



# Gamma-ray Spectrometry

## The basic equation for gamma-ray spectrometry

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# Obligatory slide

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# Basic Equation

$$C = AP_{\gamma}t\varepsilon$$

$$A = \frac{C}{t\varepsilon P_{\gamma}}$$

Peak Count

Gamma-ray emission probability

FEP (Full Energy Peak) efficiency

Measurement time (live time)

# Activity calculations & efficiency

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$$A = \frac{C_{TOT} - C_{Peak}^{Bkg} - C_{Continuum}}{\epsilon_{REF}^{Exp} \frac{\epsilon_{Sample}^{MC}}{\epsilon_{REF}^{MC}} P_{\gamma}} e^{\lambda t_d} \frac{\lambda}{(1 - e^{-\lambda t_m})} K_1 K_2 K_3$$

Reference  
sample similar  
to the measured  
one

Correction factor from  
e.g. EGS4 MonteCarlo  
code

**Uncertainties on efficiency  
down to ~ 1%**

$K_1$  = summing correction

$K_2$  = Branching correction

$K_3$  = Equilibrium correction

$t_d$  = decay time (to a reference date)

$t_m$  = measurement live time

Combined activities from  
several gamma-rays and from  
several daughters to activity  
for one radionuclide



# Basic Equation – Time corrections

Decay during measurement and decay from reference date

$$\frac{1}{t_m} \longrightarrow \frac{\lambda}{(1 - e^{-\lambda t_m})}$$

Decay from reference date  $e^{\lambda t_d}$

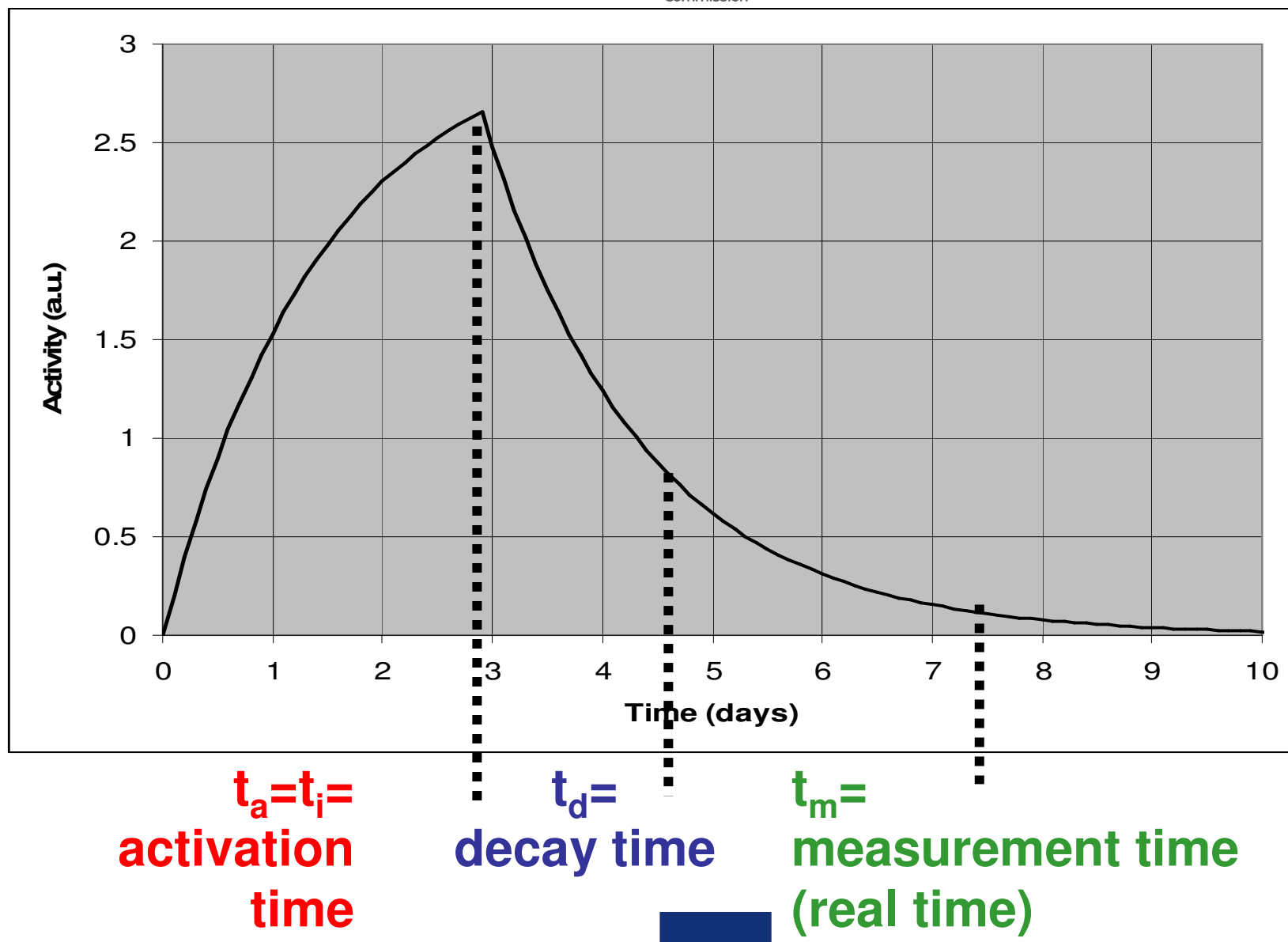


# Basic Equation – Time corrections

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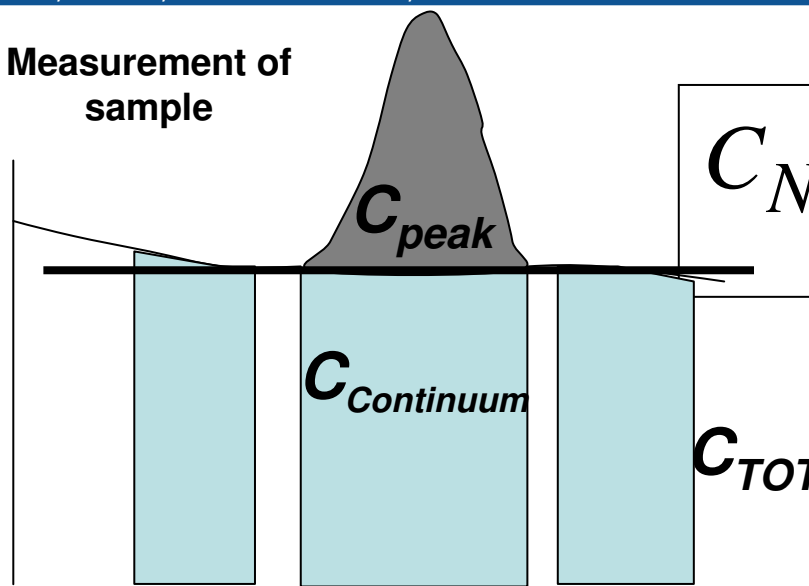


# Basic Equation – Peak Counts

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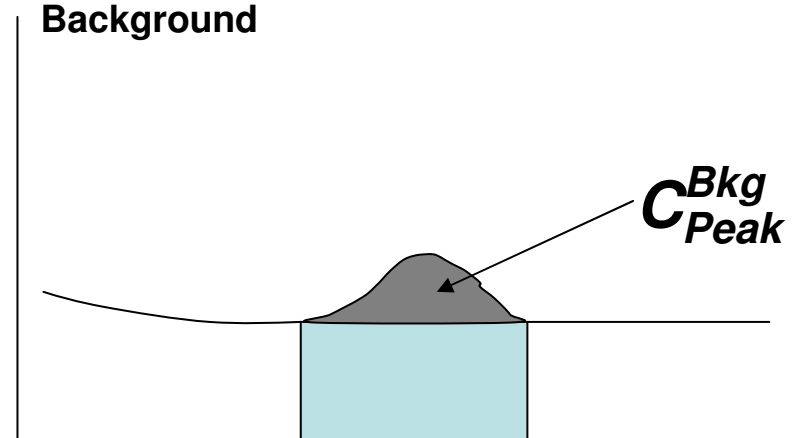
Measurement of  
sample



$$C_{NET} = C_{TOT} - C_{Peak}^{Bkg} - C_{Continuum}$$

$$C_{TOT} = C_{Continuum} + C_{peak}$$

Measurement of  
Background



# Basic Equation – Peak Counts – peak fitting

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## Different algorithms for different data analysis software

- **Genie-2000 (Canberra)**
  - **InterWinner (Eurisys)**
  - **Gamma-Vision (Ortec)**
  - **Gamma-W (Dr. Westmaier)**
  - **Fitzpeaks**
  - **HYPERMET**
- + many other more or less home-made ones**





# Basic Equation – Peak Counts – peak fitting

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## A Golden Rule

- **Don't trust an automatic fit. Use visual inspection whenever possible!**

### **Common problems:**

- **unexpected Interfering peak(s),**
- **Errors in nuclear data library,**
- **erroneous settings used,**
- **bad peak shape due to unstable electronics during measurements, etc.**

**The best QC is to visually inspect spectra!**



# Basic Equation – Peak Counts – peak fitting

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- **Start with a good energy calibration and a good calibration of the FWHM.....and good quality data (Gaussian peaks)**
- **Ideally a full-energy peak can be represented by a Gaussian function**
- **In practice there is tailing on the low-E side of the peak (also at hi-E side at e.g. high count rates)**
- **Note different shapes for X-rays peaks and annihilation peak**



# Basic Equation – Peak Counts – peak fitting

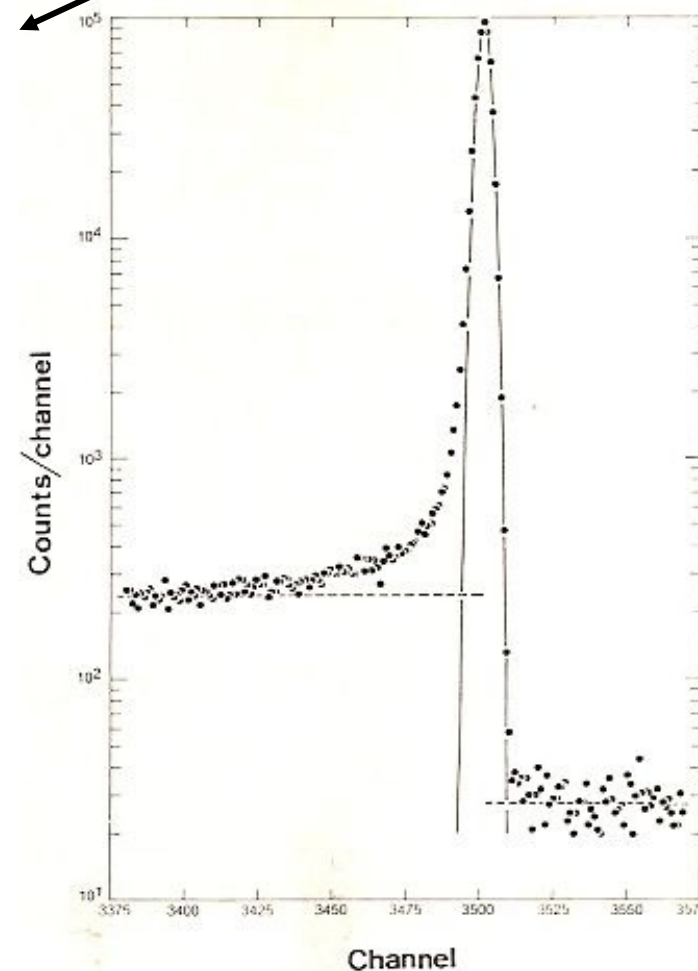
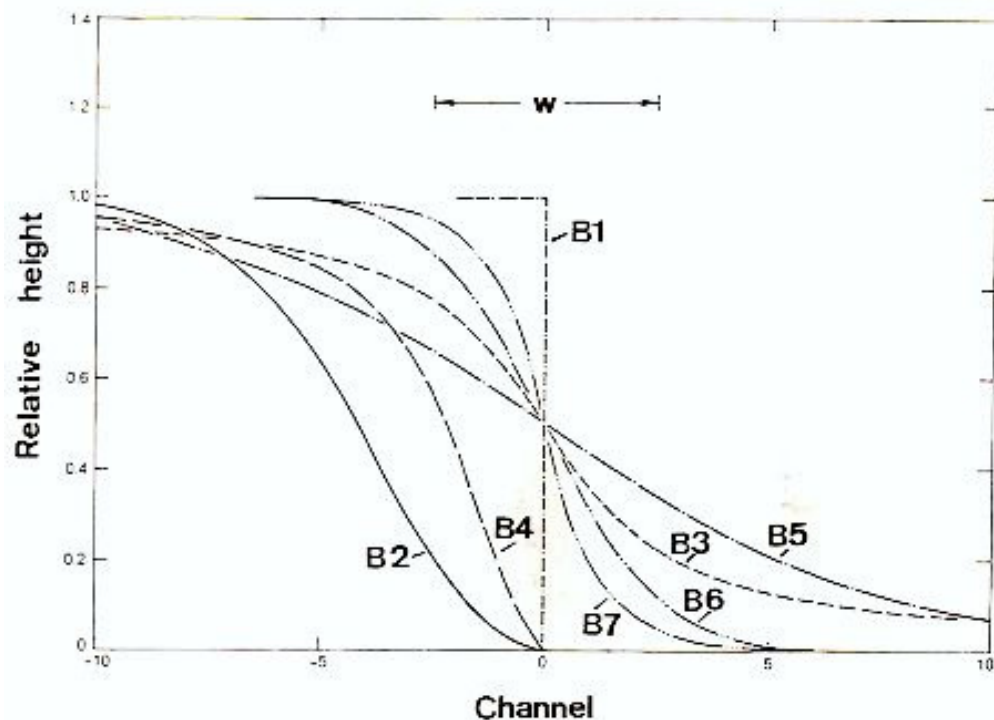
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## Most common fits include

1. A Gaussian (main feature)
2. A linear background with step-function
3. A linear low-E tail

Mainly these  
varies between  
software



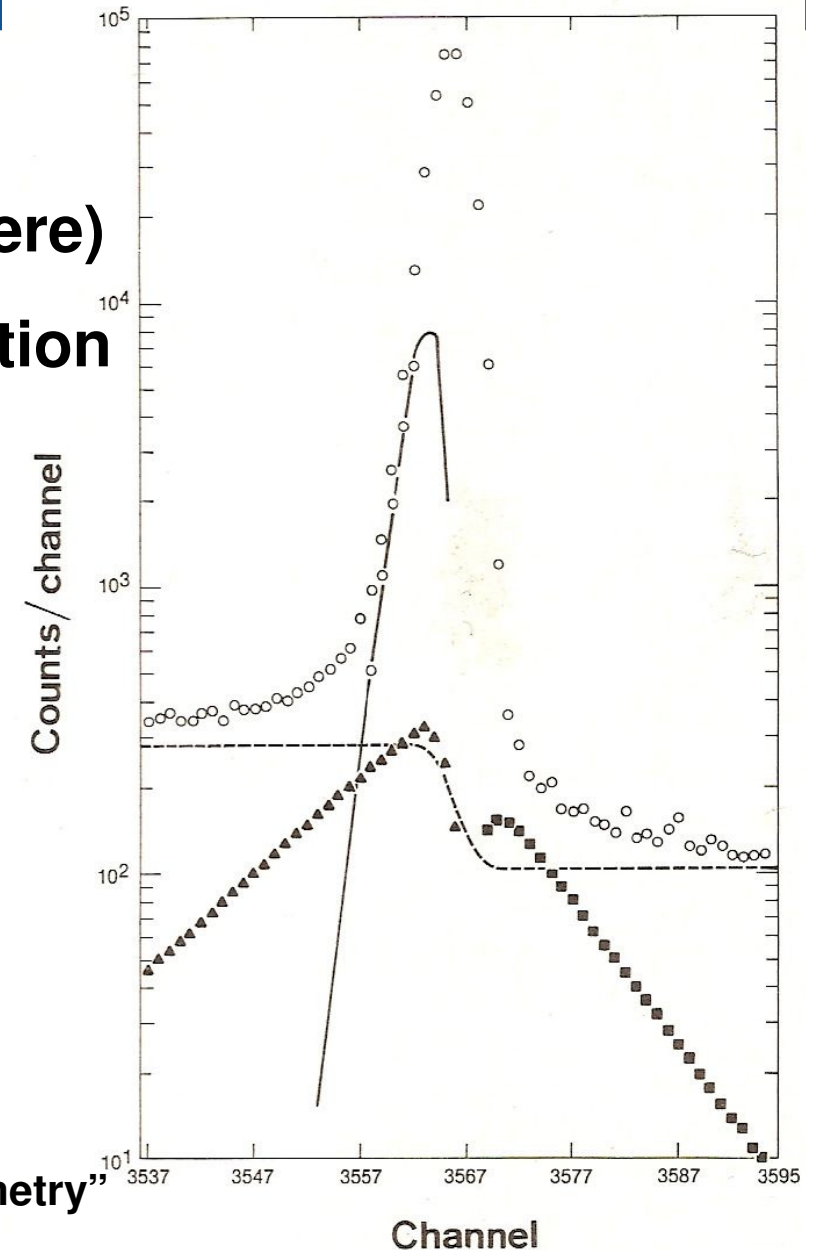
# Basic Equation – Peak Counts – peak fitting

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**More complex fits include**

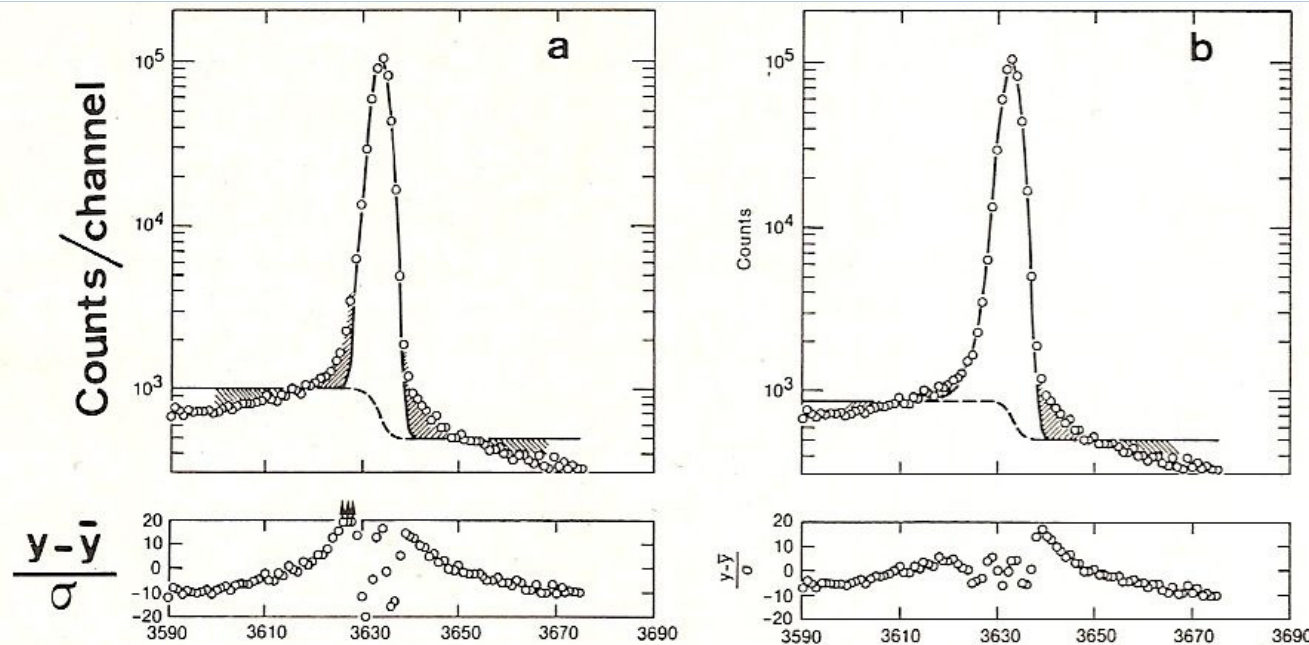
- 1. Gaussian (main feature-not shown here)**
- 2. Constant background with step-function**
- 3. Long term linear low-E tail**
- 4. Short term low-E-tail**
- 5. Hi-E tail**



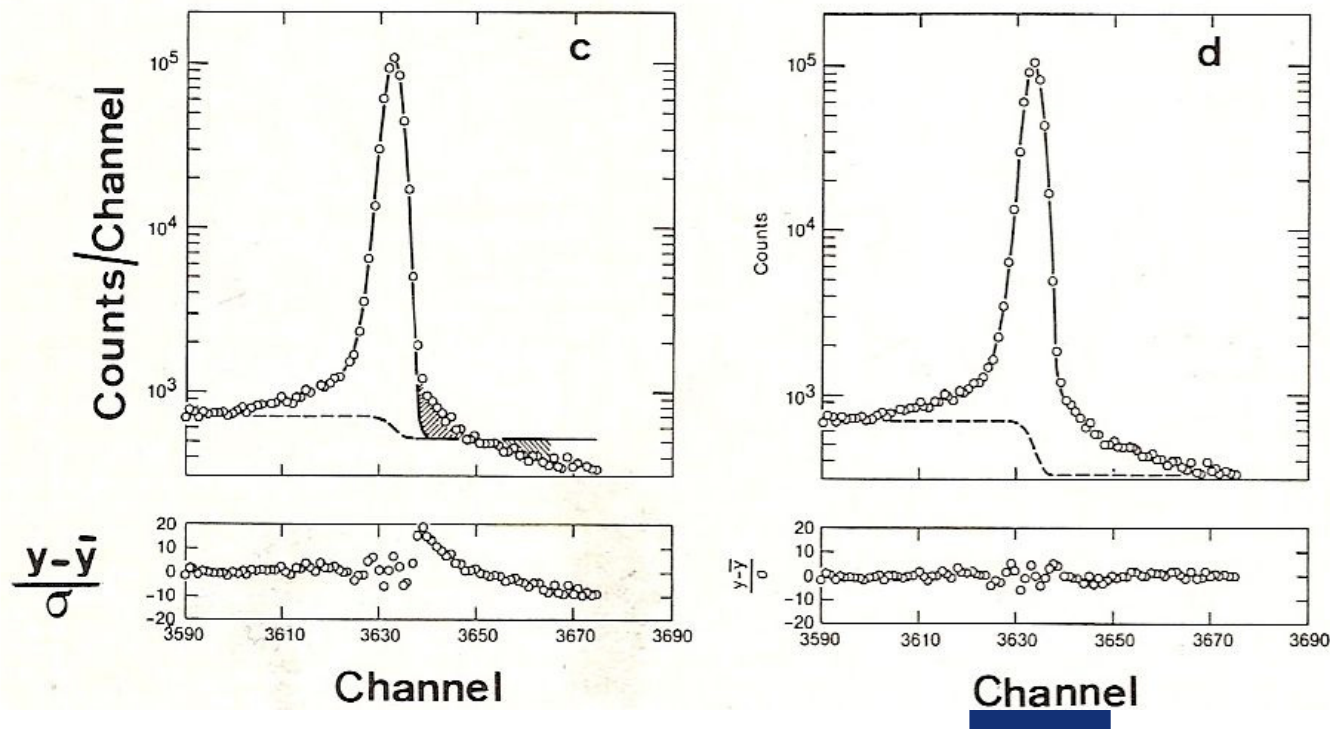
From: Debertin and Helmer, "X-ray and gamma-ray spectrometry"



**Less tailing with  
“new generation”  
HPGe-detectors!**

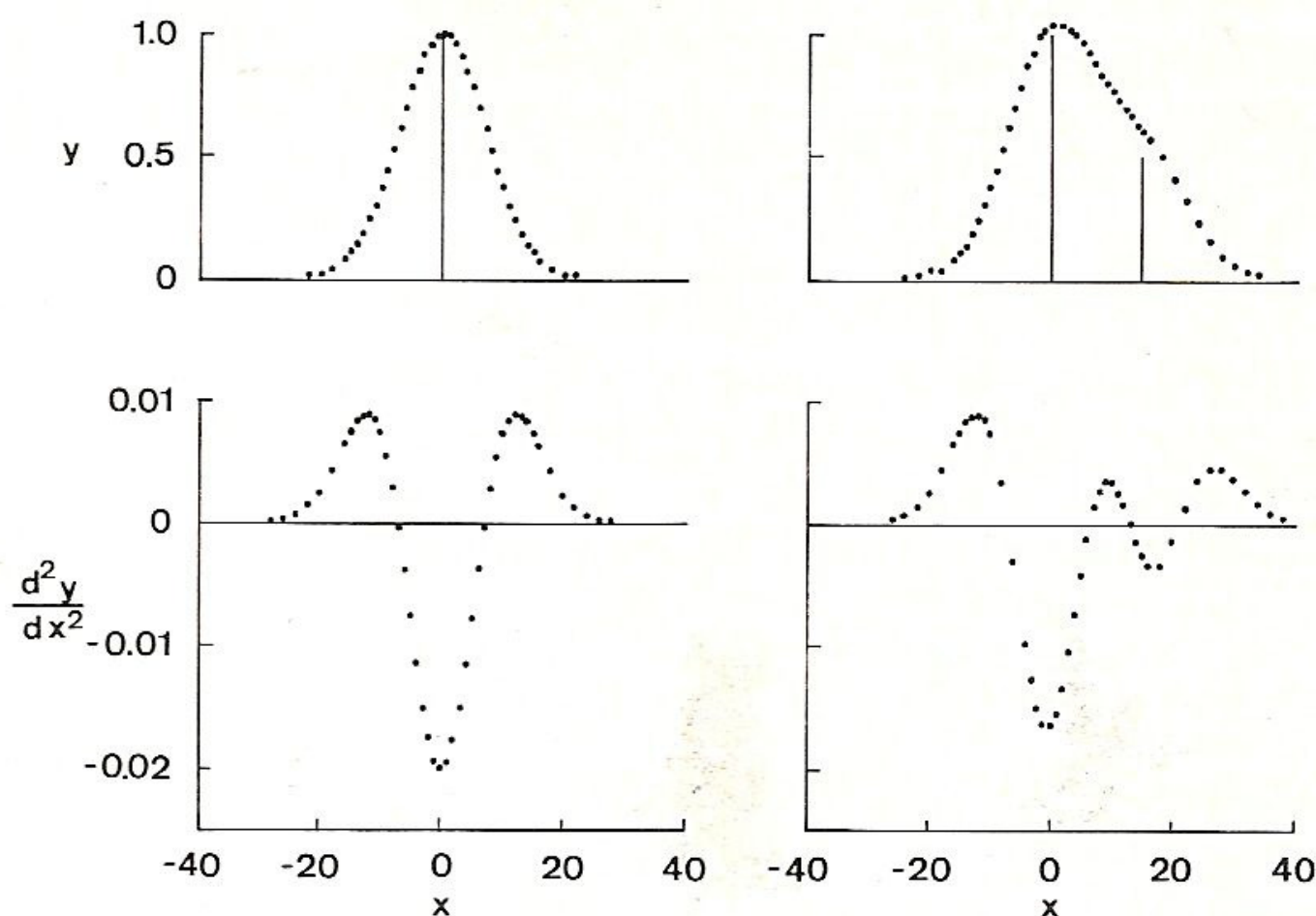


**Important to have  
“nice” well  
separated peaks to  
determine  
parameters for  
tailing**





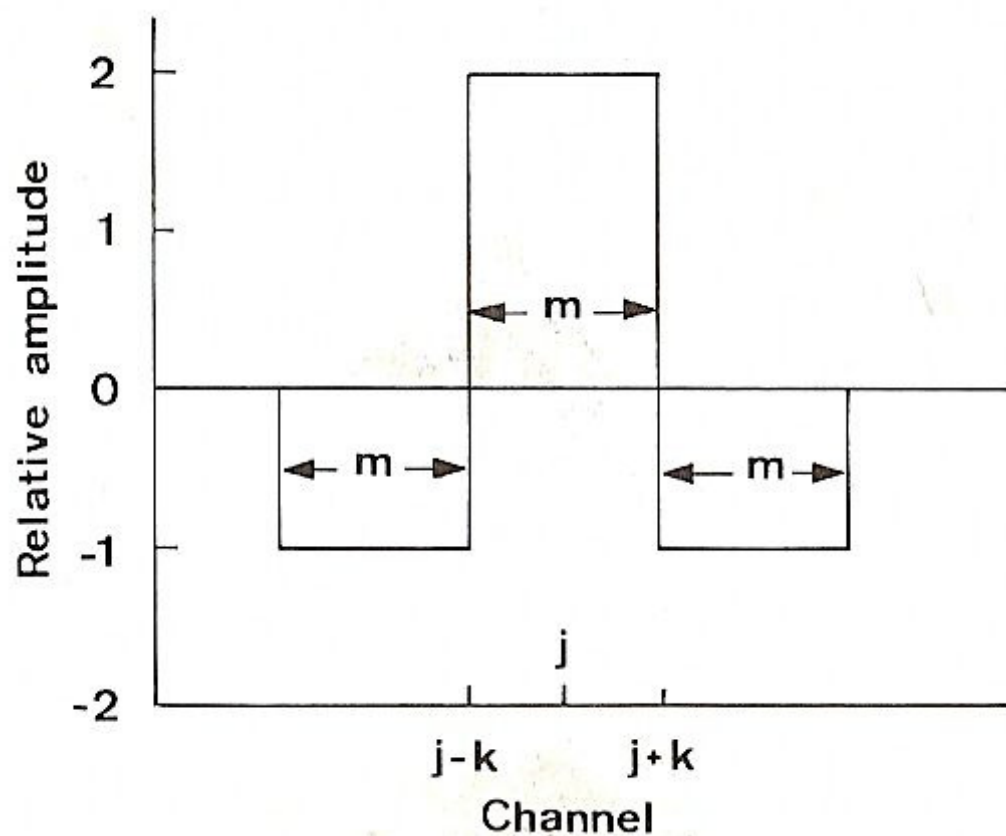
# Automatic Peak Search – 2<sup>nd</sup> derivative



- Smoothed data
- apply statistical selection criteria
- => can distinguish single peak/doublets/Compt on edges etc.



# Automatic Peak Search – Filter



- The simplest filter shown here
- Calculate C

$$C_j = \sum_{i=j-k}^{j+k} 2y_i - \sum_{i=j-k-(2k+1)}^{j-k-1} y_i - \sum_{i=j+k+1}^{j+k+(2k+1)} y_i$$

- $C > \text{given value} \Rightarrow$  peak detected



# Basic Equation – Decay Data

Referring to presentation by Vidmar (Bé)!

$P_\gamma$  (gamma-ray emission probability) and

$\lambda$  (Decay constant) can vary significantly between  
different sources of data

Pay attention to equilibrium between mother-daughter,  
which decay constant (half-life) to use. Not automatically  
correct in commercial software!

And library data may not be up to date and not from  
DDEP.







## Basic Equation – Efficiency

*Measure FEP efficiency in a calibration source with several gamma-rays*

*Fit an empirical (or semiempirical) function to the data*

*If the "real sample" is different from the calibration source in shape, mass or composition – calculate efficiency correction using an efficiency transfer code*

**And... Note that time corrections usually needed !!!**

$$\varepsilon = \frac{C}{tAP_{\gamma}}$$

**BUT: Beware of coincidence summing corrections!!!**

# Basic Equation – Efficiency

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**Lots of articles on “suitable” mathematical expressions,**  
**some general others very specific for a certain geometry and**  
**detector**

$\alpha_1$ : adjustable  
parameter

$\alpha_2$  and  $\alpha_3$ , Depends  
on crystal size

Loss of Compton  
scattered photons

**Semi-empirical  
Early attempt  
(1966):**

$$\varepsilon(E) = a_1 \left[ \tau(E) + \sigma(E) \alpha_2 e^{-\alpha_3 E} \right]$$

Photoelectric  
cross-section of  
Ge

Compton cross-  
section of Ge

Only useful for small detectors and energies  
from 500 keV to 1500 keV

# Basic Equation – Efficiency

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**Empirical attempts with polynomial liner or non-linear functions dominate today, e.g.:**

$$\varepsilon = \frac{1}{E} \sum_{i=1}^8 a_i (\log E / E_0)^{i-1}$$

$$\log \varepsilon = \sum_{i=1}^n a_i (\log E / E_0)^{i-1}$$

**Used e.g. in Genie 2000**

**But needs a second polynomial to handle the low energy part (below ~200 keV)**

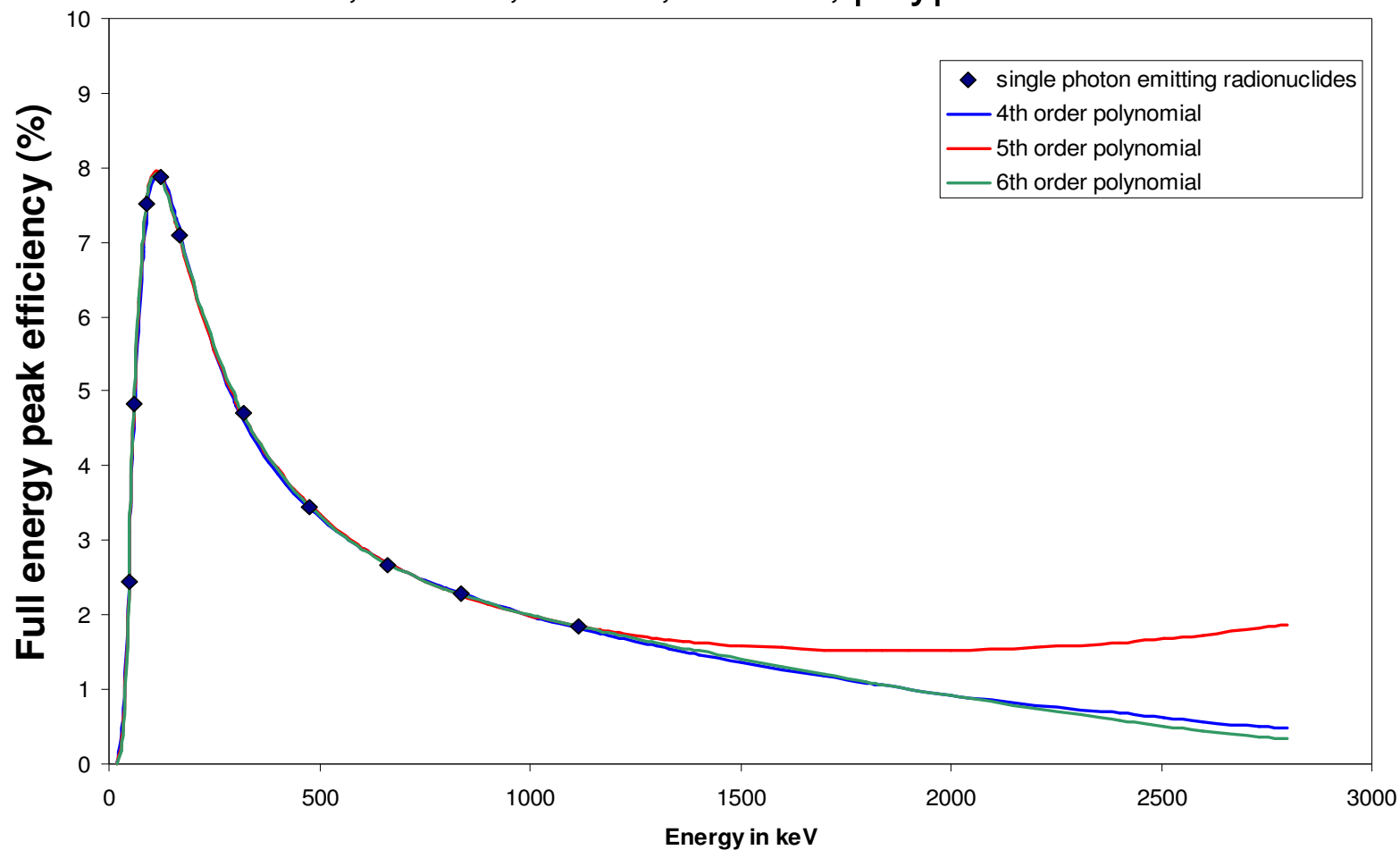


# Basic Equation – Efficiency

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Full energy peak efficiency for Pb-210, Am-241, Cd-109, Co-57, Be-7, Cs-137, Mn-54, Zn-65; 100 ml; p-type detector



# Basic Equation – Efficiency Transfer

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**Several codes for calculating efficiency transfer exists:  
EFFTRAN, GESPECOR, DETEFF, ETNA,.....**

**Efficiency corrections can be calculated using general  
purpose Monte Carlo codes like: GEANT, PENELOPE, EGS4  
and MCNP**

$$\mathcal{E} = \mathcal{E}_{REF}^{Exp} \frac{\mathcal{E}_{Sample}^{MC}}{\mathcal{E}_{REF}^{MC}}$$





## Absolute $\varepsilon$ -values from Monte Carlo codes

**It is possible to use the absolute efficiency calculated using a Monte Carlo code but great care should be taken since:**

- **The computer model of the sample and detector can never describe reality in detail**
- **It is easy to make a small mistake**
- **The model might well describe the reference sample used for setting up the model – but might fail when the geometry changes.**
- **It is difficult to assess the uncertainty of a Monte Carlo calculation**





# Realisation of Calculations

## 1. Making full use of commercial software

- Most suited for routine analysis (analysis of many samples of the same type (matrix, mass, radionuclides))
- Drawbacks: User has not full control of all calculations.

## 2. Home-made specially designed software

- Takes long to implement

## 3. Making tailor-made calculation sheets in general purpose codes like MS Excel

- Can get “messy” unless care is taken to “program cleanly”



# Spectrum analysis and calculations in commercial software

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**Const. T,  
etc.**

**Stable  
electricity**

**Good quality data**

**E-cal.  
FWHM cal.**

**Peak locate**

**Peak fitting**

**Visual inspection  
of fitting**

**Bkg subtr. &  
Nuclide id.**

**Efficiency**

**Summing  
correction**

**Parent daughter  
correction**

**Reporting**

**Many steps => easy  
To get one wrong**

**Take care when using  
Automatic systems**

**Non-automatic =>  
Problems with  
integrity of data**

**Uncertainty? Often  
only counting  
statistics**





$$A = \frac{C_{TOT} - C_{Peak}^{Bkg} - C_{Continuum}}{\epsilon_{REF}^{Exp} \frac{\epsilon_{Sample}^{MC}}{\epsilon_{REF}^{MC}} P_{\gamma}} e^{\lambda t_d} \frac{\lambda}{(1 - e^{-\lambda t_m})} K_1 K_2 K_3$$

**$K_1$  = summing correction**

**$K_2$  = Branching correction**

**$K_3$  = Equilibrium correction**

**Combine activity from  
several gamma-rays to  
activity for one  
radionuclide**

**Combine activity from several  
daughters to one parent**

**Ex.:  $^{226}\text{Ra}$  from  $^{214}\text{Bi}$  and  $^{214}\text{Pb}$**