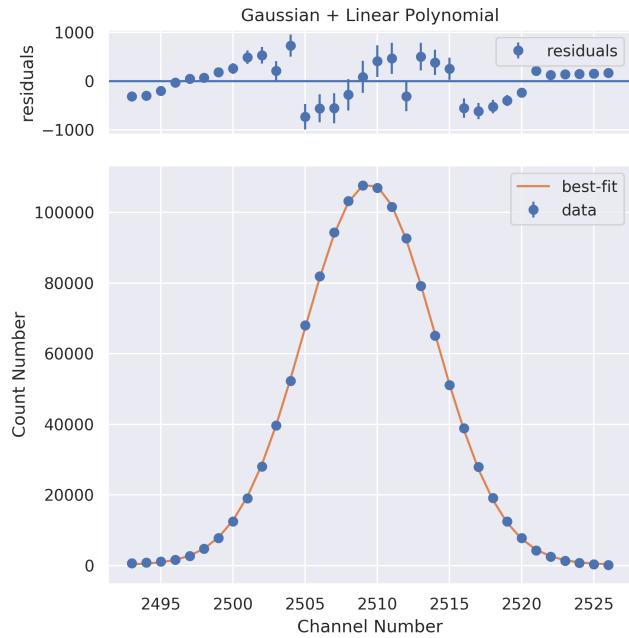


(a) Full peak with fit.  $\chi^2 = 131.45$ ,  $\chi^2_\nu = 4.24$ ,  
Prob = 0%,  $\mu = 2134.55$ ,  $\sigma = 4.45$

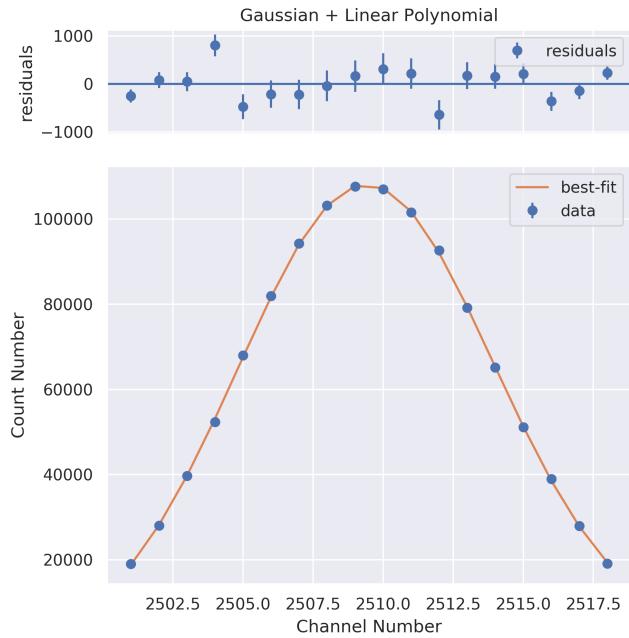


(b) Zoomed in peak with fit.  $\chi^2 = 21.84$ ,  $\chi^2_\nu = 1.36$ ,  
Prob = 14.85%,  $\mu = 2134.49$ ,  $\sigma = 4.45$

**Figure A.13:** Fit of full & zoomed in peak of  $^{133}\text{Ba}$  303 keV peak

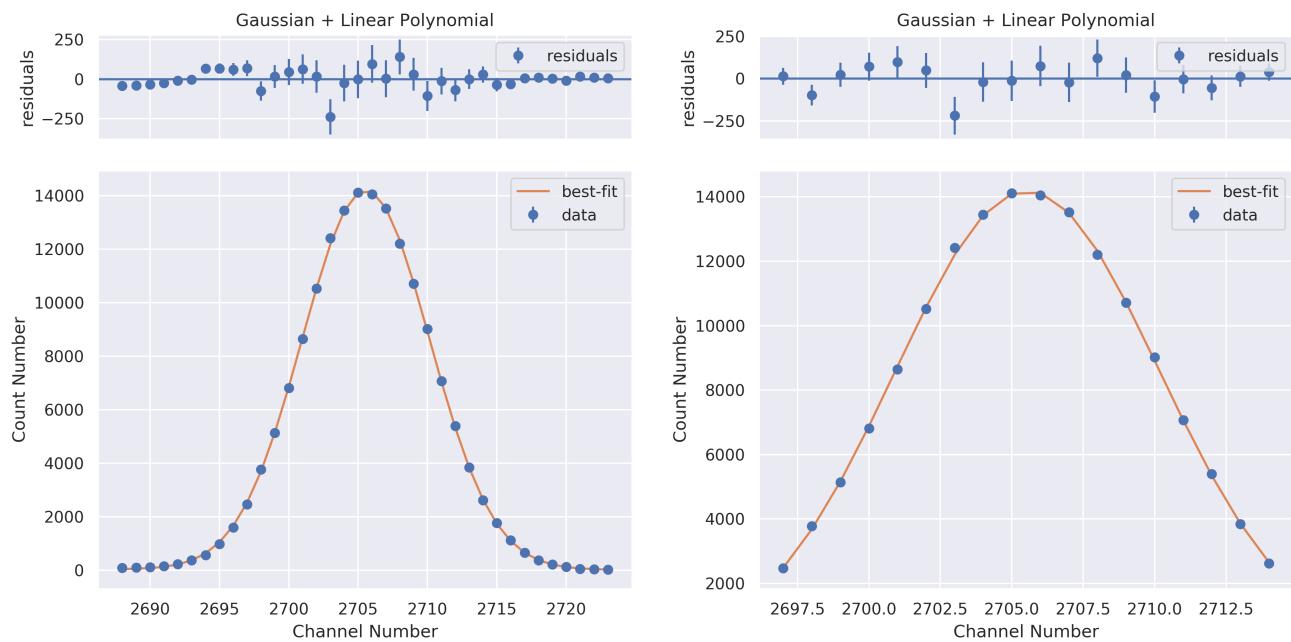


(a) Full peak with fit.  $\chi^2 = 658.02$ ,  $\chi^2_\nu = 20.56$ ,  
Prob = 0%,  $\mu = 2509.44$ ,  $\sigma = 4.55$



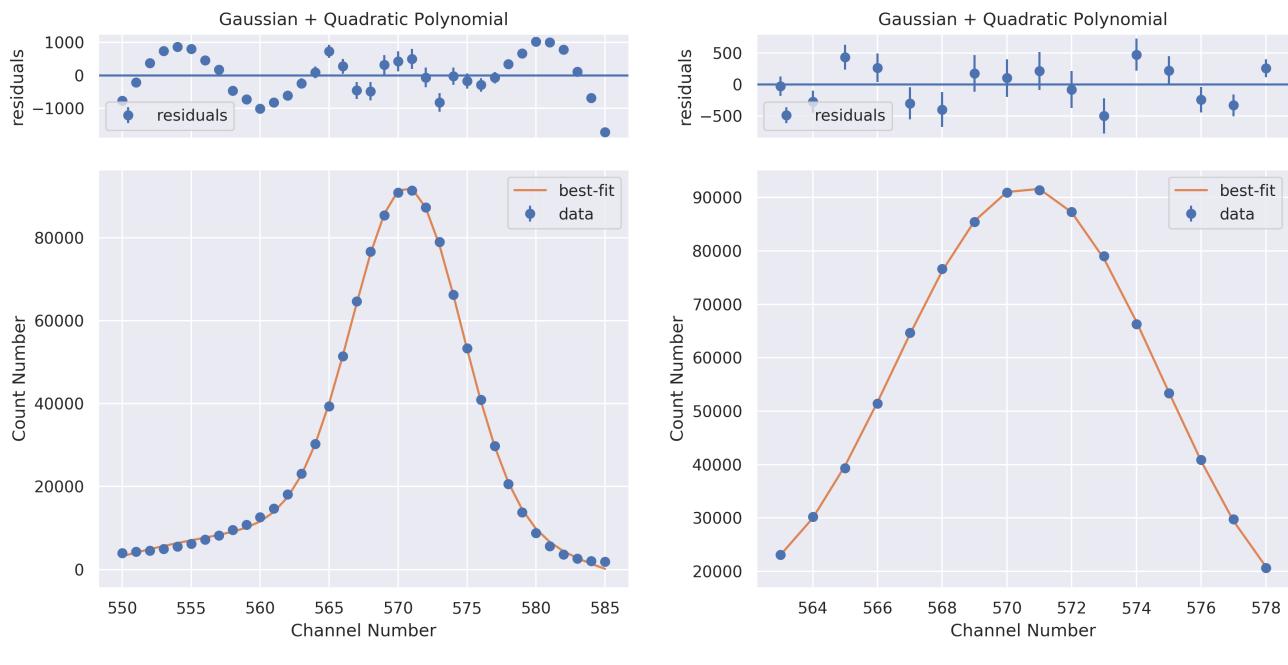
(b) Zoomed in peak with fit.  $\chi^2 = 34.92$ ,  $\chi^2_\nu = 2.18$ ,  
Prob = 0.41%,  $\mu = 2509.38$ ,  $\sigma = 4.55$

**Figure A.14:** Fit of full & zoomed in peak of  $^{133}\text{Ba}$  356 keV peak

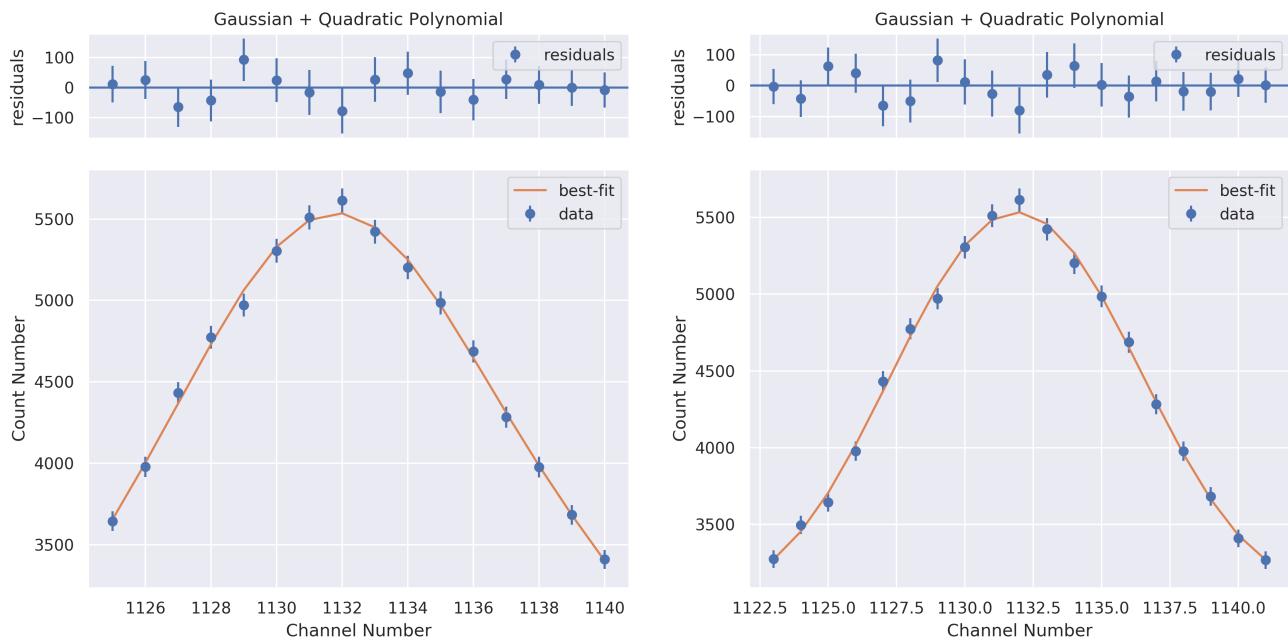


**Figure A.15:** Fit of full & zoomed in peak of  $^{133}\text{Ba}$  384 keV peak

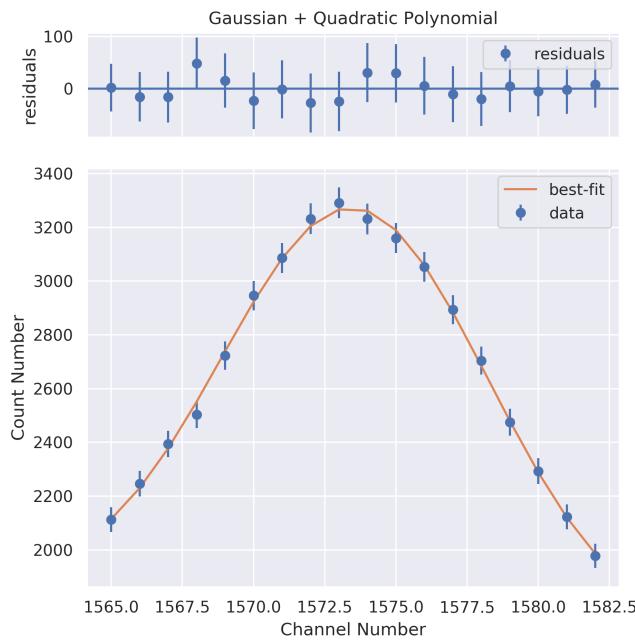
### A1.1.3 GAUSSIAN + QUADRATIC FIT



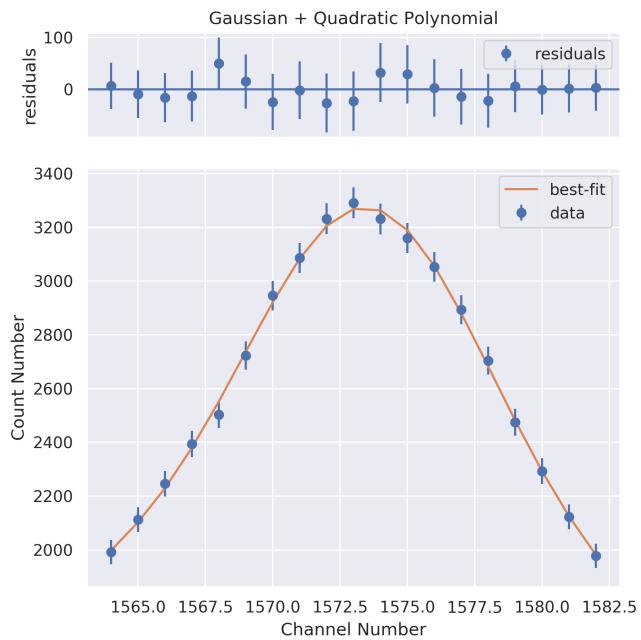
**Figure A.16:** Fit of full & zoomed in peak of  $^{133}\text{Ba}$  81 keV peak



**Figure A.17:** Fit of full & zoomed in peak of  $^{133}\text{Ba}$  161 keV peak

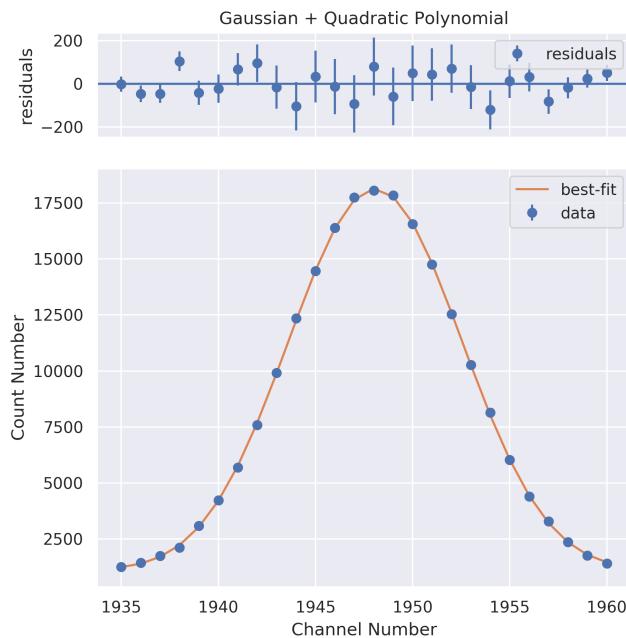


(a) Full peak with fit.  $\chi^2 = 2.66$ ,  $\chi^2_\nu = 0.17$ ,  
Prob = 99.99%,  $\mu = 1573.55$ ,  $\sigma = 4.82$

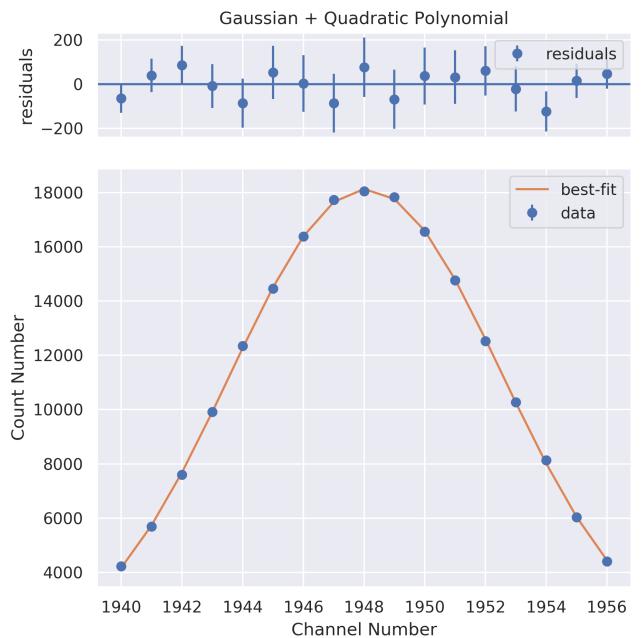


(b) Zoomed in peak with fit.  $\chi^2 = 2.77$ ,  $\chi^2_\nu = 0.16$ ,  
Prob = 100%,  $\mu = 1573.53$ ,  $\sigma = 4.34$

**Figure A.18:** Fit of full & zoomed in peak of  $^{133}\text{Ba}$  223 keV peak

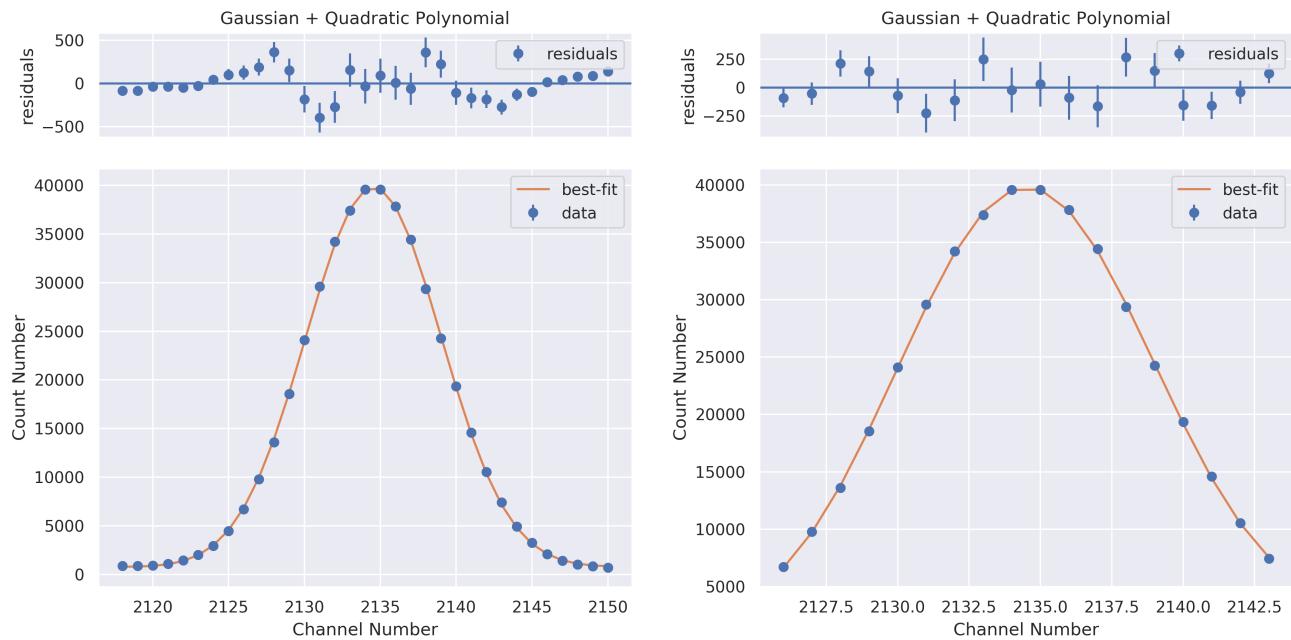


(a) Full peak with fit.  $\chi^2 = 19.63$ ,  $\chi^2_\nu = 0.82$ ,  
Prob = 71.76%,  $\mu = 1948.07$ ,  $\sigma = 4.49$



(b) Zoomed in peak with fit.  $\chi^2 = 6.91$ ,  $\chi^2_\nu = 0.46$ ,  
Prob = 96.00%,  $\mu = 1948.06$ ,  $\sigma = 4.54$

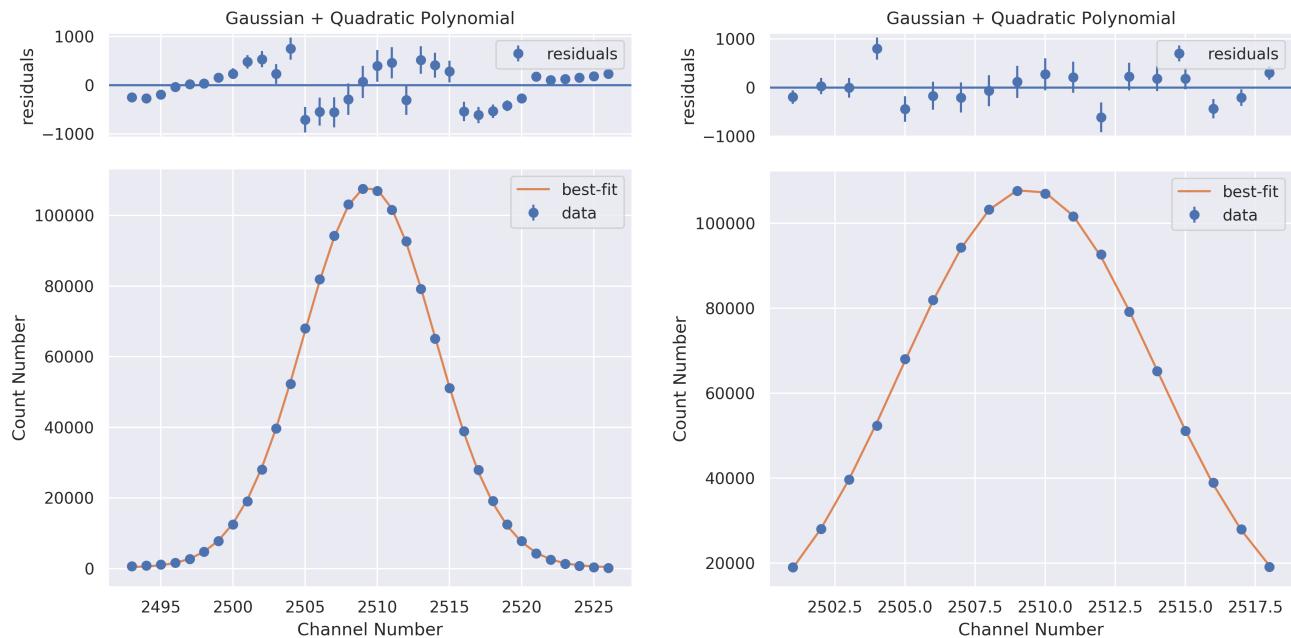
**Figure A.19:** Fit of full & zoomed in peak of  $^{133}\text{Ba}$  276 keV peak



(a) Full peak with fit  $\chi^2 = 123.06$ ,  $\chi_\nu^2 = 3.97$ , Prob = 0.00%,  $\mu = 2134.55$ ,  $\sigma = 4.47$

(b) Zoomed in peak with fit.  $\chi^2 = 19.63$ ,  $\chi_\nu^2 = 1.23$ , Prob = 23.74%,  $\mu = 2134.49$ ,  $\sigma = 4.73$

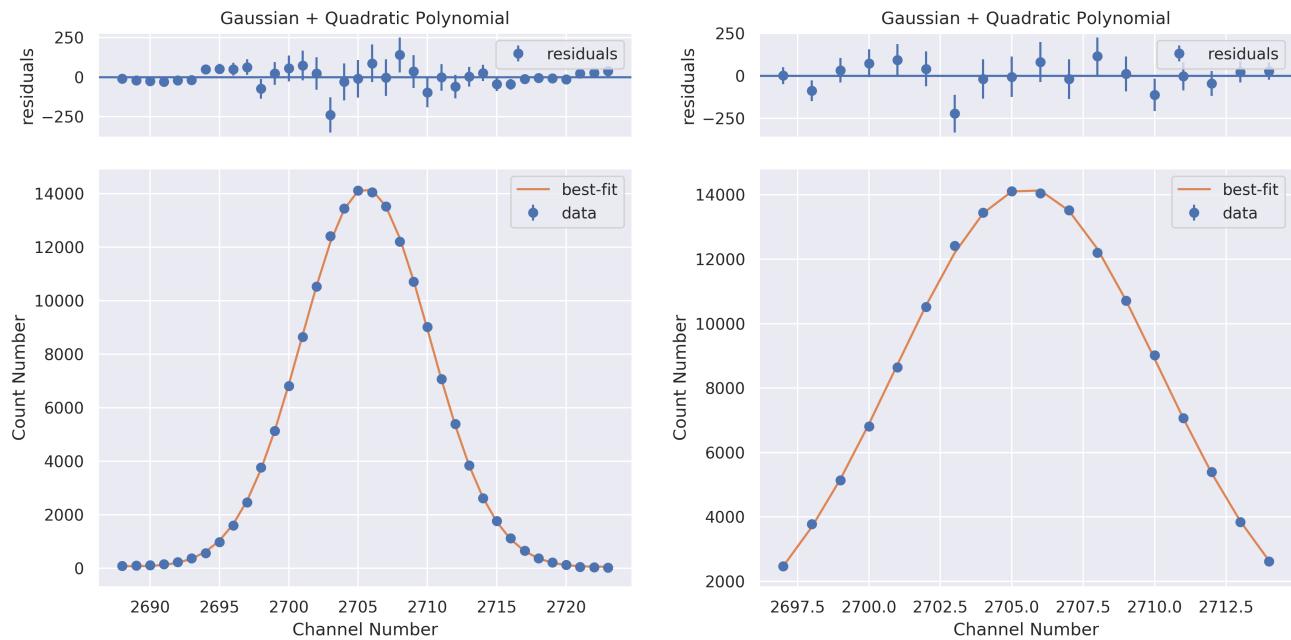
**Figure A.20:** Fit of full & zoomed in peak of  $^{133}\text{Ba}$  303 keV peak



(a) Full peak with fit.  $\chi^2 = 722.64$ ,  $\chi_\nu^2 = 22.58$ , Prob = 0.00%,  $\mu = 2509.44$ ,  $\sigma = 4.56$

(b) Zoomed in peak with fit.  $\chi^2 = 3.14$ ,  $\chi_\nu^2 = 2.26$ , Prob = 0.28%,  $\mu = 2509.38$ ,  $\sigma = 4.65$

**Figure A.21:** Fit of full & zoomed in peak of  $^{133}\text{Ba}$  356 keV peak



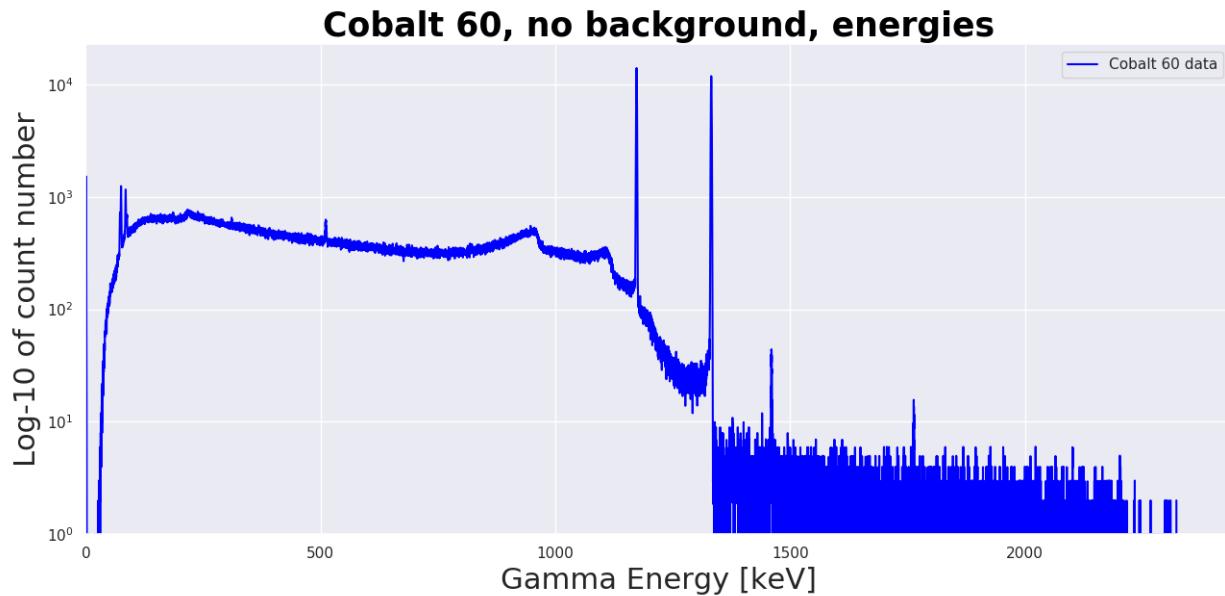
(a) Full peak with fit.  $\chi^2 = 119.72$ ,  $\chi^2_\nu = 3.52$ ,  
Prob = 0.00%,  $\mu = 2705.56$ ,  $\sigma = 4.62$

(b) Zoomed in peak with fit.  $\chi^2 = 12.07$ ,  $\chi^2_\nu = 0.75$ ,  
Prob = 73.93%,  $\mu = 2705.53$ ,  $\sigma = 4.50$

**Figure A.22:** Fit of full & zoomed in peak of  $^{133}\text{Ba}$  384 keV peak

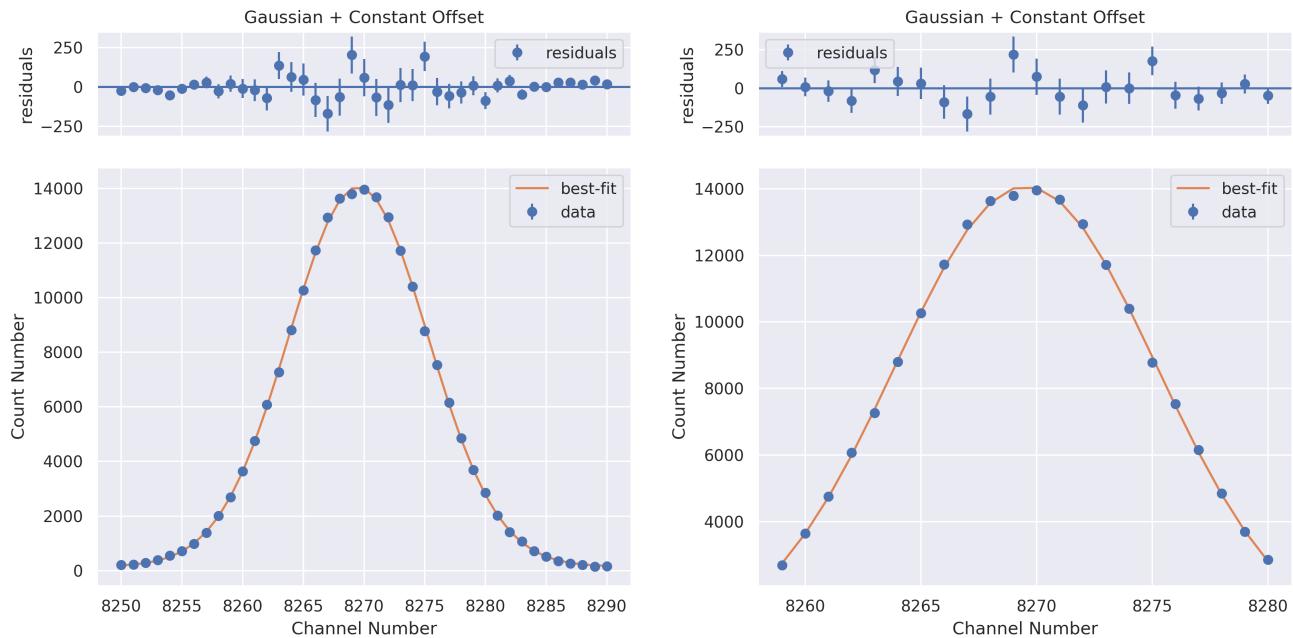
## A1.2 COBALT-60

The  $^{60}\text{Co}$  source had two high-energy peaks which were very distinctive, and so quite easy to fit for all functions.



**Figure A.23:** The final spectrum obtained for the  $^{60}\text{Co}$  source run for 24 hours. This has been converted to energies by using the quadratic energy calibration.

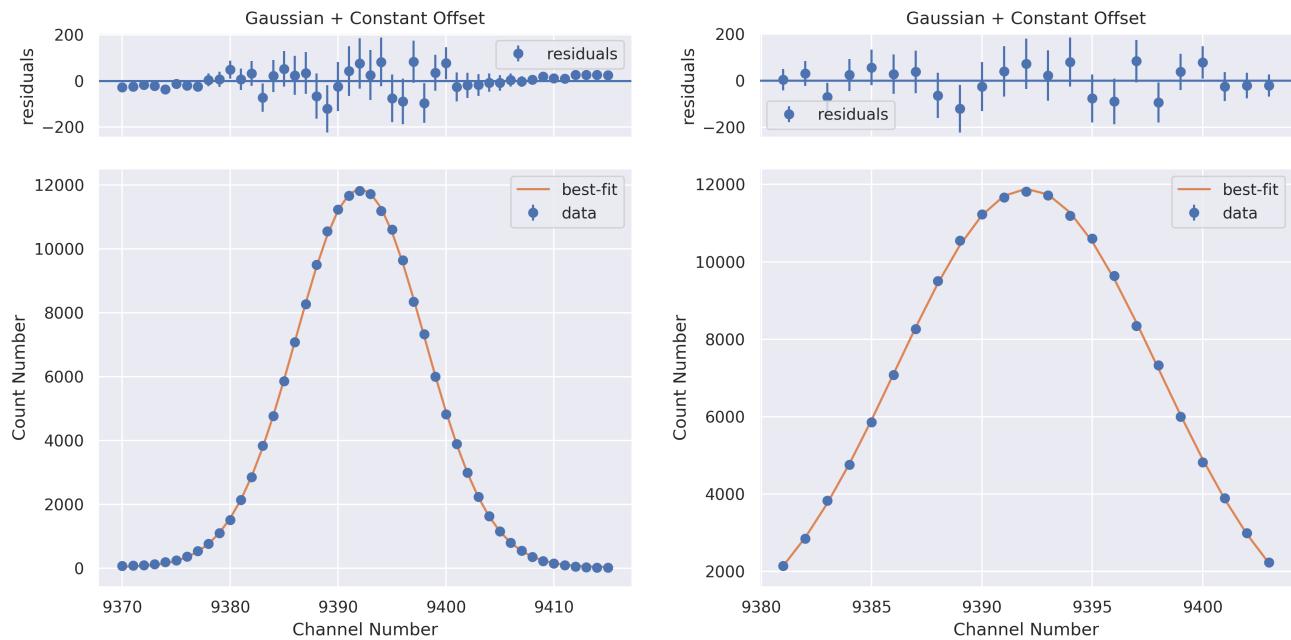
### A1.2.1 GAUSSIAN + OFFSET FIT



(a) Full peak with fit.  $\chi^2 = 54.04$ ,  $\chi^2_\nu = 1.39$ ,  
Prob = 5.52%,  $\mu = 8269.53$ ,  $\sigma = 5.73$

(b) Zoomed in peak with fit.  $\chi^2 = 18.66$ ,  $\chi^2_\nu = 0.93$ ,  
Prob = 54.42%,  $\mu = 8269.54$ ,  $\sigma = 5.67$

**Figure A.24:** Fit of full & zoomed in peak of  $^{60}\text{Co}$  1173 keV peak

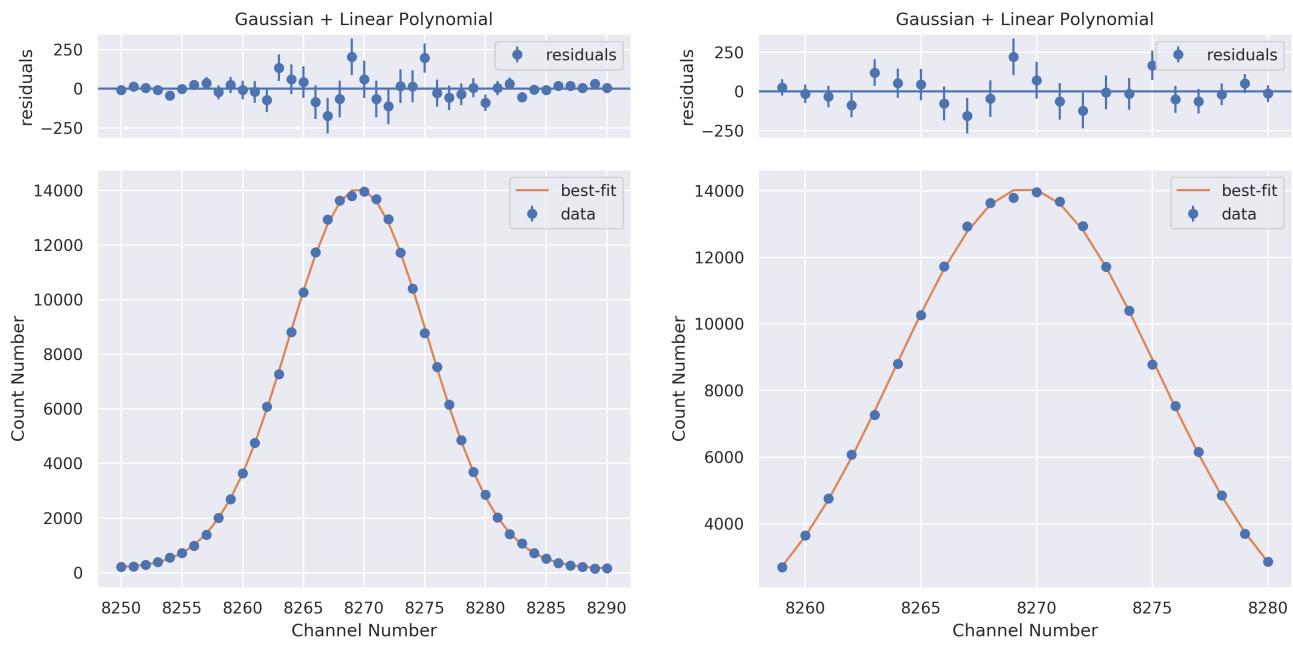


(a) Full peak with fit.  $\chi^2 = 173.52$ ,  $\chi^2_\nu = 3.94$ ,  
Prob = 0.00%,  $\mu = 9392.05$ ,  $\sigma = 5.95$

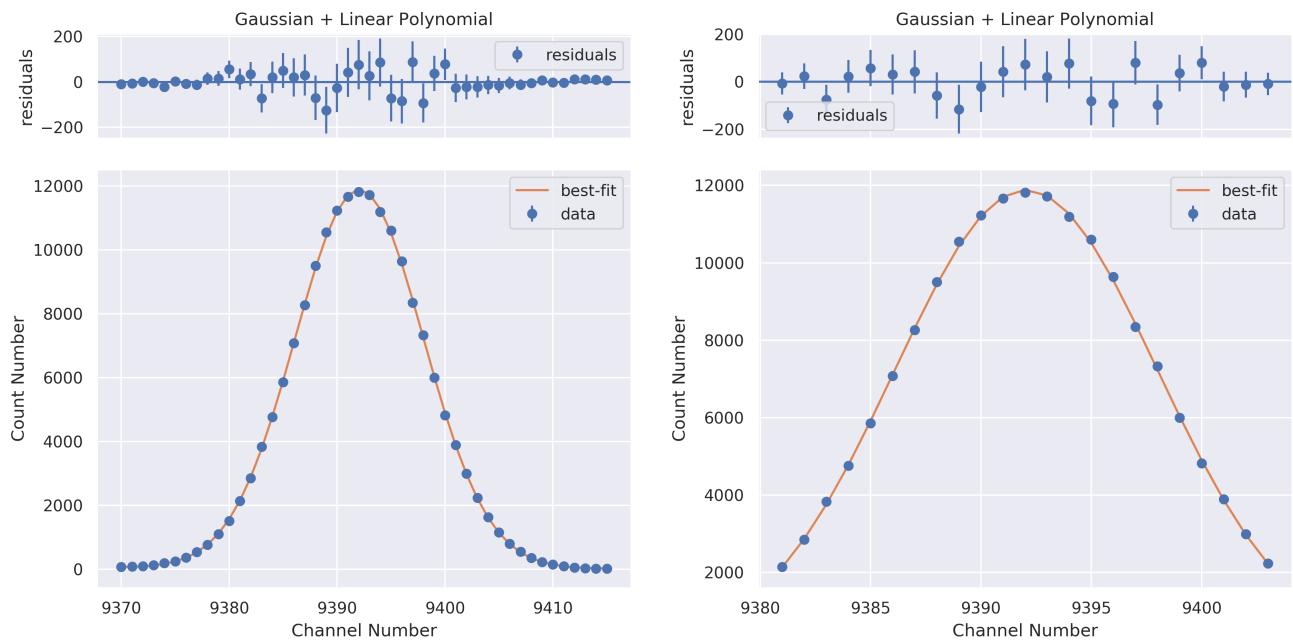
(b) Zoomed in peak with fit.  $\chi^2 = 11.12$ ,  $\chi^2_\nu = 0.53$ ,  
Prob = 96.04%,  $\mu = 9392.05$ ,  $\sigma = 5.96$

**Figure A.25:** Fit of full & zoomed in peak of  $^{60}\text{Co}$  1332 keV peak

### A1.2.2 GAUSSIAN + LINEAR FIT

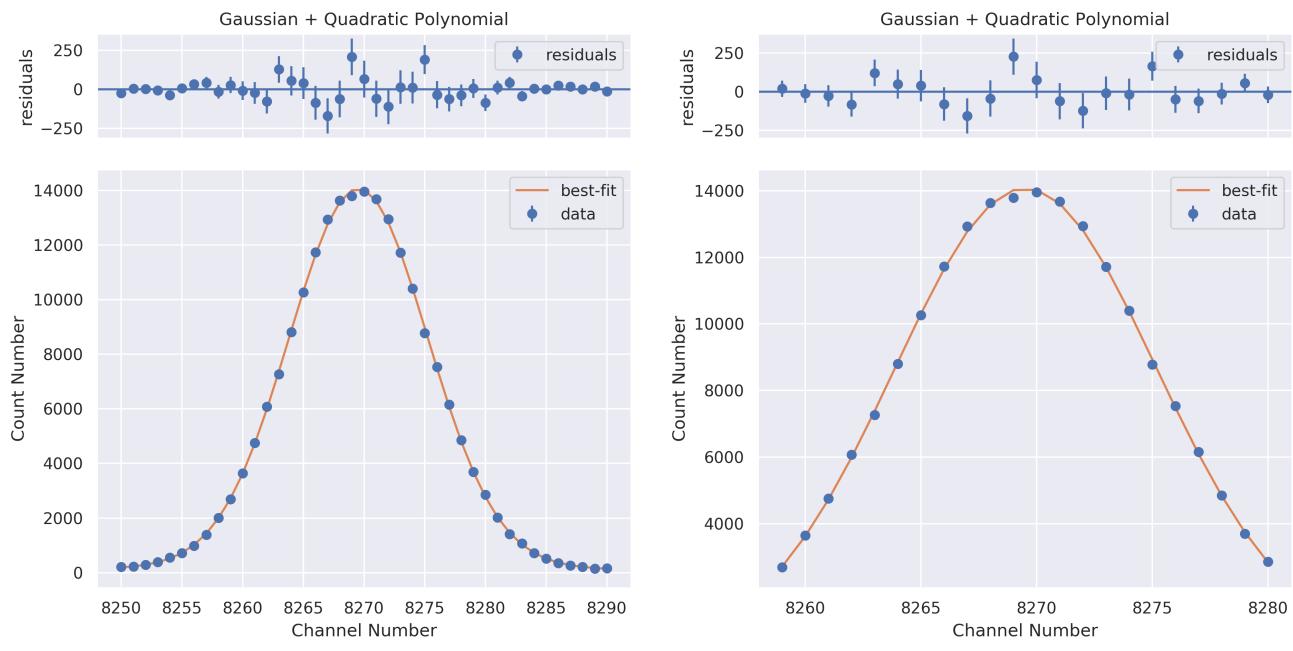


**Figure A.26:** Fit of full & zoomed in peak of  $^{60}\text{Co}$  1173 keV peak

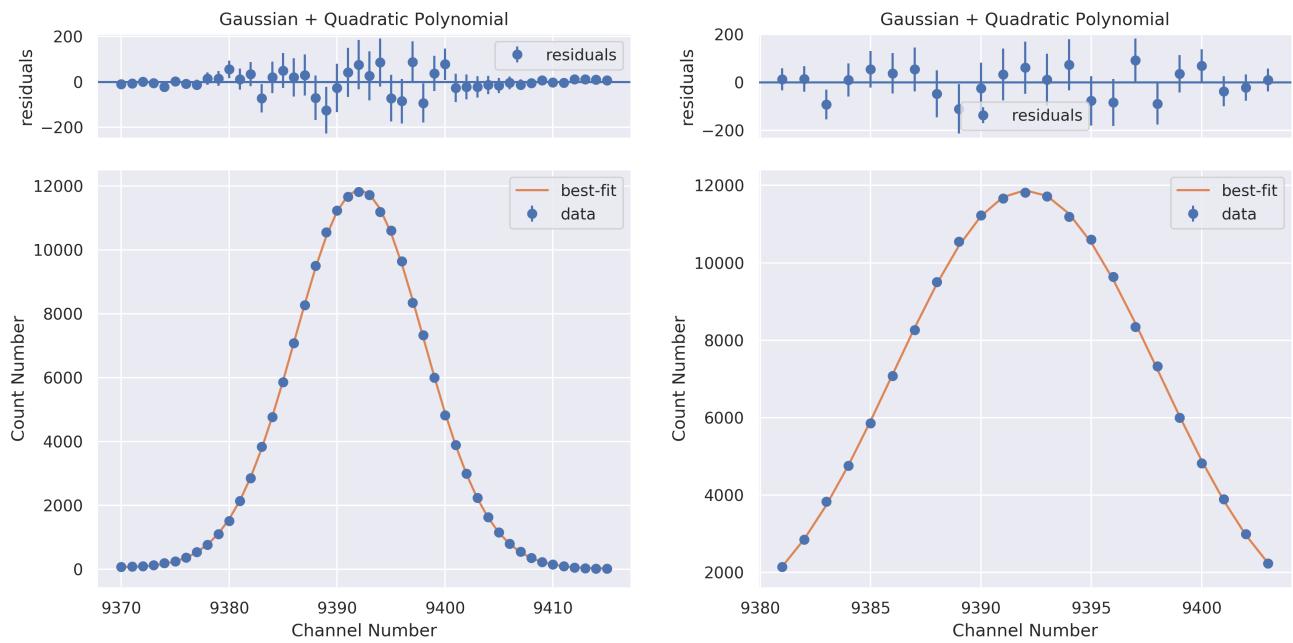


**Figure A.27:** Fit of full & zoomed in peak of  $^{60}\text{Co}$  1332 keV peak

### A1.2.3 GAUSSIAN + QUADRATIC FIT



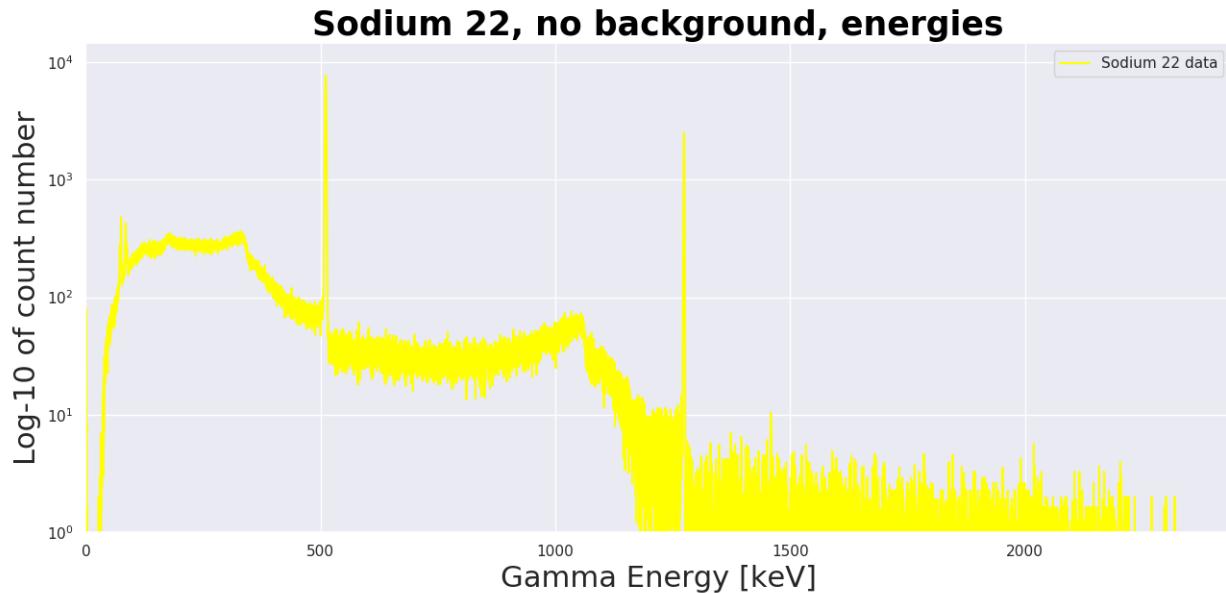
**Figure A.28:** Fit of full & zoomed in peak of  $^{60}\text{Co}$  1173 keV peak



**Figure A.29:** Fit of full & zoomed in peak of  $^{60}\text{Co}$  1332 keV peak

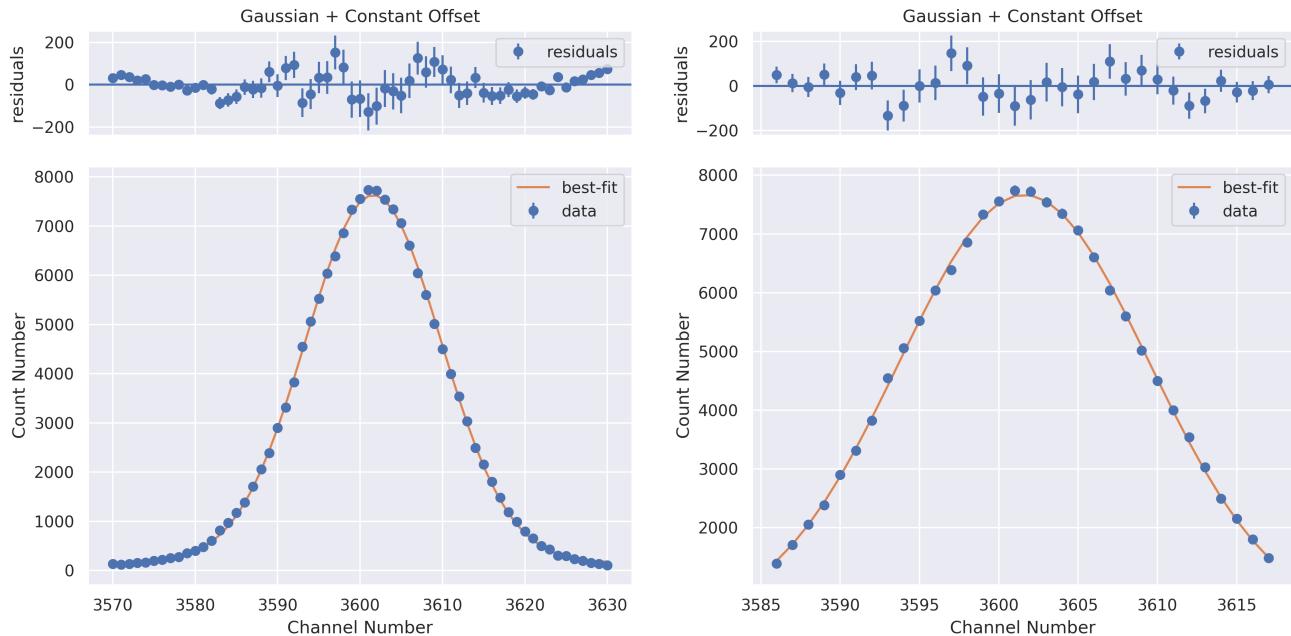
## A1.3 SODIUM-22

The  $^{22}\text{Na}$  source had one medium energy peak, and one high-energy peak. Both of which were very distinctive, and so quite easy to fit for all functions.



**Figure A.30:** The final spectrum obtained for the  $^{22}\text{Na}$  source run for 24 hours. This has been converted to energies by using the quadratic energy calibration.

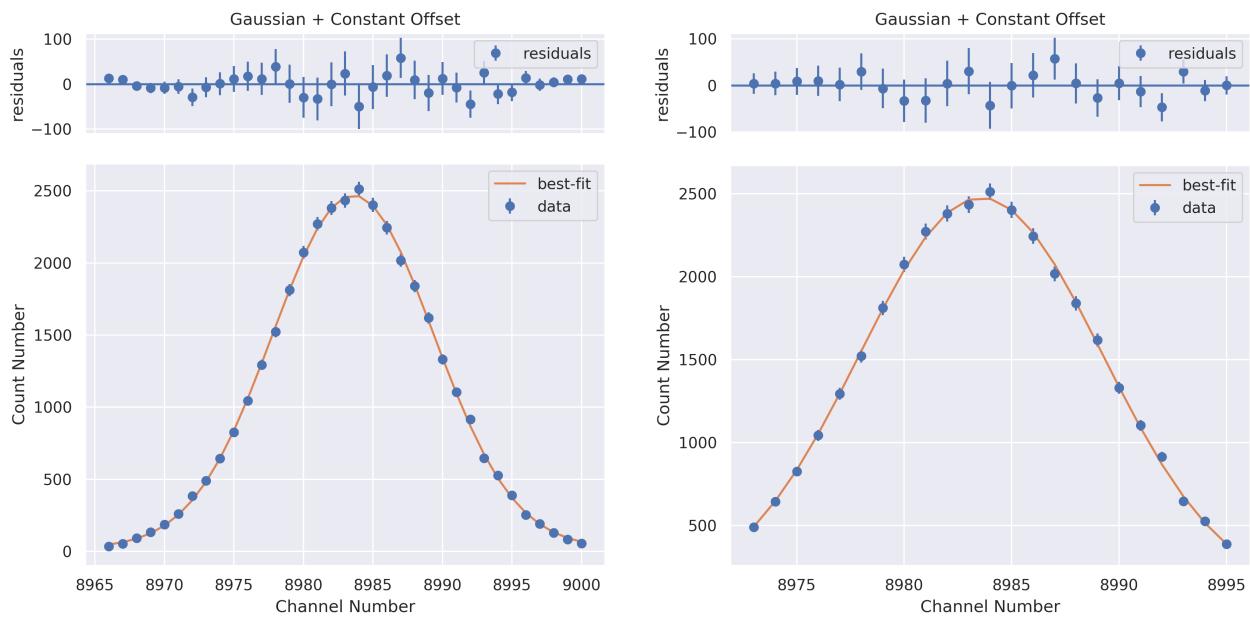
### A1.3.1 GAUSSIAN + OFFSET FIT



(a) Full peak with fit.  $\chi^2 = 202.42$ ,  $\chi^2_\nu = 3.43$ , Prob = 0.00%,  $\mu = 3601.60$ ,  $\sigma = 8.18$

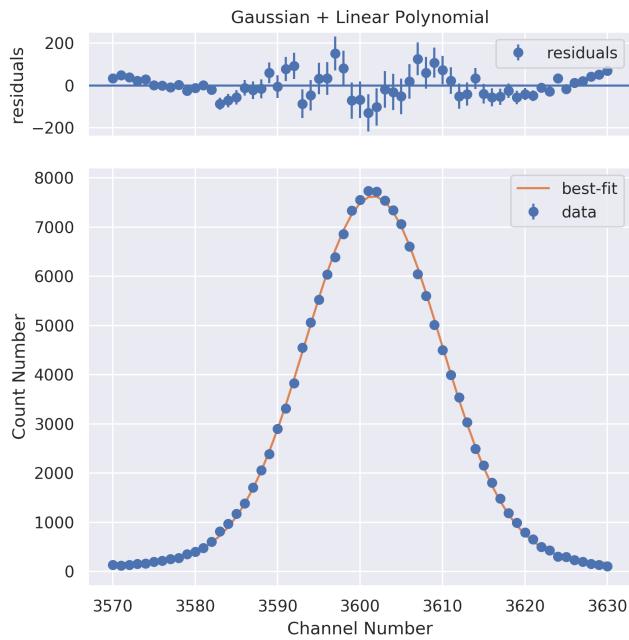
(b) Zoomed in peak with fit.  $\chi^2 = 24.72$ ,  $\chi^2_\nu = 0.82$ , Prob = 73.83%,  $\mu = 3601.60$ ,  $\sigma = 7.88$

**Figure A.31:** Fit of full & zoomed in peak of  $^{22}\text{Na}$  511 keV peak

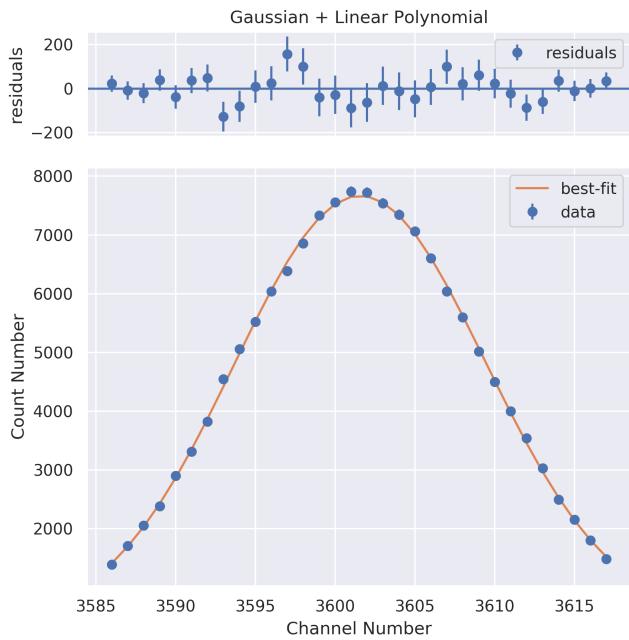


**Figure A.32:** Fit of full & zoomed in peak of  $^{22}\text{Na}$  1274 keV peak

### A1.3.2 GAUSSIAN + LINEAR FIT

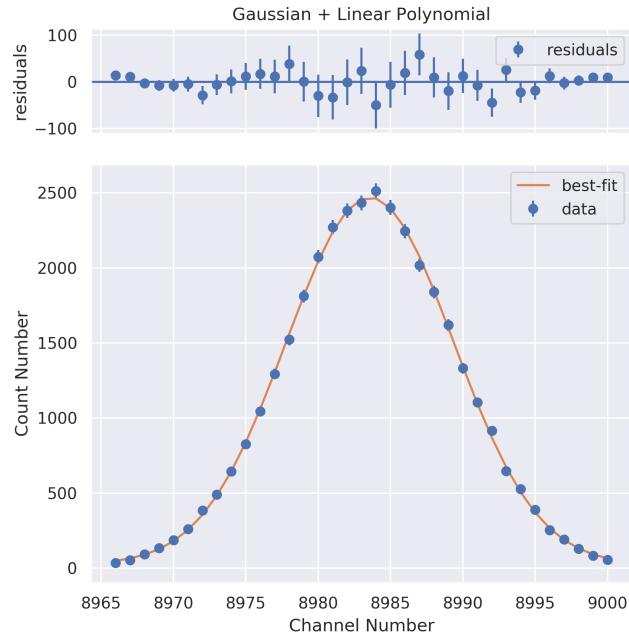


(a) Full peak with fit.  $\chi^2 = 197.63$ ,  $\chi^2_\nu = 3.35$ ,  
Prob = 0.00%,  $\mu = 3601.60$ ,  $\sigma = 8.18$

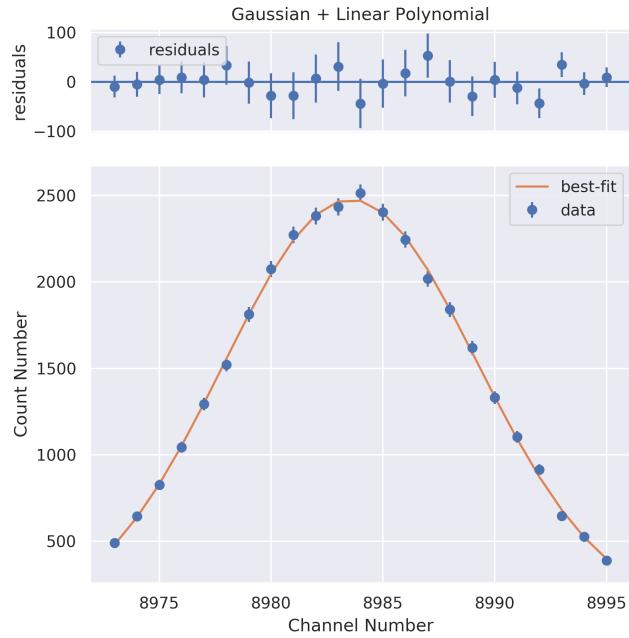


(b) Zoomed in peak with fit.  $\chi^2 = 22.76$ ,  $\chi^2_\nu = 0.76$ ,  
Prob = 82.51%,  $\mu = 3601.55$ ,  $\sigma = 7.88$

**Figure A.33:** Fit of full & zoomed in peak of  $^{60}\text{Na}$  511 keV peak



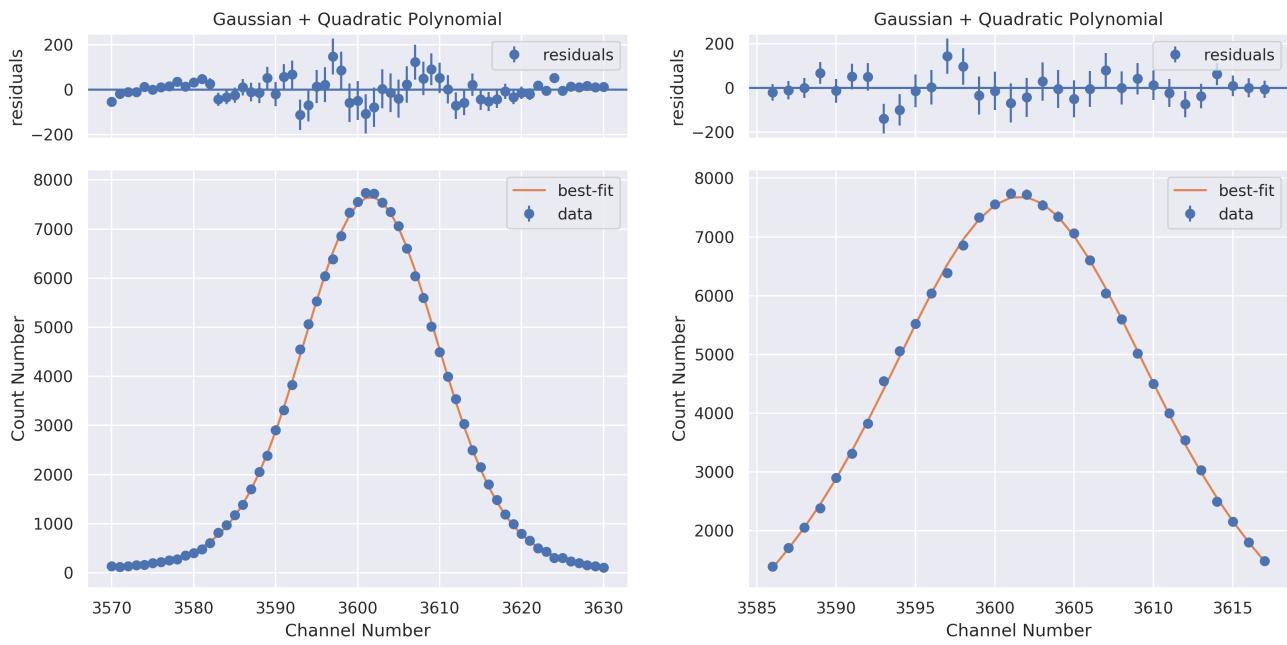
(a) Full peak with fit.  $\chi^2 = 26.34$ ,  $\chi^2_\nu = 0.80$ ,  
Prob = 78.77%,  $\mu = 8983.58$ ,  $\sigma = 5.79$



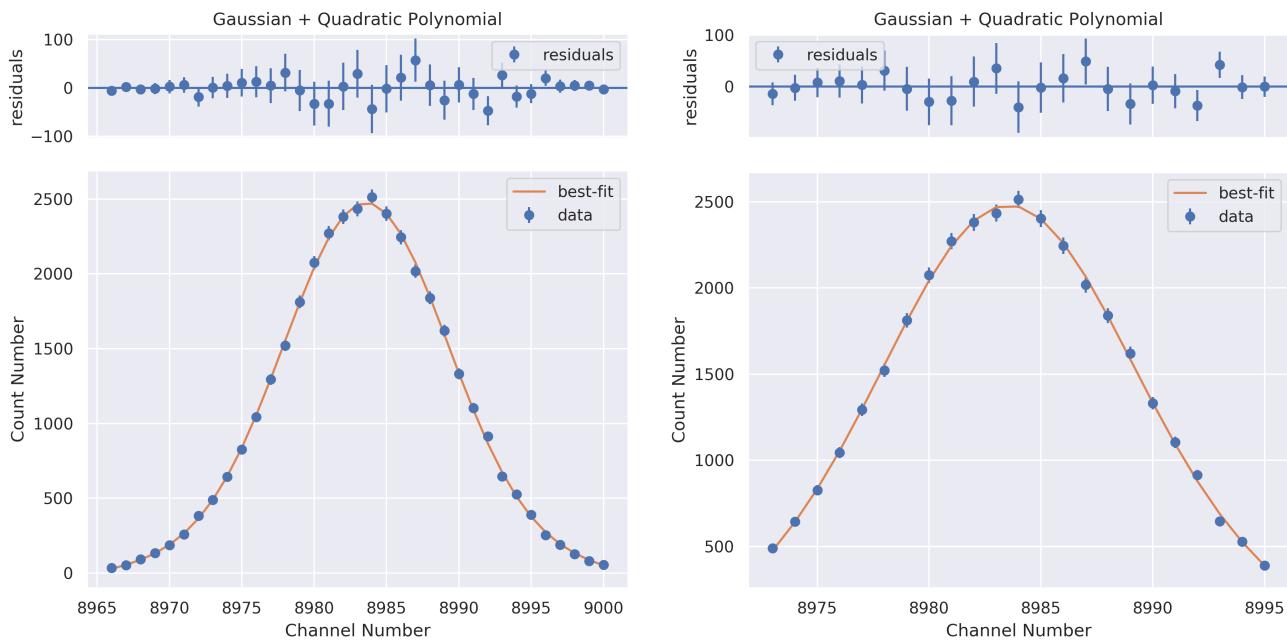
(b) Zoomed in peak with fit.  $\chi^2 = 9.41$ ,  $\chi^2_\nu = 0.45$ ,  
Prob = 98.56%,  $\mu = 8983.53$ ,  $\sigma = 5.67$

**Figure A.34:** Fit of full & zoomed in peak of  $^{22}\text{Na}$  1274 keV peak

### A1.3.3 GAUSSIAN + QUADRATIC FIT



**Figure A.35:** Fit of full & zoomed in peak of  $^{60}\text{Na}$  511 keV peak



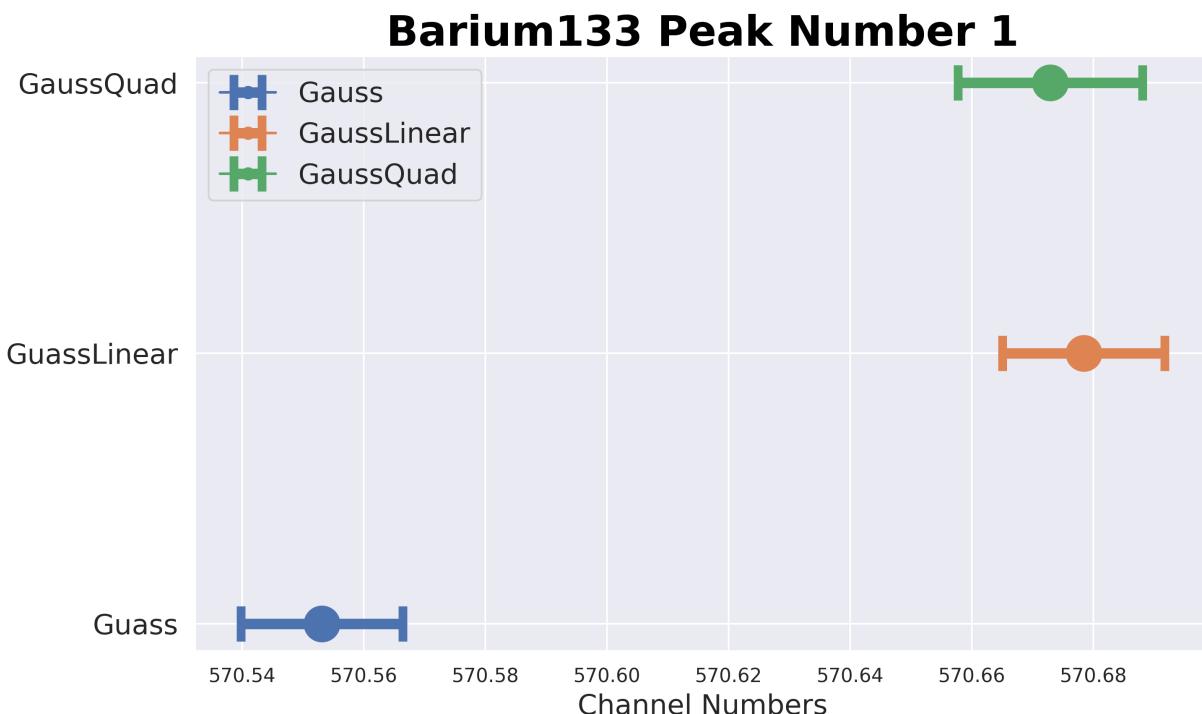
**Figure A.36:** Fit of full & zoomed in peak of  $^{22}\text{Na}$  1274 keV peak

## A2 FIT COMPARISONS

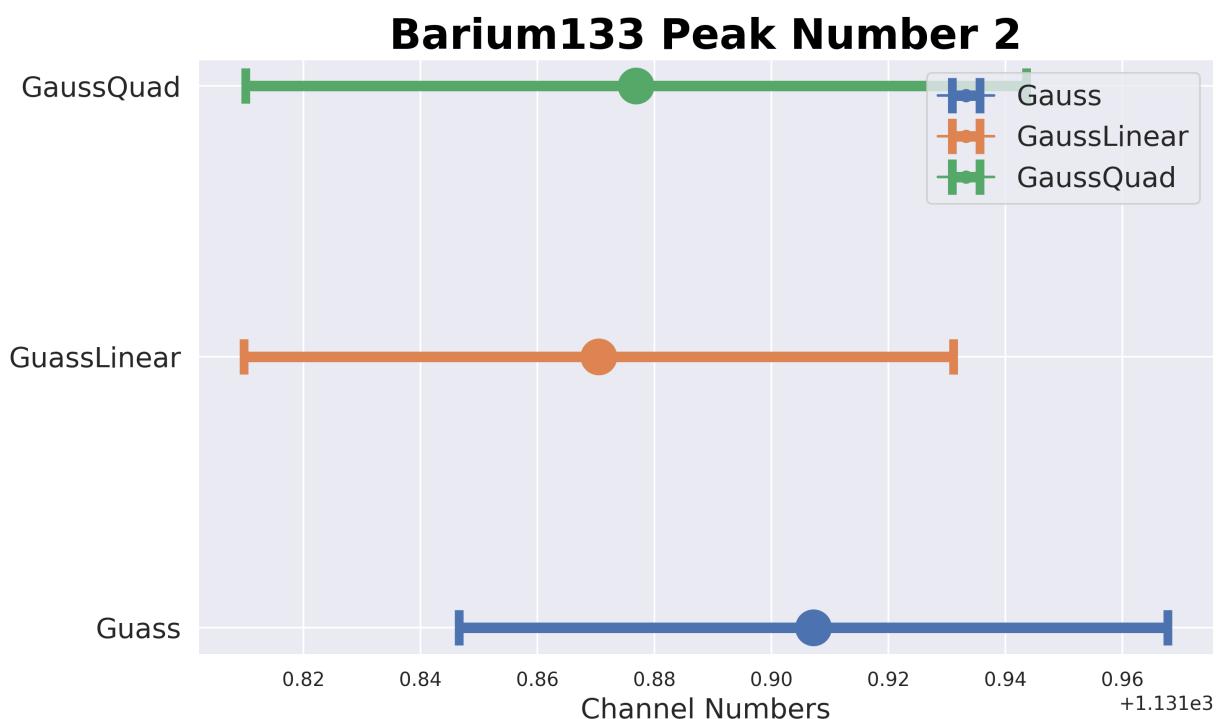
Since there were three types of function fitted to the data, it is important that we identify which function both matched the data the best, and that is the simplest function. From functional analysis, we know that any function can be represented as an arbitrary sum of polynomials, and so by increasing the order of the fitted function will naturally cause better fit statistics, even though there is less physical reason to do so. Therefore, to find the ideal fitting function, we wanted to identify the lowest-order polynomial that correctly fitted the data, by showing that when we increase the order the fit remains the same, within errors.

This can be represented graphically by plotting the means produced for the same peaks of the different fits, with the error bars representing the standard deviation on the mean from the fit. If two error bars from different fits are overlapping with the means, then this shows that these fits are compatible with each other, and so the lowest order should be taken as the true fit. However, if they do not overlap, then this shows that they are not compatible, and so the higher order fit should be taken.

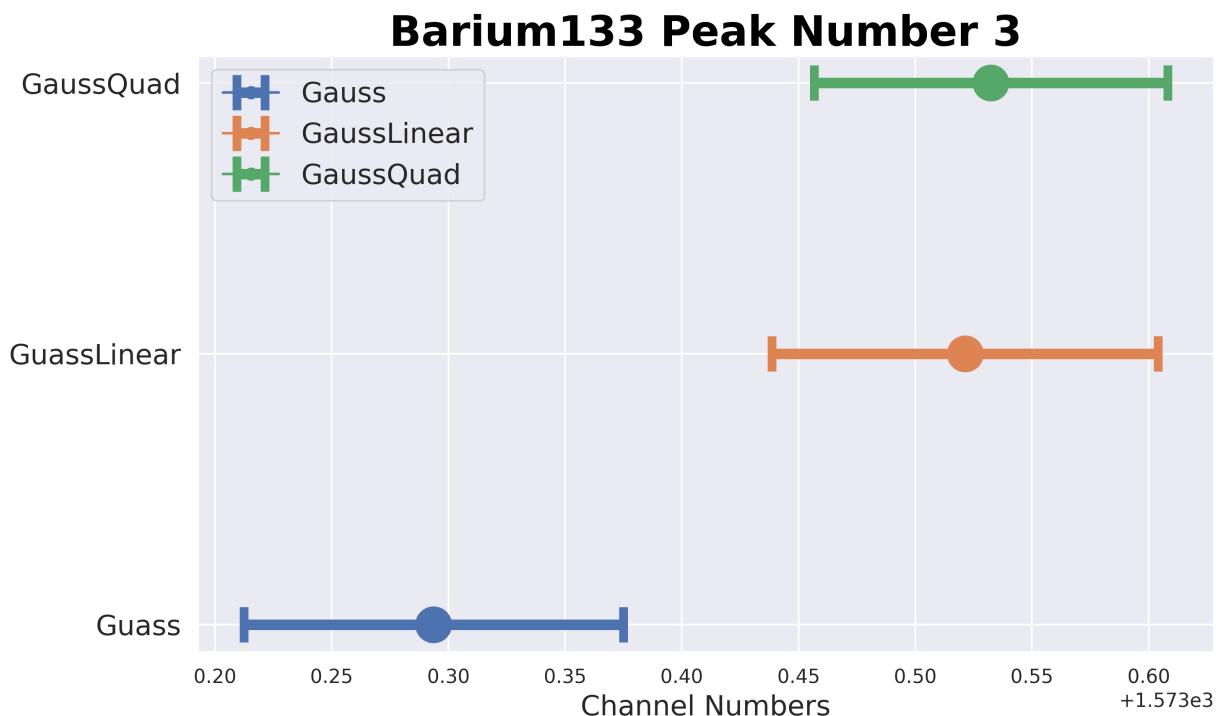
### A2.1 BARIUM-133



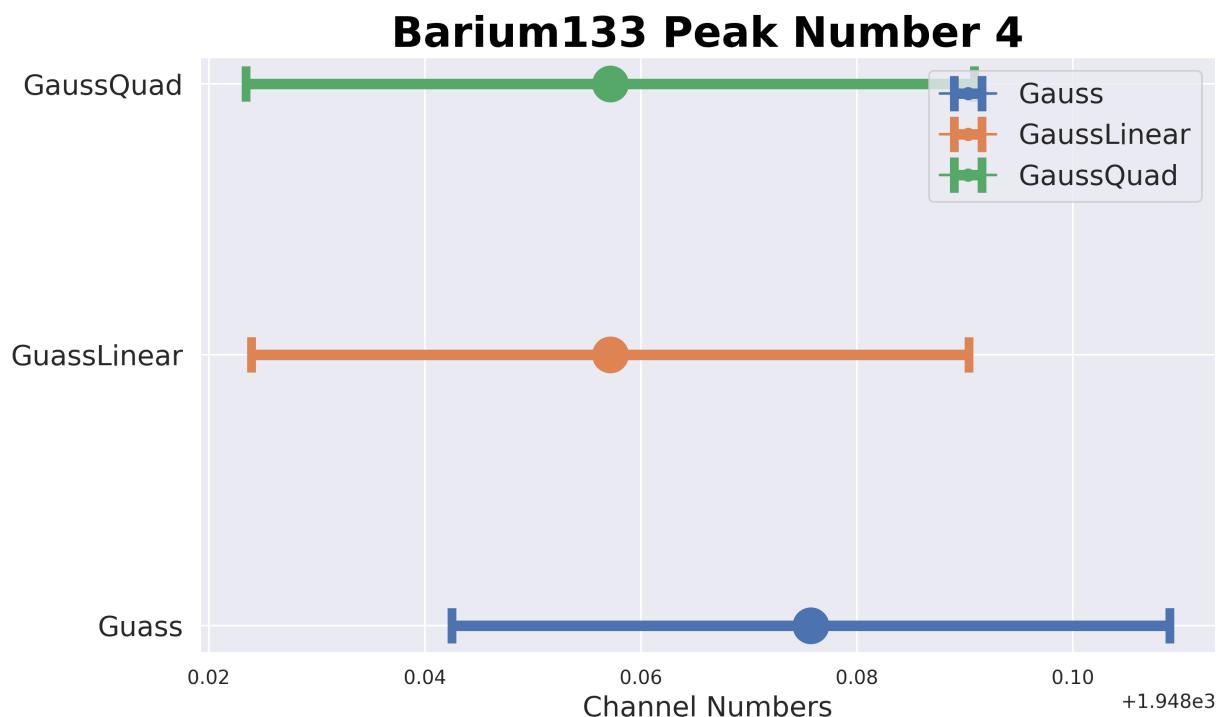
**Figure A.37:** Fit comparisons for  $^{133}\text{Ba}$  81 keV peak. Here is clear evidence that a simple Gaussian + offset fit is not sufficient, as the mean from the fit is significantly outside the acceptable area of the other two fits. This also shows that the linear polynomial provides no additional gain over the quadratic polynomial as both of their means coincide with each others error-bars, and so are statistically insignificant from each other.



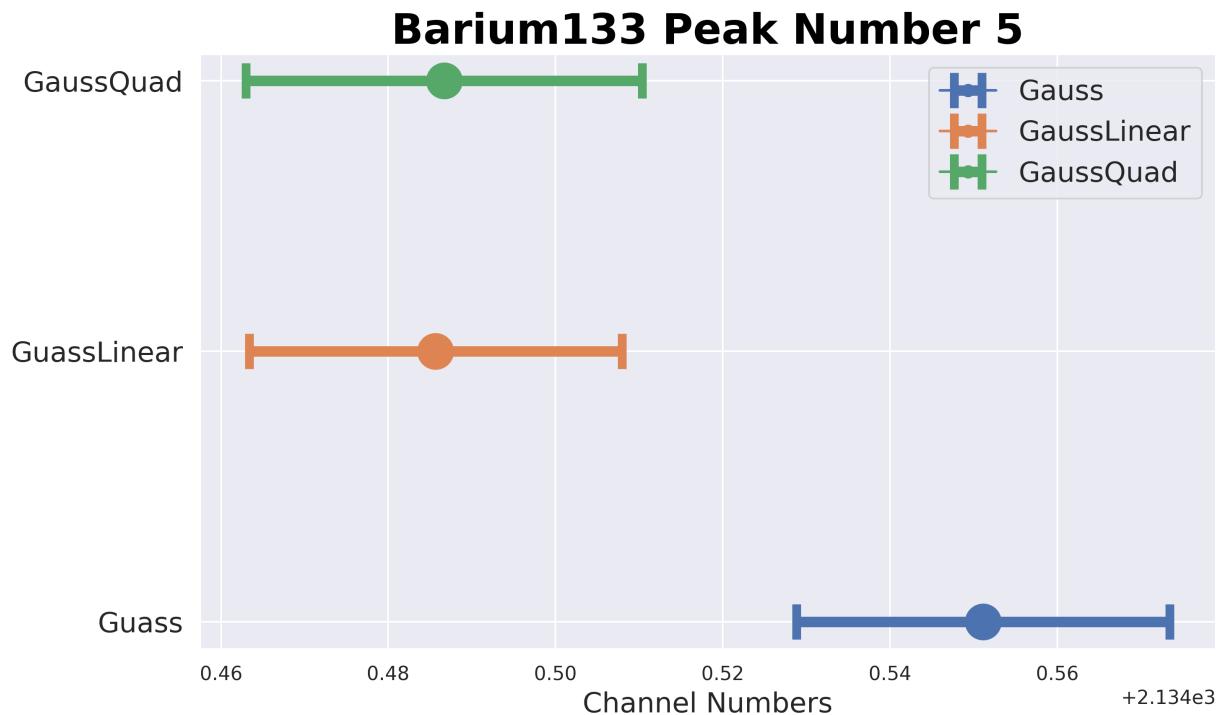
**Figure A.38:** Fit comparisons for  $^{133}\text{Ba}$  161 keV peak. Here, this shows all three means coincide with each other, and so for this particular peak going to anything above an offset has no advantage, however as other plots show that Gaussian + linear polynomial is best, that is what was used throughout.



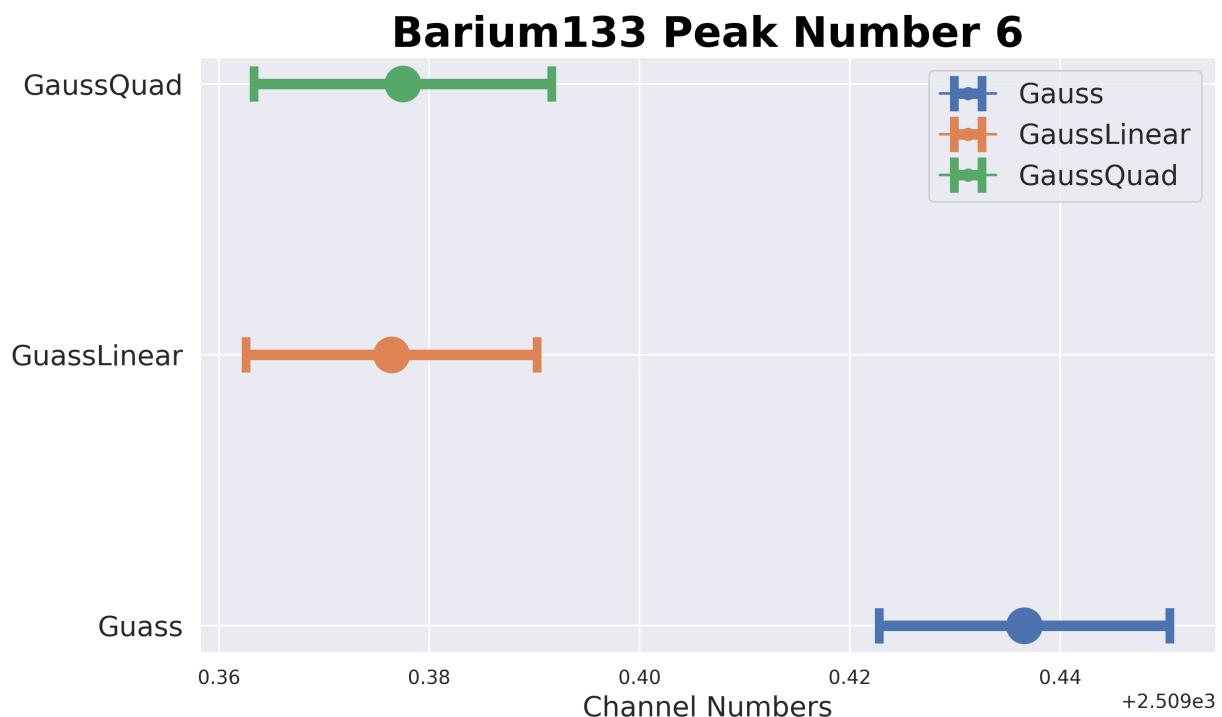
**Figure A.39:** Fit comparisons for  $^{133}\text{Ba}$  223 keV peak



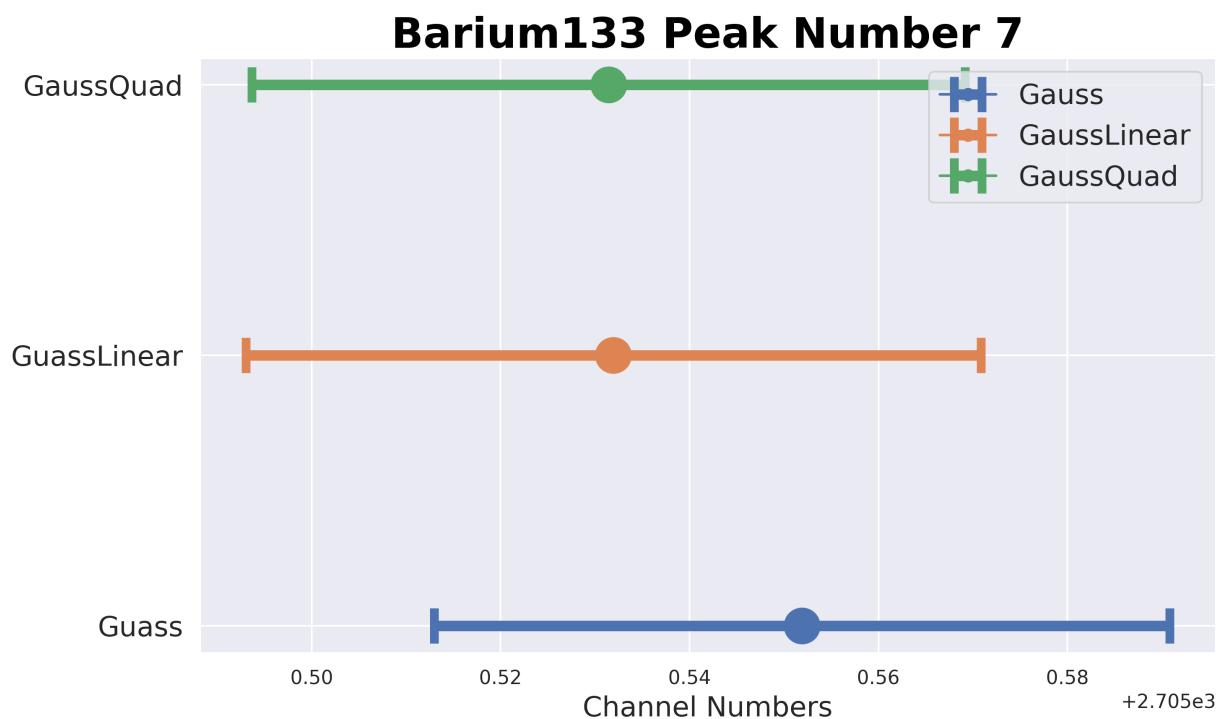
**Figure A.40:** Fit comparisons for  $^{133}\text{Ba}$  276 keV peak



**Figure A.41:** Fit comparisons for  $^{133}\text{Ba}$  303 keV peak

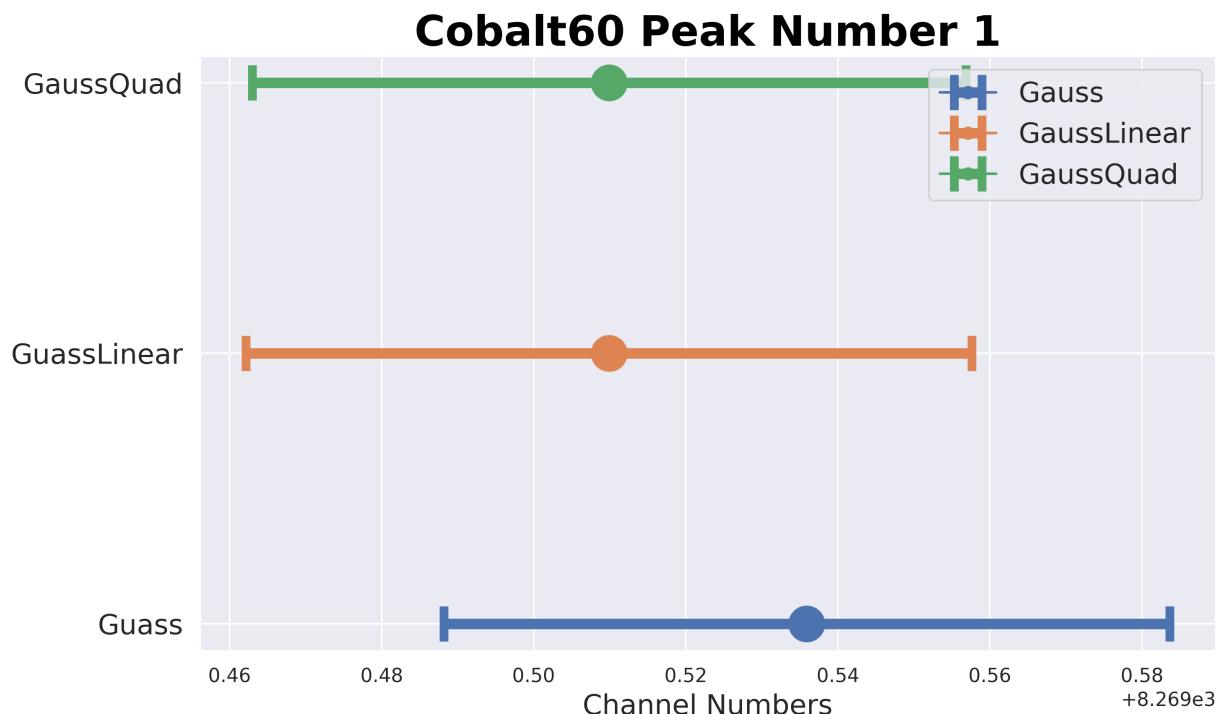


**Figure A.42:** Fit comparisons for  $^{133}\text{Ba}$  356 keV peak

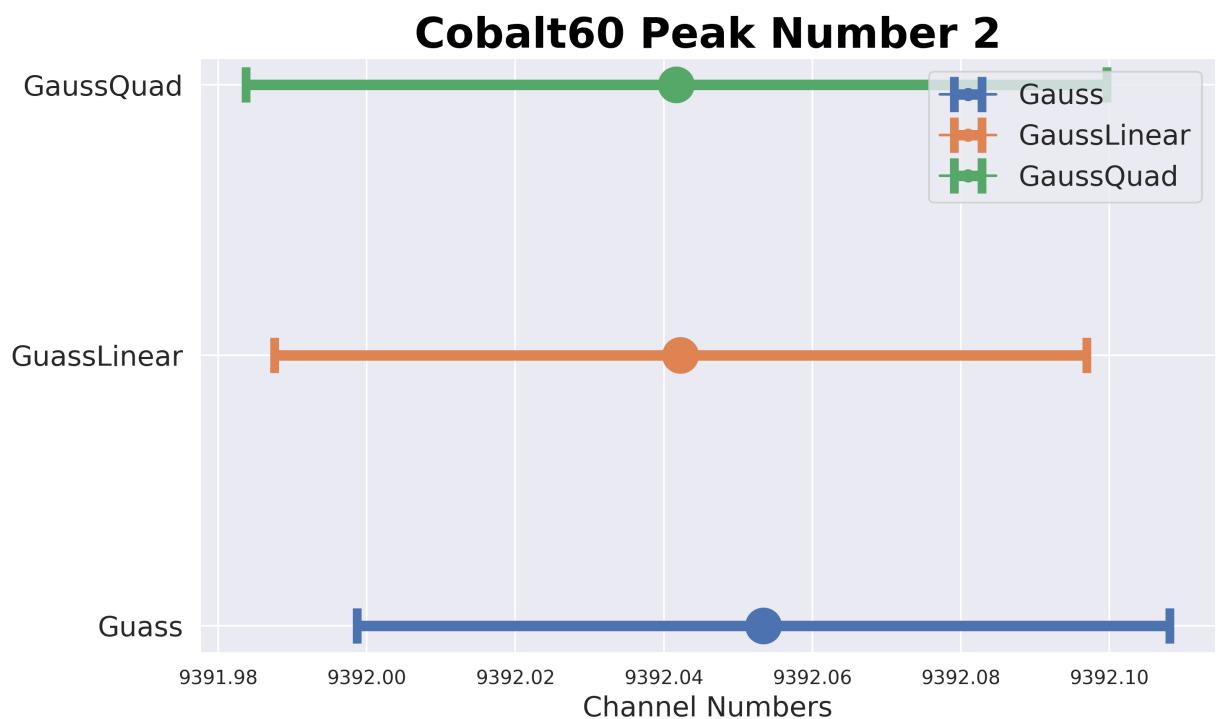


**Figure A.43:** Fit comparisons for  $^{133}\text{Ba}$  384 keV peak

## A2.2 COBALT-60

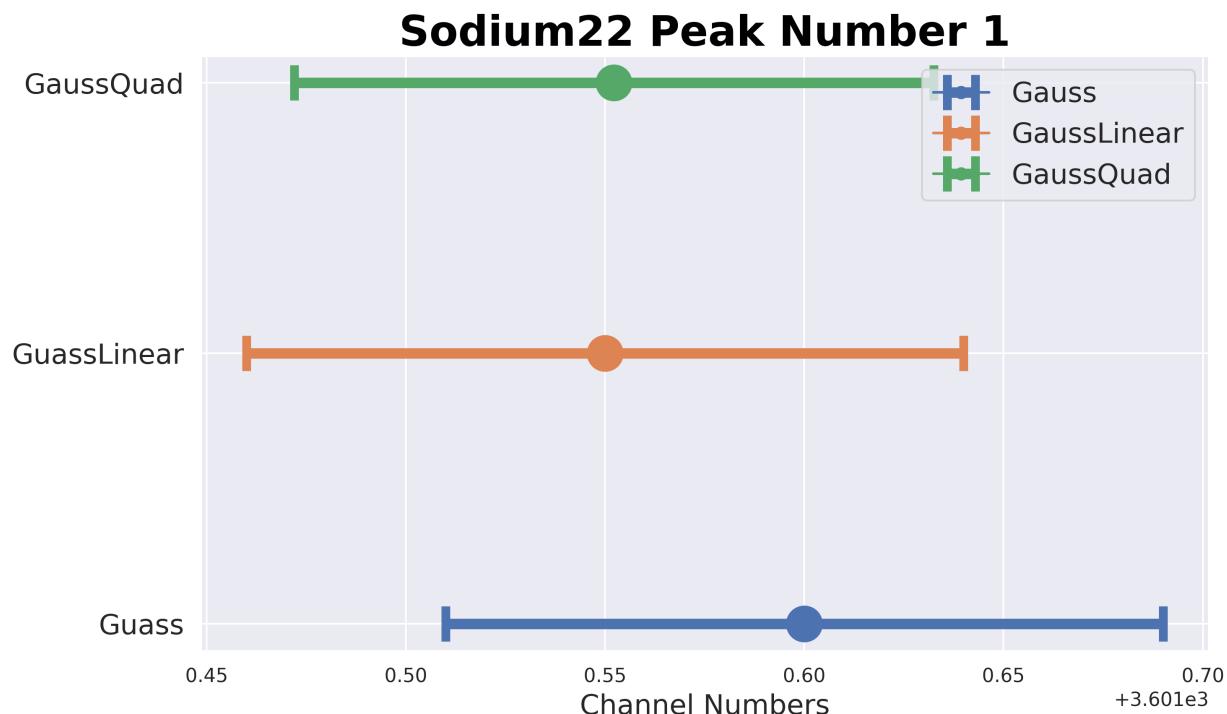


**Figure A.44:** Fit comparisons for  $^{60}\text{Co}$  1173 keV peak

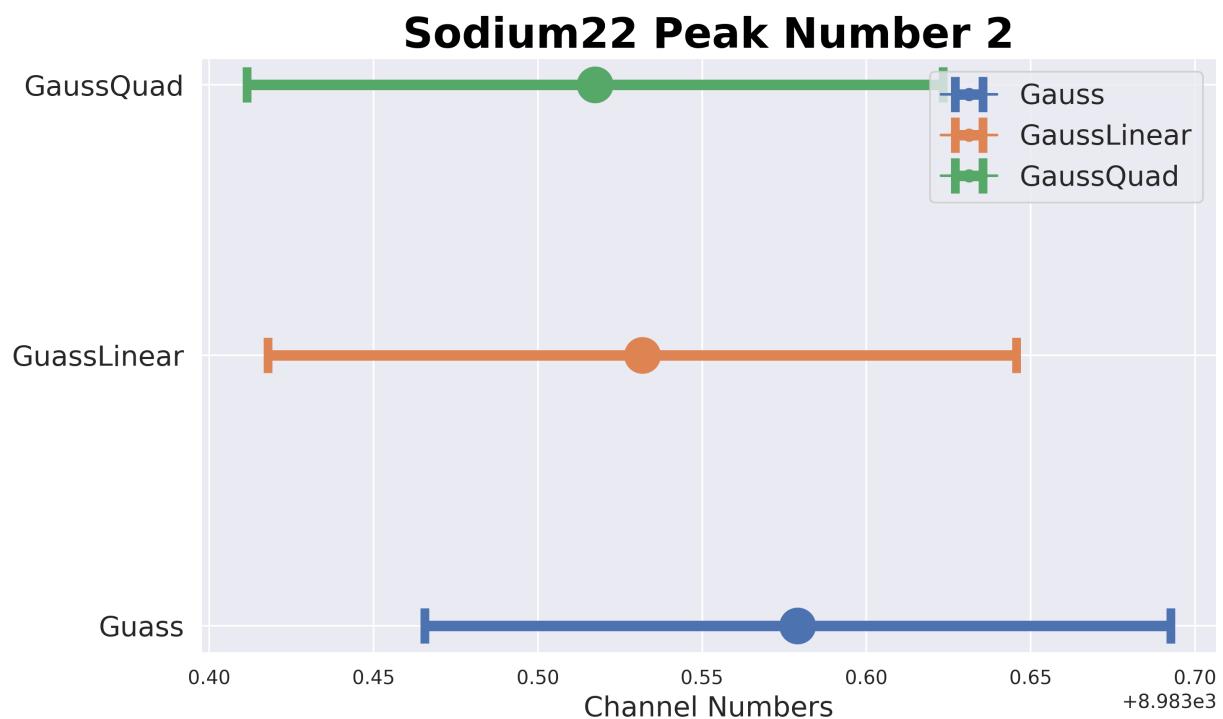


**Figure A.45:** Fit comparisons for  $^{60}\text{Co}$  1332 keV peak

## A2.3 SODIUM-22



**Figure A.46:** Fit comparisons for  $^{22}\text{Na}$  511 keV peak



**Figure A.47:** Fit comparisons for  $^{22}\text{Na}$  1274 keV peak

## A3 PYTHON CODE

Below is the complete collection of our Python codes that we wrote during the four weeks of our experiment to run various data analysis tools on the raw data, including curve fitting, energy calibration and detector resolution finding. They are included in this appendix so that way we have them on records, in case something tragic happens to the digital copies stored on the central university file system.

However, we have also mitigated against file loss by uploading our Python files, raw data, and this appendix to a GitHub repository that is hosted externally. In addition to being used as a version tracker tool helping us identify changes made to the codes over time (as we add, change & remove features), it also provides mitigation against file loss as the current builds can simply be downloaded from the repo again. This also provides an easy way for others to double check figures or results, if they so desire – improving transparency for our experiment, which is something that all scientists should be striving to do.

GitHub: <https://github.com/AlexMaraio/HPGeDetector>

### A3.1 GAUSSIANFIT.PY

This file is the file responsible for the fitting of the individual Gaussians to the peaks in the data. Here, we also add various order polynomials in addition to the Gaussian, starting from a simple offset to a quadratic order polynomial. This adjusts the way the fit works in subtle ways and we needed to see how fitting the various functions worked. Of course any arbitrary function can be approximated as a linear polynomial of arbitrary degree, so naturally going to a higher order polynomial will increase the goodness of fit, however detract from what can be extracted from the fit.

By looking how the mean changes between the different fit types we can determine if we are justified in increasing the order of the polynomial, and after our analysis thanks to the following code, we can not justify going beyond a linear polynomial in addition to the Gaussian function.

```

1 import numpy as np
2 import matplotlib.pyplot as plt
3 import seaborn as sns
4 sns.set()
5 import scipy.optimize as sciopt
6 import scipy.stats as scistats
7 from lmfit import Model, Parameters
8 import pandas as pd
9
10 def ChiSqFunc(Measured,Fitted,Errors):
11     ChiSquared = 0
12     for i in range(len(Measured)):
13         ChiSquared += ((Measured[i] - Fitted[i])**2.0) / ((Errors[i])**2.0)
14     ReducedChiSq = ChiSquared/(len(Measured)-2)
15     ProbChiSq = (1.0 - scistats.chi2.cdf(ChiSquared,len(Measured)-2) ) * 100.0
16     return ChiSquared, ReducedChiSq, ProbChiSq
17
18 def Gauss(x,A,mean,sigma,a):

```

```

19     return (A/(np.sqrt(2*np.pi) * sigma )) *np.exp(-((x-mean)**2.0)/(2.0*(sigma**2.0)))
20     ↵ + a
21
22 def GaussLinear(x,A,mean,sigma,a,b):
23     return (A/(np.sqrt(2*np.pi) * sigma ))*np.exp(-((x-mean)**2.0)/(2.0*(sigma**2.0)))
24     ↵ + a + b*x
25
26 def GaussQuad(x,A,mean,sigma,a,b,c):
27     return (A/(np.sqrt(2*np.pi) * sigma ))*np.exp(-((x-mean)**2.0)/(2.0*(sigma**2.0)))
28     ↵ + a + b*x + c*x**2.0
29
30
31 def ChToEnergy(Ch,a,b,c):
32     Energy = np.sqrt( ( (Ch - c + (b**2.0 / (4*a))) )/a ) - (b / (2*a))
33     return Energy
34
35 SetSigma = 2
36
37 BariumList = [ [[550,585,570,81],[1125,1140,1132,161],[1565,1582,1574,223],[1935,1960]
38     ↵ ,1948,276],[2118,2150,2135,303],[2493,2526,2509,356],[2688,2723,2705,384]] , []
39     ↵ ]
40 SodiumList = [ [ [3570,3630,3600,511] , [8966,9000,8984,1274]] , [] ]
41 CobaltList = [ [ [8250,8290,8270,1173] , [9370,9415,9392,1332]] , [] ]
42
43 SourceList = ["Barium133-24HrRun_006_eh_1","Sodium22-24HrData_002_eh_1","Cobalt60-24H
44     ↵ rData_004_eh_1"]
45
46 PlotResolution = 300
47
48 PeakNo = int(1)
49
50 datadict = {'Element':[],'Fit type':[], 'Peak number':[], 'Peak type':[], 'Energy
51     ↵ (keV)': [], 'Resolution':[], 'Min of range':[], 'Max of range':[], 'Mean':[],
52     ↵ 'A':[], 'Sigma':[], 'Error on mean':[], 'a':[], 'b':[], 'c':[], 'chisq':[],
53     ↵ 'Reduced chisq':[], 'Probchisq':[]}
54
55 # Fit parameters
56 a = 2.63977565e-06
57 b = 7.04767648
58 c = -1.91414134e-01
59
60 for source in SourceList:
61     if source == "Barium133-24HrRun_006_eh_1":
62         ElementList = BariumList
63         Element = "Barium133"
64     elif source == "Sodium22-24HrData_002_eh_1":
65         ElementList = SodiumList
66         Element = "Sodium22"
67     else:
68         ElementList = CobaltList
69         Element = "Cobalt60"
70
71     for item in ElementList:
72         for peak in item:

```

```

63      #! Initialises the source and ranges for run
64
65      if item == ElementList[0]:
66          PeakType = "Full"
67      else:
68          PeakType = "Zoom"
69
70      df = pd.read_csv(source + ".dat", sep = r"\s+", names = ['channel
    ↵   number','count number'])
71
72      df['count errors'] = np.sqrt(df['count number'])
73      df['count errors'] = df['count errors'].replace(0,1)
74
75      MinValue = peak[0]
76      MaxValue = peak[1]
77      MeanValue = peak[2]
78      PeakEnergy = peak[3]
79
80      data = df[(MinValue<=df['channel number']) & (df['channel number']<=MaxValue)]
81      #-----
82      #! Gaussian + Offset model only
83
84      GaussModel = Model(Gauss)
85      Params = Parameters()
86
87      Params.add('A',value=max(data['count number']),vary=True,min=0)
88      Params.add('mean',value=MeanValue,vary=True)
89      Params.add('sigma',value=1,vary=True)
90      Params.add('a',value=1,vary=True)
91
92      FitResult = GaussModel.fit(data['count number'],params=Params,x=data['channel
    ↵   number'])
93      TempList = [ FitResult.best_values['mean'] - SetSigma
    ↵   *FitResult.best_values['sigma'] , FitResult.best_values['mean'] + SetSigma
    ↵   *FitResult.best_values['sigma'] , FitResult.best_values['mean'] ,
    ↵   PeakEnergy ]
94      if item == ElementList[0]:
95          ElementList[1].append(TempList)
96      #print(FitResult.best_values)
97      FitResult.plot(yerr=data['count errors'],xlabel='Channel Number',ylabel='Count
    ↵   Number',title='Gaussian + Constant Offset')
98
99      ChiStats = ChiSqFunc(list(data['count
    ↵   number']),list(FitResult.best_fit),list(data['count errors']))
100
101     plt.tight_layout()
102     plt.savefig(f'Plots/{Element}/Gauss/Gauss_{PeakNo}_{PeakType}.png',
    ↵   format='png', dpi=PlotResolution)
103     plt.close('all')
104
105     CountMax1 = FitResult.best_values['A'] / ( FitResult.best_values['sigma'] *
    ↵   np.sqrt(2 * np.pi ) ) + FitResult.best_values['a']
106     ErrorOnMean1 = FitResult.best_values['sigma'] / np.sqrt(CountMax1)

```

```

107     datadict['Element'].append(Element)
108     datadict['Fit type'].append('Gauss')
109     datadict['Energy (keV)'].append(PeakEnergy)
110     FWHM_Energy = 2 * np.sqrt(2*np.log(2)) *
111         → ChToEnergy(FitResult.best_values['sigma'],a,b,c)
112     datadict['Resolution'].append( FWHM_Energy / PeakEnergy)
113     #datadict['Resolution'].append( ( 2 *
114         → np.sqrt(2*np.log(2))*FitResult.best_values['sigma'] - b ) / a ) /
115         → PeakEnergy)
116     datadict['Peak number'].append(PeakNo)
117     datadict['Peak type'].append(PeakType)
118     datadict['Min of range'].append(MinValue)
119     datadict['Max of range'].append(MaxValue)
120     datadict['Mean'].append(FitResult.best_values['mean']))
121     datadict['A'].append(FitResult.best_values['A'])
122     datadict['Sigma'].append(FitResult.best_values['sigma'])
123     datadict['Error on mean'].append(1)
124     datadict['a'].append(FitResult.best_values['a'])
125     datadict['b'].append(0)
126     datadict['c'].append(0)
127     datadict['chisq'].append(ChiStats[0])
128     datadict['Reduced chisq'].append(ChiStats[1])
129     datadict['Probchisq'].append(ChiStats[2])
130
131     #-----#
132     #! Gaussian + Linear model
133
134     GaussModel2 = Model(GaussLinear)
135     Params2 = Parameters()
136
137     Params2.add('A',value=max(data['count number']),vary=True,min=0)
138     Params2.add('mean',value=MeanValue,vary=True)
139     Params2.add('sigma',value=1,vary=True)
140     Params2.add('a',value=1,vary=True)
141     Params2.add('b',value=1,vary=True)
142
143     FitResult2 = GaussModel2.fit(data['count
144         → number'],params=Params2,x=data['channel number'])
145     FitResult2.plot(yerr=data['count errors'],xlabel='Channel Number',ylabel='Count
146         → Number',title='Gaussian + Linear Polynomial')
147
148     ChiStats2 = ChiSqFunc(list(data['count
149         → number']),list(FitResult2.best_fit),list(data['count errors']))
150
151     plt.tight_layout()
152     plt.savefig(f'Plots/{Element}/Linear/Linear_{PeakNo}_{PeakType}.png',
153         → format='png', dpi=PlotResolution)
154     plt.close('all')
155
156     CountMax2 = FitResult2.best_values['A'] / ( FitResult2.best_values['sigma'] *
157         → np.sqrt(2 * np.pi ) ) + FitResult2.best_values['a'] +
158         → FitResult2.best_values['b'] * FitResult2.best_values['mean']

```

```

152     ErrorOnMean2 = FitResult2.best_values['sigma'] / np.sqrt(CountMax2)
153
154     datadict['Element'].append(Element)
155     datadict['Fit type'].append('Linear')
156     datadict['Energy (keV)'].append(PeakEnergy)
157     FWHM_Energy = 2 * np.sqrt(2*np.log(2)) *
158         → ChToEnergy(FitResult2.best_values['sigma'],a,b,c)
159     datadict['Resolution'].append( FWHM_Energy / PeakEnergy)
160     datadict['Peak number'].append(PeakNo)
161     datadict['Peak type'].append(PeakType)
162     datadict['Min of range'].append(MinValue)
163     datadict['Max of range'].append(MaxValue)
164     datadict['Mean'].append(FitResult2.best_values['mean'])
165     datadict['A'].append(FitResult2.best_values['A'])
166     datadict['Sigma'].append(FitResult2.best_values['sigma'])
167     datadict['Error on mean'].append(1)
168     datadict['a'].append(FitResult2.best_values['a'])
169     datadict['b'].append(FitResult2.best_values['b'])
170     datadict['c'].append(0)
171     datadict['chisq'].append(ChiStats2[0])
172     datadict['Reduced chisq'].append(ChiStats2[1])
173     datadict['Probchisq'].append(ChiStats2[2])
174
175 #-----
176 #! Gaussian + Quadratic model
177
178 GaussModel3 = Model(GaussQuad)
179 Params3 = Parameters()
180
181 Params3.add('A',value=max(data['count number']),vary=True,min=0)
182 Params3.add('mean',value=MeanValue,vary=True)
183 Params3.add('sigma',value=1,vary=True)
184 Params3.add('a',value=1,vary=True)
185 Params3.add('b',value=1,vary=True)
186 Params3.add('c',value=1,vary=True)
187
188 FitResult3 = GaussModel3.fit(data['count
189     → number'],params=Params3,x=data['channel number'])
190 FitResult3.plot(yerr=data['count errors'],xlabel='Channel Number',ylabel='Count
191     → Number',title='Gaussian + Quadratic Polynomial')
192
193 ChiStats3 = ChiSqFunc(list(data['count
194     → number']),list(FitResult3.best_fit),list(data['count errors']))
195
196 plt.tight_layout()
197 plt.savefig(f'Plots/{Element}/Quad/Quad_{PeakNo}_{PeakType}.png', format='png',
198     → dpi=PlotResolution)
199 plt.close('all')
200
201 CountMax3 = FitResult3.best_values['A'] / ( FitResult3.best_values['sigma'] *
202     → np.sqrt(2 * np.pi ) ) + FitResult3.best_values['a'] +
203     → FitResult3.best_values['b'] * FitResult3.best_values['mean'] +
204     → FitResult3.best_values['c'] * (FitResult3.best_values['mean'])** 2.0

```

```

197     ErrorOnMean3 = FitResult3.best_values['sigma'] / np.sqrt(CountMax3)
198
199     datadict['Element'].append(Element)
200     datadict['Fit type'].append('Quad')
201     datadict['Energy (keV)'].append(PeakEnergy)
202     FWHM_Energy = 2 * np.sqrt(2*np.log(2)) *
203         → ChToEnergy(FitResult3.best_values['sigma'],a,b,c)
204     datadict['Resolution'].append( FWHM_Energy / PeakEnergy)
205     datadict['Peak number'].append(PeakNo)
206     datadict['Peak type'].append(PeakType)
207     datadict['Min of range'].append(MinValue)
208     datadict['Max of range'].append(MaxValue)
209     datadict['Mean'].append(FitResult3.best_values['mean'])
210     datadict['A'].append(FitResult3.best_values['A'])
211     datadict['Sigma'].append(FitResult3.best_values['sigma'])
212     datadict['Error on mean'].append(1)
213     datadict['a'].append(FitResult3.best_values['a'])
214     datadict['b'].append(FitResult3.best_values['b'])
215     datadict['c'].append(FitResult3.best_values['c'])
216     datadict['chisq'].append(ChiStats3[0])
217     datadict['Reduced chisq'].append(ChiStats3[1])
218     datadict['Probchisq'].append(ChiStats3[2])
219
220
221     Delta12 = np.abs(FitResult.best_values['mean'] - FitResult2.best_values['mean'])
222     ErrorDelta12 = np.sqrt( ErrorOnMean1**2.0 + ErrorOnMean2**2.0 )
223
224     Delta23 = np.abs(FitResult2.best_values['mean'] -
225         → FitResult3.best_values['mean'])
226     ErrorDelta23 = np.sqrt( ErrorOnMean2**2.0 + ErrorOnMean3**2.0 )
227
228     if PeakType == "Zoom":
229         if abs(ErrorDelta23) > abs(Delta23):
230             print('best fit')
231         else:
232             print('not best fit')
233
234     TitleFont = {'size':'24', 'color':'black', 'weight':'bold'}
235     AxTitleFont = {'size':'16'}
236     plt.figure(figsize=(10,6))
237     plt.errorbar(FitResult.best_values['mean'],1,xerr=ErrorOnMean1,elinewidth=6,c_
238         → apsize=10,capthick=5,marker='o',markersize=20,label="Gauss")
239     plt.errorbar(FitResult2.best_values['mean'],1.25,xerr=ErrorOnMean2,elinewidth_
240         → =6,capsize=10,capthick=5,marker='o',markersize=20,label="GaussLinear")
241     plt.errorbar(FitResult3.best_values['mean'],1.5,xerr=ErrorOnMean3,elinewidth=_
242         → 6,capsize=10,capthick=5,marker='o',markersize=20,label="GaussQuad")
243     plt.legend(fontsize=16,markerscale=0.33)
244     plt.title(str(Element) + ' Peak Number ' + str(PeakNo),**TitleFont)
245     plt.xlabel('Channel Numbers',**AxTitleFont)
246     plt.yticks([1,1.25,1.5],['Guass','GuassLinear','GaussQuad'],**AxTitleFont)
247     plt.tight_layout()
248     plt.savefig(f'Plots/{Element}/FitComparison_Peak{PeakNo}.png', format='png',
249         → dpi=PlotResolution)

```

```
245     plt.close('all')
246
247     PeakNo +=1
248
249     PeakNo = 1
250
251 Fitdf = pd.DataFrame(datadict)
252 Fitdf.to_csv('Gamma_Peak_Stats_and_Params.csv')
```

## A3.2 CALIBRATION.PY

This Python file was responsible for reading in the fitted Gaussians that were generated by `GaussainFit.py` and turning them into the fit between channel number and energy. First, we used a linear fit as it was assumed that the equipment would have a linear response, however, it was observed that for large energies, there was a large discrepancy between our measured value and the predicted value for the peaks. This lead us to believe that the energy calibration might have to be done with a quadratic function, not just a simple straight line.

When we performed this new quadratic fit, we found that the energies matched much more closely to what we expected them to be, and so we can determine that the system overall has a non-linear response between energy and channel number.

```

1 import numpy as np
2 import matplotlib.pyplot as plt
3 import scipy.optimize as sciopt
4 import seaborn as sns
5 import pandas as pd
6 sns.set()
7 import scipy.stats as scistats
8
9 TitleFont = {'size':'25', 'color':'black', 'weight':'bold'}
10 AxTitleFont = {'size':'22'}
11
12 def Linear(E,a,b):
13     return E*a + b
14
15 def Quadratic(E,a,b,c):
16     return a*E**2.0 + E*b + c
17
18 def ChiSqFunc(Measured,Fitted,Errors,Params):
19     ChiSquared = 0
20     for i in range(len(Measured)):
21         ChiSquared += ((Measured[i] - Fitted[i])**2.0) / ((Errors[i])**2.0)
22     ReducedChiSq = ChiSquared/(len(Measured)-len(Params))
23     ProbChiSq = (1.0 - scistats.chi2.cdf(ChiSquared,len(Measured)-len(Params))) *
24     ↪ 100.0
25     return ChiSquared, ReducedChiSq, ProbChiSq
26
27 #? Imports the csv to a Pandas DataFrame
28 FitDF = pd.read_csv('Gamma_Peak_Stats_and_Params.csv', names=['Element', 'Fit type',
29 ↪ 'Peak number', 'Peak type', 'Energy (keV)', 'Resolution', 'Min', 'Max', 'Mean',
30 ↪ 'A', 'Sigma', 'Error', 'a', 'b', 'chisq', 'Reduced chisq',
31 ↪ 'Probchisq'],usecols=list(range(1,18)))
32 #? Ensures that only the linear and zoomed peaks are fitted
33 FitDF = FitDF[FitDF['Fit type'] == 'Linear']
34 FitDF = FitDF[FitDF['Peak type'] == 'Zoom']
35 #? Puts the DF as floats otherwise it crashes
36 FitDF['Mean'] = FitDF['Mean'].astype(float)
37 FitDF['Error'] = FitDF['Error'].astype(float)
38 FitDF['Energy (keV)'] = FitDF['Energy (keV)'].astype(float)

```

```

36 DataBa = FitDF[FitDF['Element'] == 'Barium133']
37 DataCo = FitDF[FitDF['Element'] == 'Cobalt60']
38 DataNa = FitDF[FitDF['Element'] == 'Sodium22']
39
40 #? Fits the quadratic to the data
41 ParamsLinear, ErrorsLinear = sciopt.curve_fit(Quadratic,FitDF['Energy
   ↵ (keV)'],FitDF['Mean'])
42 print(ParamsLinear)
43
44 a = ParamsLinear[0]
45 b = ParamsLinear[1]
46 c = ParamsLinear[2]
47
48 Energies = np.sqrt( (FitDF['Mean'] - c + (b**2.0/ (4 * a)))/a ) - (b/(2*a))
49 #Energies = (FitDF['Mean'] - b)/(a) # keV
50
51 #? Calibration Plot
52 plt.errorbar(DataBa['Energy (keV)'],DataBa['Mean'], fmt='ro',label="Barium-133",
   ↵ markersize=10.5,yerr=DataBa['Error'].values)
53 plt.errorbar(DataCo['Energy (keV)'],DataCo['Mean'],fmt='bo',label="Cobalt-60",
   ↵ markersize=10.5,yerr=DataCo['Error'].values)
54 plt.errorbar(DataNa['Energy (keV)'],DataNa['Mean'],fmt='yo',label="Sodium-22",
   ↵ markersize=10.5,yerr=DataNa['Error'].values)
55 plt.plot(Energies,Quadratic(Energies,*ParamsLinear),label='Fit')
56 plt.title('Energy calibration plot with fit',**TitleFont)
57 plt.xlabel('Gamma Energy [keV]',**AxTitleFont)
58 plt.ylabel('Channel Number',**AxTitleFont)
59 plt.legend()
60 plt.show()
61
62 #? Residual plot
63 plt.errorbar(DataBa['Energy (keV)'],[(DataBa['Mean'].iloc[i] -
   ↵ Quadratic(DataBa['Energy (keV)'].iloc[i],*ParamsLinear)) for i in
   ↵ range(len(DataBa['Energy (keV)']))],fmt='ro',label="Barium-133",
   ↵ markersize=7.5,yerr=DataBa['Error'].values)
64 plt.errorbar(DataCo['Energy (keV)'],[(DataCo['Mean'].iloc[i] -
   ↵ Quadratic(DataCo['Energy (keV)'].iloc[i],*ParamsLinear)) for i in
   ↵ range(len(DataCo['Energy (keV)']))],fmt='bo',label="Cobalt-60",
   ↵ markersize=7.5,yerr=DataCo['Error'].values)
65 plt.errorbar(DataNa['Energy (keV)'],[(DataNa['Mean'].iloc[i] -
   ↵ Quadratic(DataNa['Energy (keV)'].iloc[i],*ParamsLinear)) for i in
   ↵ range(len(DataNa['Energy (keV)']))],fmt='yo',label="Sodium-22",
   ↵ markersize=7.5,yerr=DataNa['Error'].values)
66 plt.title('Energy calibration residual plot',**TitleFont)
67 plt.xlabel('Gamma Energy [keV]',**AxTitleFont)
68 plt.ylabel('Channel Number Residual',**AxTitleFont)
69 plt.legend()
70 plt.show()
71
72 ChiStats = ChiSqFunc(list(FitDF['Mean']),list(Quadratic(FitDF['Energy
   ↵ (keV)'],*ParamsLinear)),list(FitDF['Error']),ParamsLinear)
73 print(ChiStats)

```

### A3.3 SUBTRACTION.PY

This Python script was responsible for reading in the data files for the three known sources and the mystery source and subtracting the background that was there in all spectra. This background had to be correctly normalised to the same time period as the sources were taken for, so that way the correct amount of background is subtracted.

```

1 import numpy as np
2 import matplotlib.pyplot as plt
3 import seaborn as sns
4 sns.set()
5 import scipy as sc
6 import pandas as pd
7
8 # If statements that control what the code does!
9 PlotMysterySource = True
10 PlotKnownSources = True
11
12 # Plot controls
13 TitleFont = {'size':'23', 'color':'black', 'weight':'bold'}
14 AxTitleFont = {'size':'20'}
15
16 if PlotMysterySource:
17     BgWithShield = "WeekendBgWithShield_003_eh_1"
18     BgWithShieldDF = pd.read_csv(BgWithShield+ ".dat", sep=r"\s+", names = ['channel
19     ↵ number','count number'])
20
21 MysterySource = "MysterySource-24HrRun_001_eh_1"
22 MysterySourceDF = pd.read_csv(MysterySource+ ".dat", sep=r"\s+", names = ['channel
23     ↵ number','count number'])
24
25 # Normalises the timings of the two runs to 24 hours each
26 BgWithShieldDF['count number2'] = BgWithShieldDF['count number']
27 BgWithShieldDF['count number'] = BgWithShieldDF['count number'] / ( 239068 ) *
28     ↵ 86400
29
30 MysterySourceDF['count number'] = MysterySourceDF['count number'] -
31     ↵ BgWithShieldDF['count number']
32
33 # Fit parameters
34 a = 2.63977565e-06
35 b = 7.04767648
36 c = -1.91414134e-01
37
38 # Old fit params for comparison
39 aold = 7.012
40 bold = 8.170
41
42 MysterySourceDF['Energies'] = np.sqrt( (MysterySourceDF['channel number'] - c +
43     ↵ (b**2.0/ (4 * a))/a ) - (b/(2*a)))
44 MysterySourceDF['Energies2'] = (MysterySourceDF['channel number'] -bold)/aold

```

```

41     BgWithShieldDF['Energies'] = np.sqrt( (BgWithShieldDF['channel number'] - c +
42                                         ↵ (b**2.0/ (4 * a)))/a ) - (b/(2*a))
43
44     # Plotting the mystery source on a log-y axis with nice formatting
45     plt.semilogy(MysterySourceDF['Energies'],MysterySourceDF['count
46                                         ↵ number'],label="Mystery Source")
47     plt.semilogy(BgWithShieldDF['Energies'],BgWithShieldDF['count
48                                         ↵ number2'],label="Background With Shield")
49     plt.xlabel('Gamma Energy [keV]',**AxTitleFont)
50     plt.ylabel('Log-10 of count number',**AxTitleFont)
51     plt.title('Mystery Source With Background Subtracted',**TitleFont)
52     plt.xlim(0,2216)
53     plt.ylim(1)
54     plt.legend()
55     plt.show()
56
57
58 if PlotKnownSources:
59     BgWithShield = "WeekendBgWithShield_003_eh_1"
60     BgWithShieldDF = pd.read_csv(BgWithShield+ ".dat", sep=r"\s+",names = ['channel
61                                         ↵ number','count number'])
62
63     Sodium = 'Sodium22-24HrData_002_eh_1'
64     SodiumDF = pd.read_csv(Sodium+ ".dat", sep=r"\s+",names = ['channel number','count
65                                         ↵ number'])
66
67     Barium = 'Barium133-24HrRun_006_eh_1'
68     BariumDF = pd.read_csv(Barium+ ".dat", sep=r"\s+",names = ['channel number','count
69                                         ↵ number'])
70
71     Cobalt = 'Cobalt60-24HrData_004_eh_1'
72     CobaltDF = pd.read_csv(Cobalt+ ".dat", sep=r"\s+",names = ['channel number','count
73                                         ↵ number'])
74
75     # Normalises the background to the correct timings
76     BgWithShieldDFCobalt = BgWithShieldDF.copy()
77     BgWithShieldDF['count number'] = BgWithShieldDF['count number'] / ( 239068 ) *
78                                         ↵ 86400
79     BgWithShieldDFCobalt['count number'] = BgWithShieldDF['count number'] / ( 239068 )
80                                         ↵ * (86400/12.0)
81
82     # Subtracts the background from the original data
83     SodiumDF['count number'] = SodiumDF['count number'] - BgWithShieldDF['count number']
84     BariumDF['count number'] = BariumDF['count number'] - BgWithShieldDF['count number']
85     CobaltDF['count number'] = CobaltDF['count number'] - BgWithShieldDFCobalt['count
86                                         ↵ number']
87
88     # Fitting parameters
89     a = 2.63977565e-06
90     b = 7.04767648
91     c = -1.91414134e-01
92
93     # Converting channel numbers to energies
94     SodiumDF['Energies'] = np.sqrt( (SodiumDF['channel number'] - c + (b**2.0/ (4 *
95                                         ↵ a)))/a ) - (b/(2*a))

```

```

84 BariumDF['Energies'] = np.sqrt( (BariumDF['channel number'] - c + (b**2.0/ (4 *
85   ↵ a)))/a ) - (b/(2*a))
86 CobaltDF['Energies'] = np.sqrt( (CobaltDF['channel number'] - c + (b**2.0/ (4 *
87   ↵ a)))/a ) - (b/(2*a))
88
89 # Now plotting the different elements on different log-y plots.
90 plt.figure(1)
91 plt.semilogy(SodiumDF['Energies'],SodiumDF['count number'],label="Sodium 22
92   ↵ data",color='yellow')
93 TitleFont = {'size':25, 'color':'black', 'weight':'bold'}
94 AxTitleFont = {'size':22}
95 plt.xlabel('Gamma Energy [keV]',**AxTitleFont)
96 plt.ylabel('Log-10 of count number',**AxTitleFont)
97 plt.title('Sodium 22, no background, energies',**TitleFont)
98 plt.xlim(0)
99 plt.ylim(1)
100 plt.legend()
101
102 plt.figure(2)
103 plt.semilogy(BariumDF['Energies'],BariumDF['count number'],label="Barium 133
104   ↵ data",color='green')
105 TitleFont = {'size':25, 'color':'black', 'weight':'bold'}
106 AxTitleFont = {'size':22}
107 plt.xlabel('Gamma Energy [keV]',**AxTitleFont)
108 plt.ylabel('Log-10 of count number',**AxTitleFont)
109 plt.title('Barium 133, no background, energies',**TitleFont)
110 plt.xlim(0)
111 plt.ylim(1)
112 plt.legend()
113
114 plt.figure(3)
115 plt.semilogy(CobaltDF['Energies'],CobaltDF['count number'],label="Cobalt 60
116   ↵ data",color='blue')
117 TitleFont = {'size':25, 'color':'black', 'weight':'bold'}
118 AxTitleFont = {'size':22}
119 plt.xlabel('Gamma Energy [keV]',**AxTitleFont)
120 plt.ylabel('Log-10 of count number',**AxTitleFont)
121 plt.title('Cobalt 60, no background, energies',**TitleFont)
122 plt.xlim(0)
123 plt.ylim(1)
124 plt.legend()
125 plt.show()

```

## A3.4 RESOLUTION.PY

This Python file was responsible for plotting, and fitting, the energy resolution of each peak as a function of that peaks energy.

```

1 import numpy as np
2 import matplotlib.pyplot as plt
3 import scipy.optimize as sciopt
4 import seaborn as sns
5 import pandas as pd
6 sns.set()
7 import scipy.stats as scistats
8
9 def ResolutionFit(E,a,b):
10     return a/np.sqrt(E) + b
11
12 FitDF = pd.read_csv('Gamma_Peak_Stats_and_Parms.csv', names=['Element', 'Fit type',
13     'Peak number', 'Peak type', 'Energy (keV)', 'Resolution', 'Min', 'Max', 'Mean',
14     'A', 'Sigma', 'Error', 'a', 'b', 'chisq', 'Reduced chisq',
15     'Probchisq'],usecols=list(range(1,18)))
16
17 FitDF['Energy (keV)'] = FitDF['Energy (keV)'].astype(float)
18 FitDF['Resolution'] = FitDF['Resolution'].astype(float).multiply(100)
19 SodiumDF = FitDF[(FitDF['Energy (keV)'] == 511)]
20 FitDF = FitDF[~(FitDF['Energy (keV)'] == 511)]
21
22 DataBa = FitDF[FitDF['Element'] == 'Barium133']
23 DataCo = FitDF[FitDF['Element'] == 'Cobalt60']
24 DataNa = FitDF[FitDF['Element'] == 'Sodium22']
25
26 Fit, Errors = sciopt.curve_fit(ResolutionFit,FitDF['Energy
27     (keV)'],FitDF['Resolution'])
28 print(Fit)
29
30 TitleFont = {'size':'20', 'color':'black', 'weight':'bold'}
31 AxTitleFont = {'size':'18'}
32
33 plt.figure(1)
34 Energies = np.linspace(55,1400,14000)
35 plt.plot(Energies,ResolutionFit(Energies,*Fit),label="Fit")
36 plt.errorbar(DataBa['Energy (keV)'],DataBa['Resolution'],fmt="o",color='red',label="B
37     arium-133",markersize=7,yerr=(7.05/DataBa['Energy
38     (keV)']),elinewidth=3,capsize=5,capthick=3,marker='o')
39 plt.errorbar(DataCo['Energy (keV)'],DataCo['Resolution'],fmt="o",color='blue',label="_
40     Cobalt-60",markersize=7,yerr=(7.05/DataCo['Energy
41     (keV)']),elinewidth=3,capsize=5,capthick=3,marker='o')
42 plt.errorbar(DataNa['Energy (keV)'],DataNa['Resolution'],fmt="o",color='yellow',label_
43     ="Sodium-22",markersize=7,yerr=(7.05/DataNa['Energy
44     (keV)']),elinewidth=3,capsize=5,capthick=3,marker='o')
```

```

38 plt.errorbar(SodiumDF['Energy (keV)'],SodiumDF['Resolution'],fmt="o",color='yellow',l_
  ↵ abel="",markersize=7,yerr=(7.05/SodiumDF['Energy
  ↵ (keV)']),elinewidth=3,capsize=5,capthick=3,marker='o')
39 plt.xlabel('Peak Energy [keV]',**AxTitleFont)
40 plt.ylabel('Energy Resolution [%]',**AxTitleFont)
41 plt.title('Energy Resolution as a Function of Gamma Energy',**TitleFont)
42 plt.legend(fontsize=16,fancybox=True,shadow=False)
43
44 plt.figure(2)
45 plt.errorbar(DataBa['Energy (keV)'],DataBa['Resolution']-ResolutionFit(DataBa['Energy
  ↵ (keV)'],*Fit),fmt="o",color='red',label="Barium-133",markersize=7,yerr=(7.05/Data_
  ↵ Ba['Energy
  ↵ (keV)']),elinewidth=3,capsize=5,capthick=3,marker='o')
46 plt.errorbar(DataCo['Energy (keV)'],DataCo['Resolution']-ResolutionFit(DataCo['Energy
  ↵ (keV)'],*Fit),fmt="o",color='blue',label="Cobalt-60",markersize=7,yerr=(7.05/Data_
  ↵ Co['Energy
  ↵ (keV)']),elinewidth=3,capsize=5,capthick=3,marker='o')
47 plt.errorbar(DataNa['Energy (keV)'],DataNa['Resolution']-ResolutionFit(DataNa['Energy
  ↵ (keV)'],*Fit),fmt="o",color='yellow',label="Sodium-22",markersize=7,yerr=(7.05/Da_
  ↵ taNa['Energy
  ↵ (keV)']),elinewidth=3,capsize=5,capthick=3,marker='o')
48 plt.errorbar(SodiumDF['Energy
  ↵ (keV)'],SodiumDF['Resolution']-ResolutionFit(SodiumDF['Energy (keV)'],*Fit),fmt="_
  ↵ o",color='yellow',label="",markersize=7,yerr=(7.05/SodiumDF['Energy
  ↵ (keV)']),elinewidth=3,capsize=5,capthick=3,marker='o')
49 plt.xlabel('Peak Energy [keV]',**AxTitleFont)
50 plt.ylabel('Energy Resolution Residual [%]',**AxTitleFont)
51 plt.title('Energy Resolution Residuals',**TitleFont)
52 plt.legend(fontsize=16,fancybox=True,shadow=False)
53 plt.show()
54
55 def ChiSqFunc(Measured,Fitted,Errors,Params):
56     ChiSquared = 0
57     for i in range(len(Measured)):
58         ChiSquared += ((Measured[i] - Fitted[i])**2.0) / ((Errors[i])**2.0)
59     ReducedChiSq = ChiSquared/(len(Measured)-len(Params))
60     ProbChiSq = (1.0 - scistats.chi2.cdf(ChiSquared,len(Measured)-len(Params)) ) *
  ↵ 100.0
61     return ChiSquared, ReducedChiSq, ProbChiSq
62
63
64 MeasuredRes = list(DataBa['Resolution'])+list(DataCo['Resolution']) +
  ↵ list(DataNa['Resolution'])
65 FittedRed = list(ResolutionFit(DataBa['Energy (keV)'],*Fit)) +
  ↵ list(ResolutionFit(DataCo['Energy (keV)'],*Fit)) +
  ↵ list(ResolutionFit(DataNa['Energy (keV)'],*Fit))
66 ErrorRes = list(7.05/DataBa['Energy (keV)']) + list(7.05/DataCo['Energy (keV)']) +
  ↵ list(7.05/DataBa['Energy (keV)'])
67
68 ChiStats = ChiSqFunc(MeasuredRes, FittedRed, ErrorRes,Fit)
69 print(ChiStats)

```