

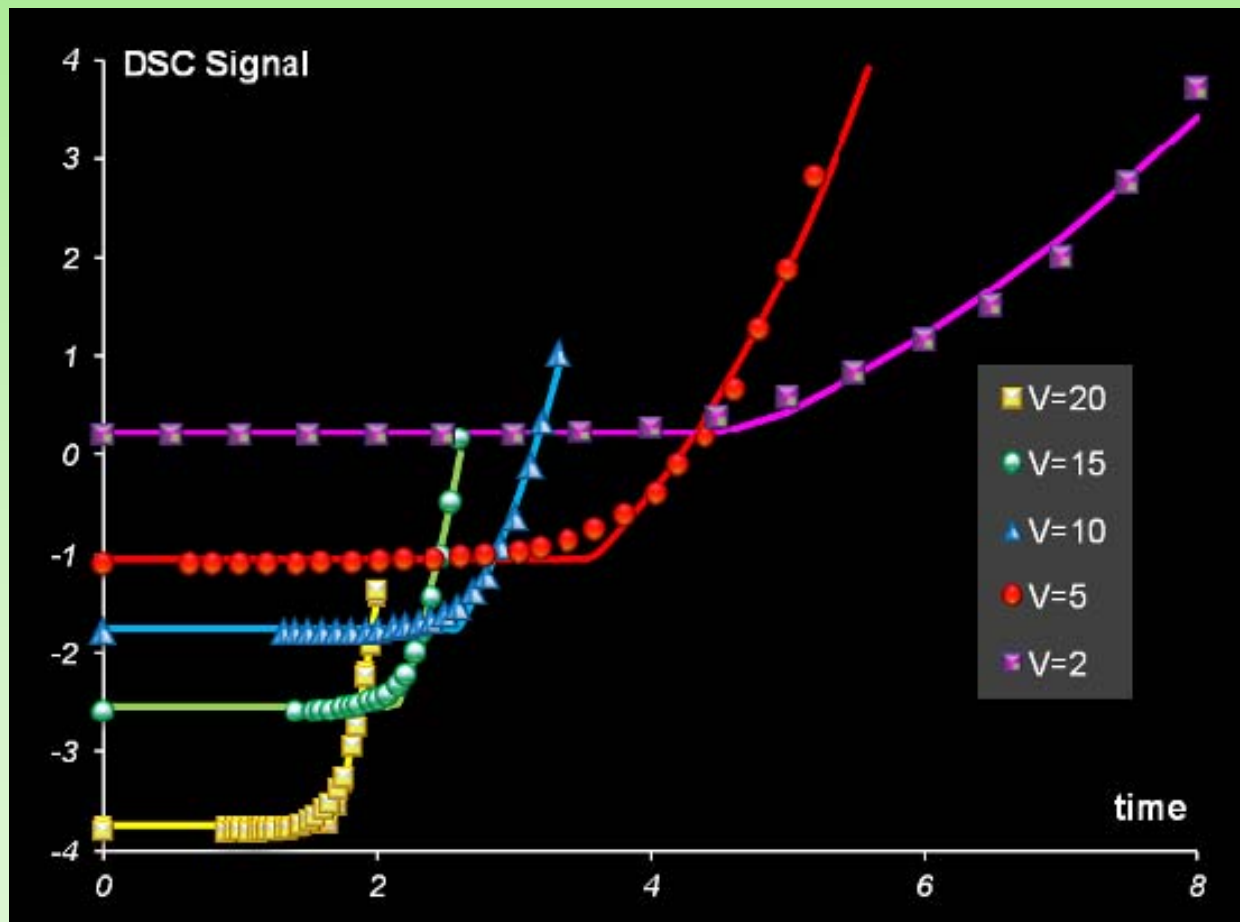
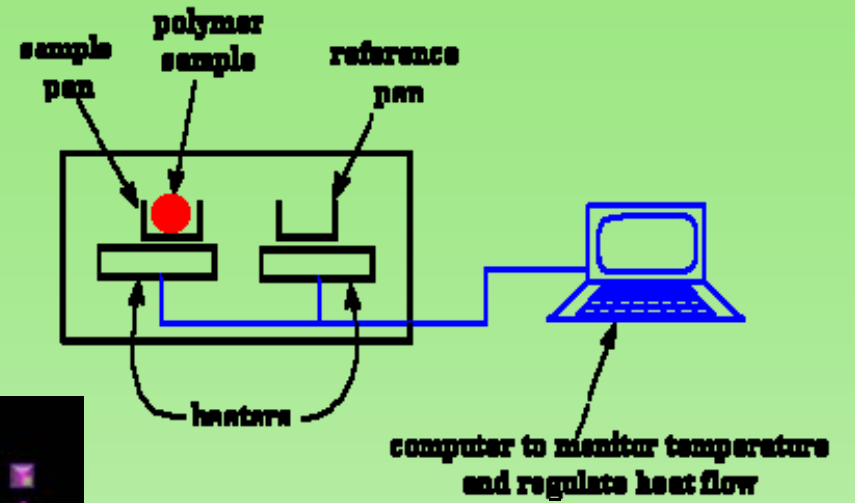
Successive Bayesian Estimation as a tool for chemometric modelling of kinetic data

Alexey Pomerantsev

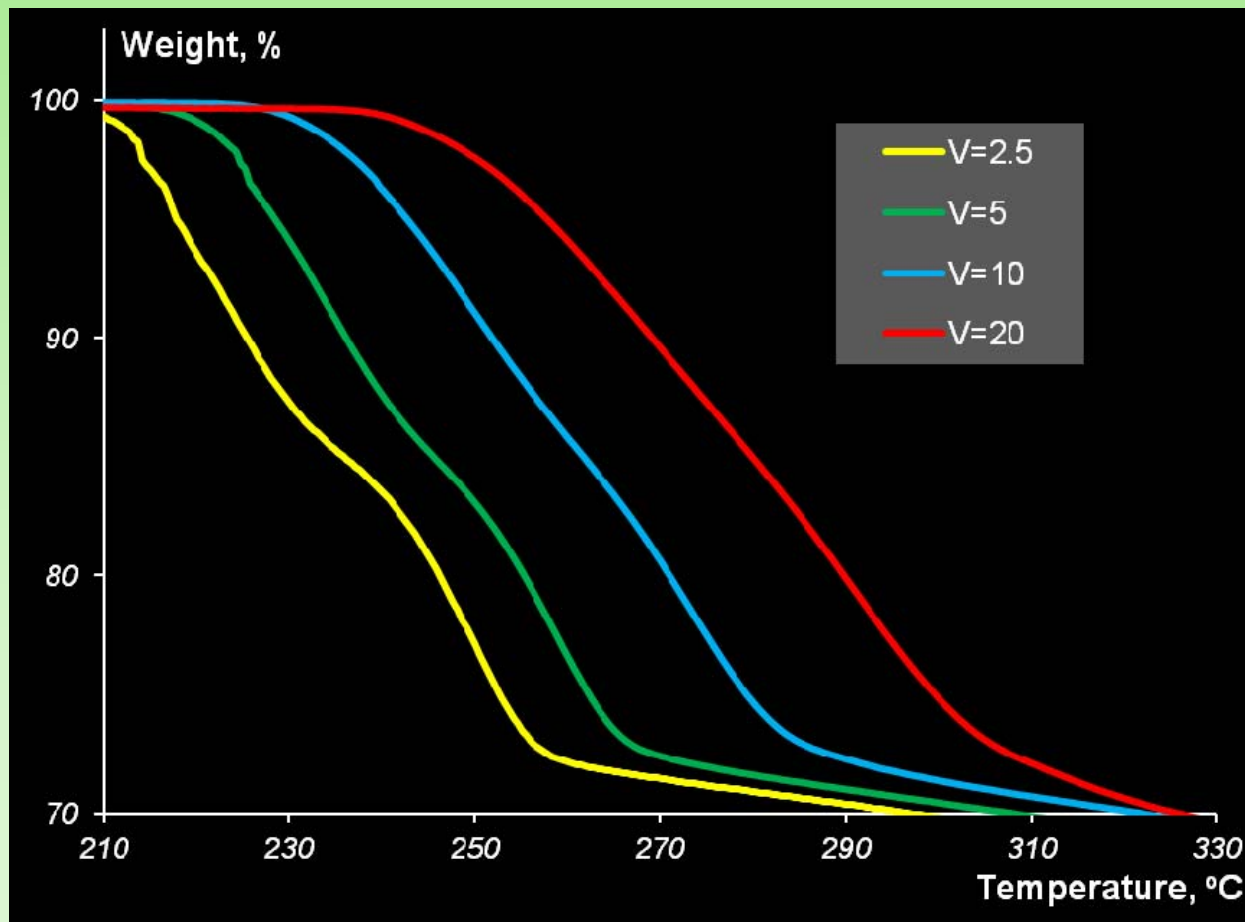
Semenov Institute of Chemical Physics

Russian Chemometrics Society

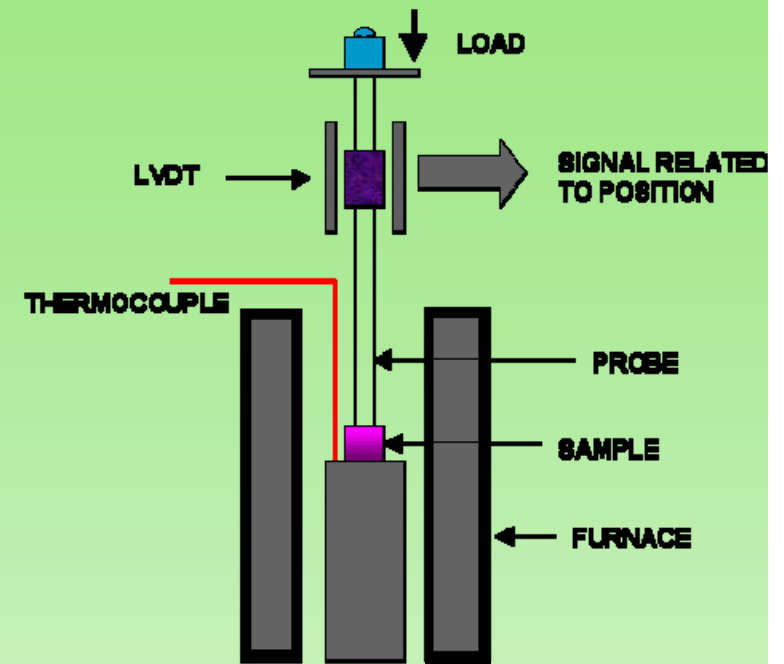
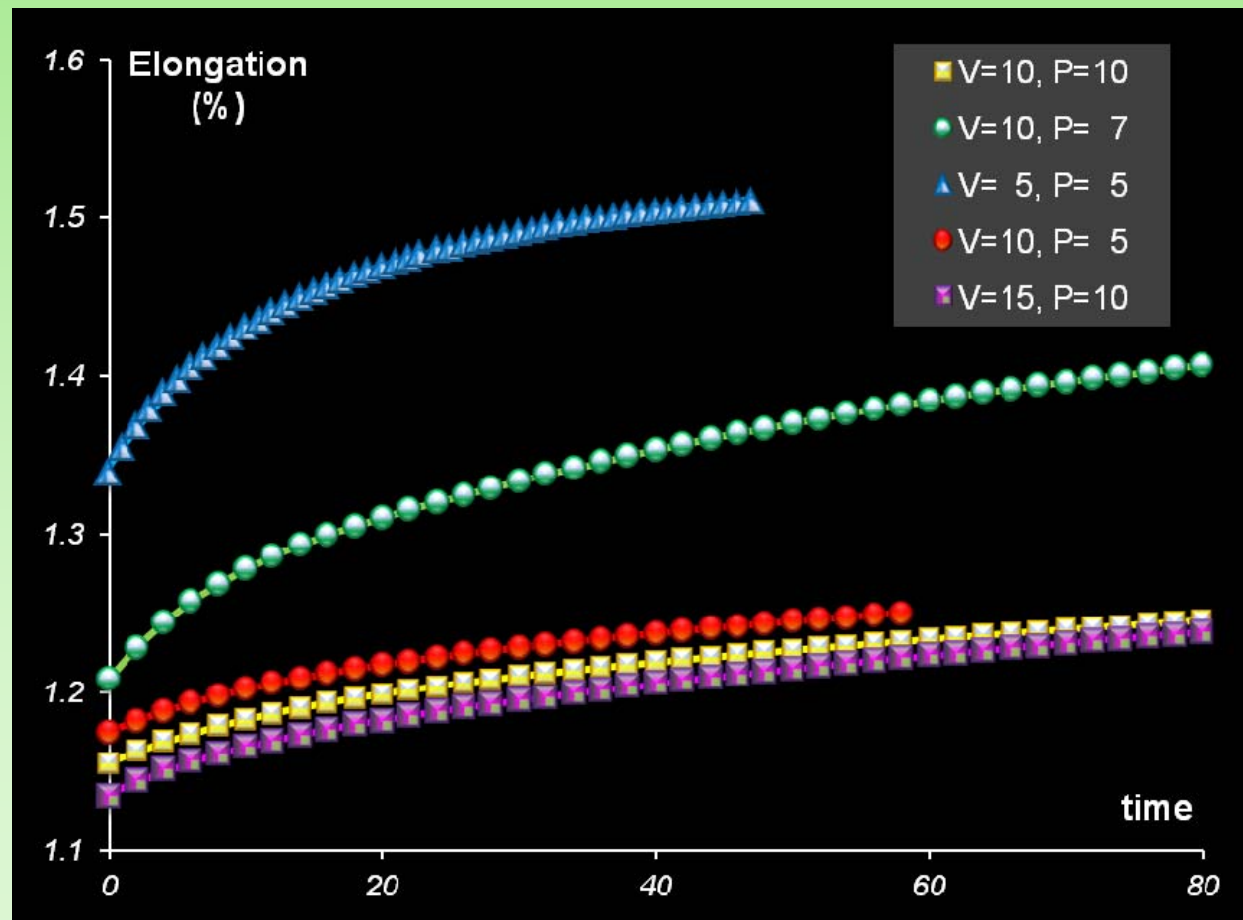
DSC data



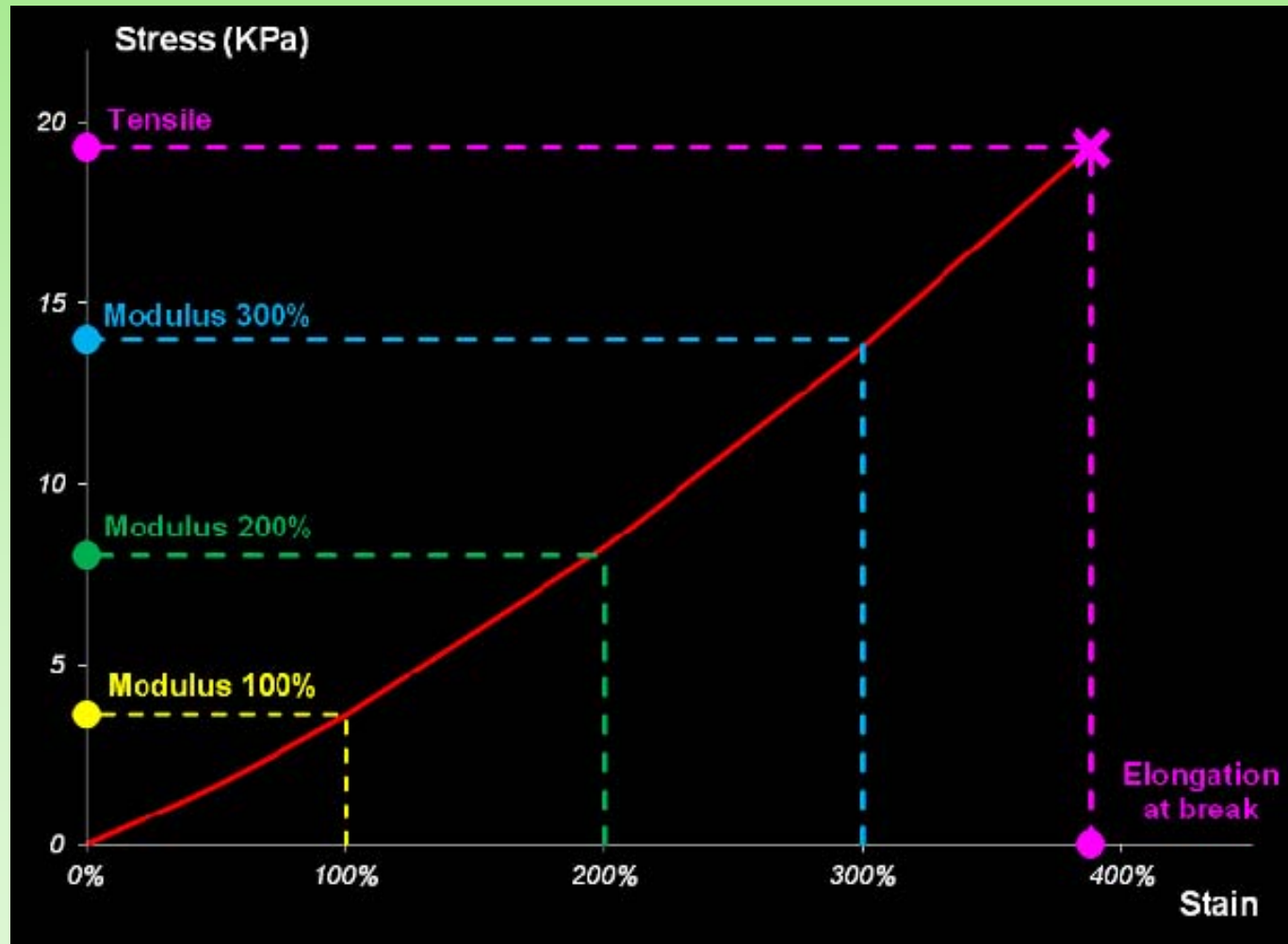
TGA Data



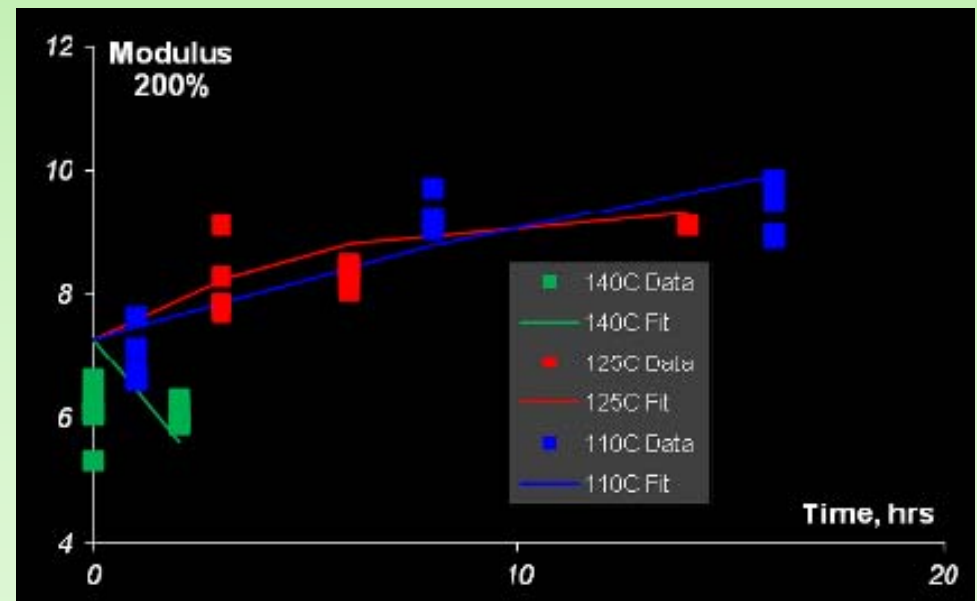
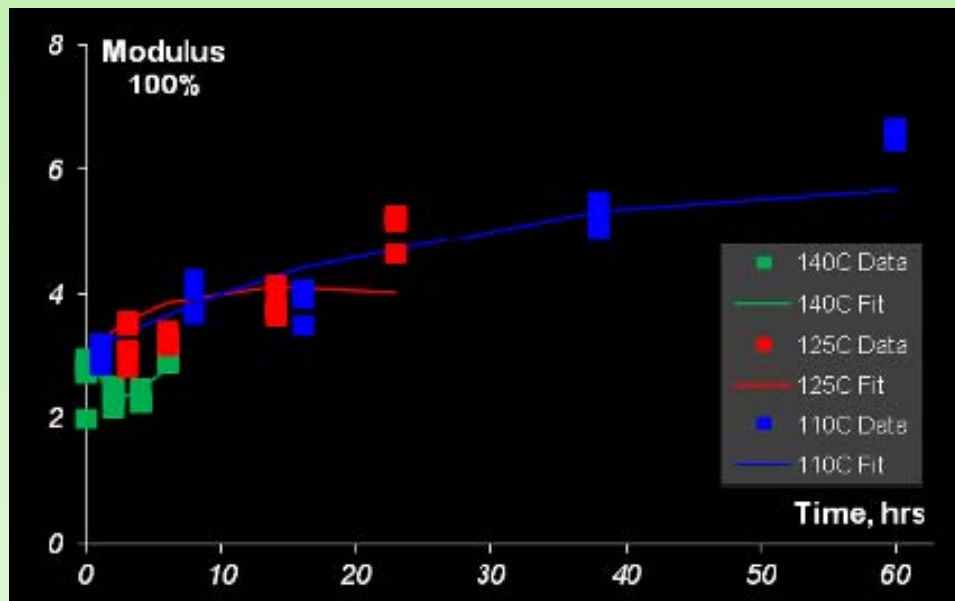
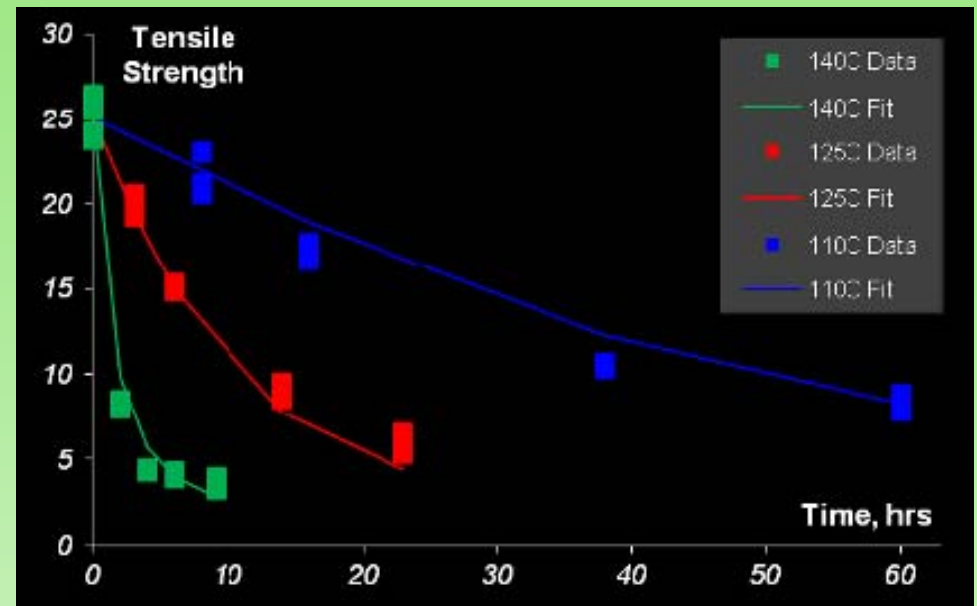
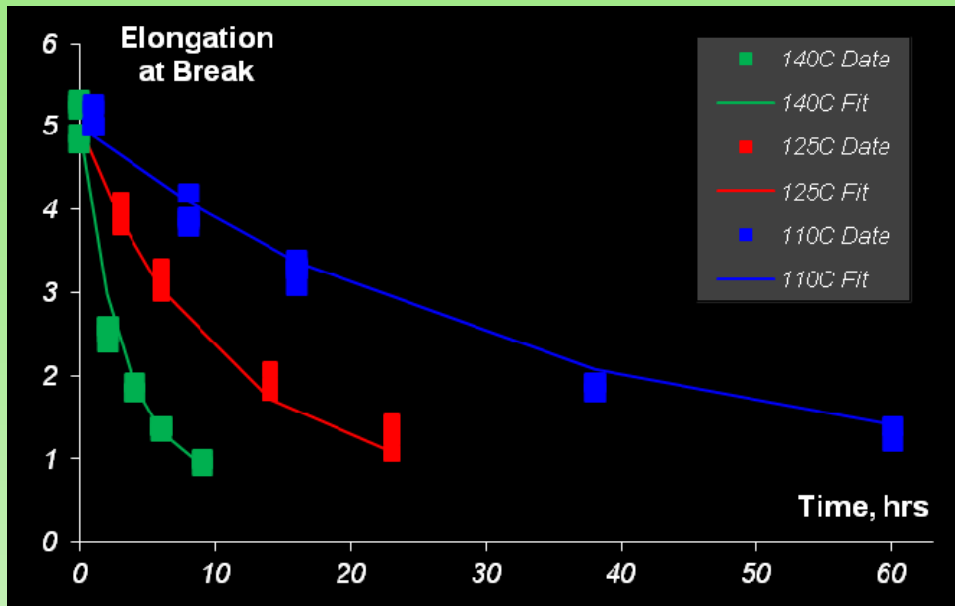
TMA data



Stress-strain experiment

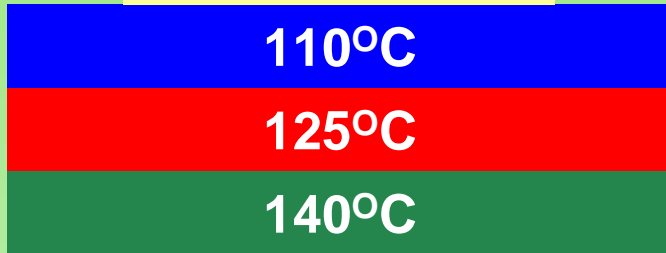


Stress-strain data

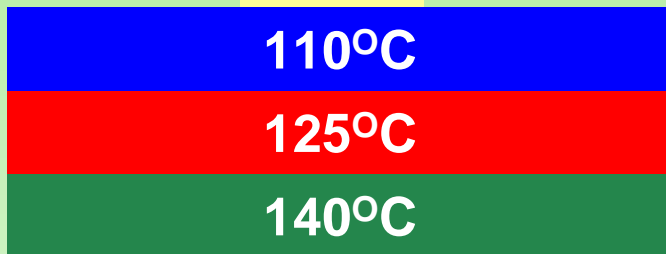


Data Structure

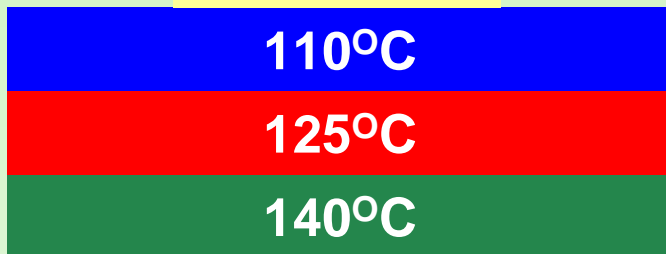
Elongation @ Break



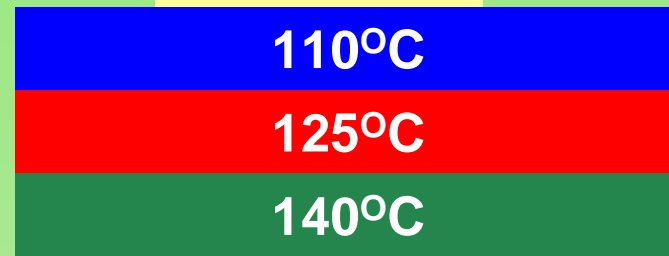
Tensile



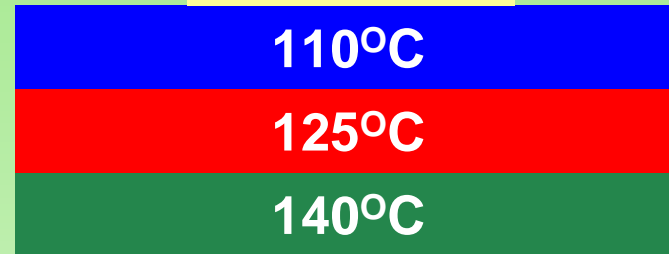
Modulus 100%



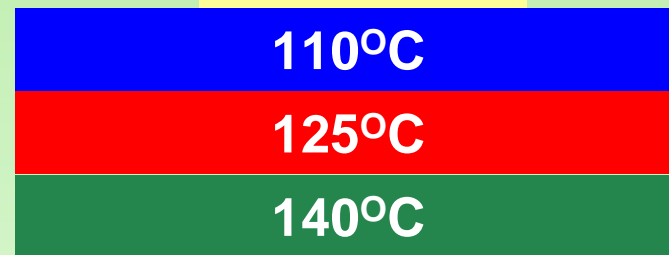
Modulus 200%



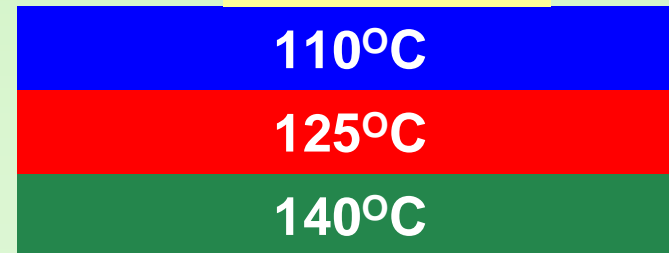
Modulus 300%



Modulus 400%



Modulus 500%



Models structure

- 3×7 data blocks = 21 models

- each model depends on m parameters

$$f(t, T | a_1, a_2, \dots, a_m) \quad m \approx 7$$

- $3 \times 7 \times m \approx 150$ parameters to estimate

- some are common, e.g. Arrhenius' parameters:

$$a = k \exp(E/R/T)$$

- finally:

7 models, 8 partial and 4 common parameters

OLS & SBE for two blocks

OLS

$$\mathbf{y} = \mathbf{y}_1 \otimes \mathbf{y}_2$$

$$\mathbf{f}(\mathbf{a}) = \mathbf{f}_1(\mathbf{a}) \otimes \mathbf{f}_2(\mathbf{a})$$

$$S(\mathbf{a}) = \|\mathbf{y} - \mathbf{f}(\mathbf{a})\|^2 = S_1(\mathbf{a}) + S_2(\mathbf{a}) = \|\mathbf{y}_1 - \mathbf{f}_1(\mathbf{a})\|^2 + \|\mathbf{y}_2 - \mathbf{f}_2(\mathbf{a})\|^2$$

$$\hat{\mathbf{a}}_{\text{OLS}} = \arg \min_{\mathbf{a}} S(\mathbf{a})$$

SBE

$$\hat{\mathbf{a}}_1 = \arg \min_{\mathbf{a}} S_1(\mathbf{a}) \quad S_1(\mathbf{a}) \approx S_1(\hat{\mathbf{a}}_1) + (\mathbf{a} - \hat{\mathbf{a}}_1)^t \mathbf{A}(\mathbf{a} - \hat{\mathbf{a}}_1) = B(\mathbf{a})$$

$$S(\mathbf{a}) \approx Q(\mathbf{a}) = S_2(\mathbf{a}) + B(\mathbf{a})$$

$$\hat{\mathbf{a}}_{\text{SBE}} = \arg \min_{\mathbf{a}} Q(\mathbf{a})$$

Likelihood approach

$$L(\mathbf{a}, \sigma^2) = \text{Const} \sigma^{-(N_1+N_2)} \exp\left(-\frac{S_1(\mathbf{a}) + S_2(\mathbf{a})}{2\sigma^2}\right)$$

$$S_i(\mathbf{a}) = \|\mathbf{y}_i - \mathbf{f}_i(\mathbf{a})\|^2, \quad i = 1, 2$$

$$L(\mathbf{a}, \sigma^2) = \text{Const} \sigma^{-(N_1+N_2-2)} \exp\left(-\frac{B_1(\mathbf{a}) + S_2(\mathbf{a})}{2\sigma^2}\right)$$

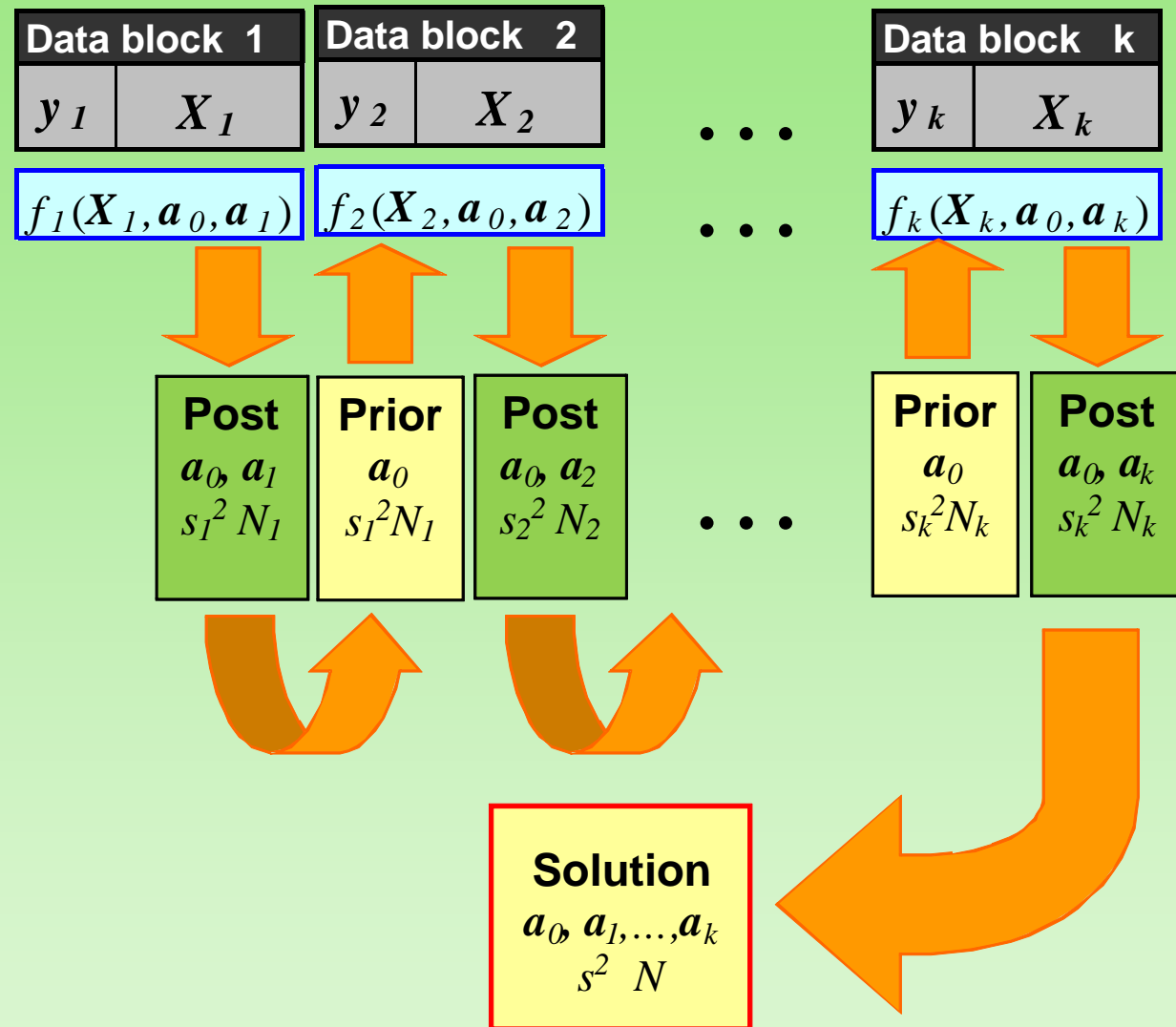
$$B_1(\mathbf{a}) = N_1 s_1^2 + (\mathbf{a} - \hat{\mathbf{a}}_1)^t \mathbf{A} (\mathbf{a} - \hat{\mathbf{a}}_1)$$

$$L(\mathbf{a}, \sigma^2) = \text{Const} \sigma^{-(N_1-2)} \exp\left(-\frac{B_1(\mathbf{a})}{2\sigma^2}\right) \times \sigma^{-N_2} \exp\left(-\frac{S_2(\mathbf{a})}{2\sigma^2}\right) =$$

$$= L_{\text{Prior}}(\mathbf{a}, \sigma^2) \times L_{\text{Data}}(\mathbf{a}, \sigma^2)$$

SBE procedure

- 1) Process first block alone
- 2) Make posterior information
- 3) Convert it to prior information
- 4) Use it for the next block
- 5) Repeat this for all blocks
- 6) Last result is the solution



Posterior & Prior Information

Block 1. Posterior Information

$$B_1(\mathbf{a}) = N_1 s_1^2 + (\mathbf{a} - \hat{\mathbf{a}}_1)^t \mathbf{A}(\mathbf{a} - \hat{\mathbf{a}}_1)$$

Rebuilding (common & partial parameters)

$$\mathbf{A} = \begin{bmatrix} \mathbf{A}_{00} & \mathbf{A}_{01} \\ \mathbf{A}_{01}^t & \mathbf{A}_{11} \end{bmatrix} \quad \mathbf{H} = \begin{bmatrix} \mathbf{A}_{00} - \mathbf{A}_{01} \mathbf{A}_{11}^{-1} \mathbf{A}_{01}^t & \mathbf{0} \\ \mathbf{0} & \mathbf{0} \end{bmatrix}$$

$$\hat{\mathbf{a}}_1 = \begin{bmatrix} \hat{\mathbf{a}}_0 \\ \hat{\mathbf{a}}_{11} \end{bmatrix} \quad \mathbf{b}_1 = \begin{bmatrix} \hat{\mathbf{a}}_0 \\ \mathbf{0} \end{bmatrix}$$

Block 2. Prior Information

$$B_2(\mathbf{a}) = N_1 s_1^2 + (\mathbf{a} - \mathbf{b}_1)^t \mathbf{H}(\mathbf{a} - \mathbf{b}_1)$$

SBE Main Theorem

Different order of blocks processing

1 → 2

$$\hat{\mathbf{a}}_{\text{SBE}}^{12} = \arg \min_{\mathbf{a}} [S_2(\mathbf{a}) + B_1(\mathbf{a})]$$

2 → 1

$$\hat{\mathbf{a}}_{\text{SBE}}^{21} = \arg \min_{\mathbf{a}} [S_1(\mathbf{a}) + B_2(\mathbf{a})]$$

Theorem (1995)

1) Linear regression

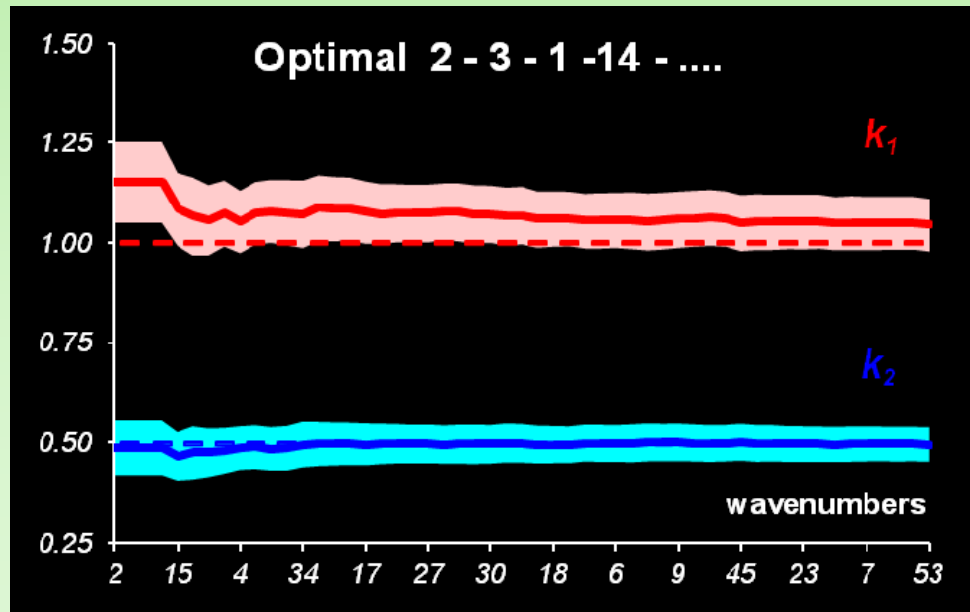
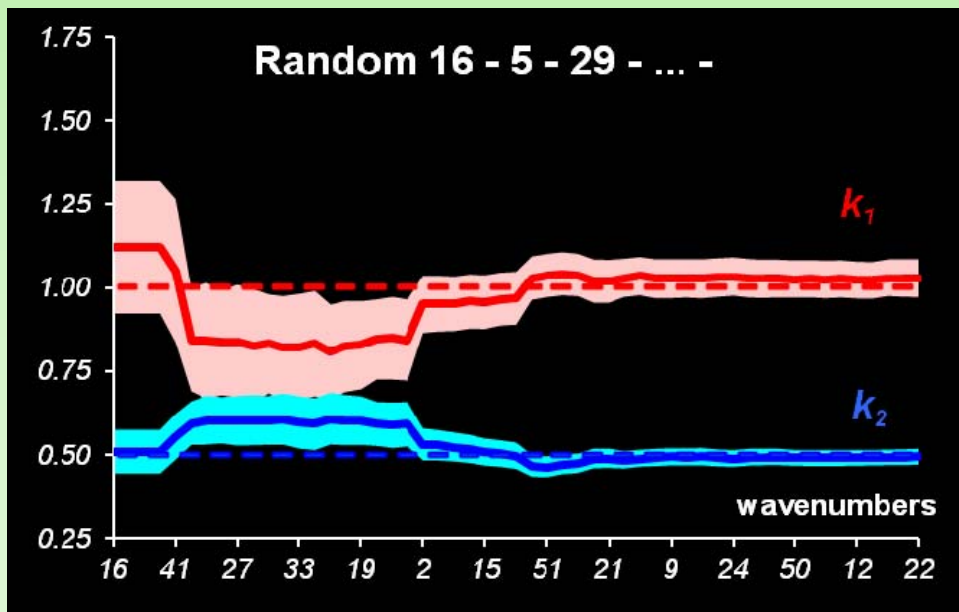
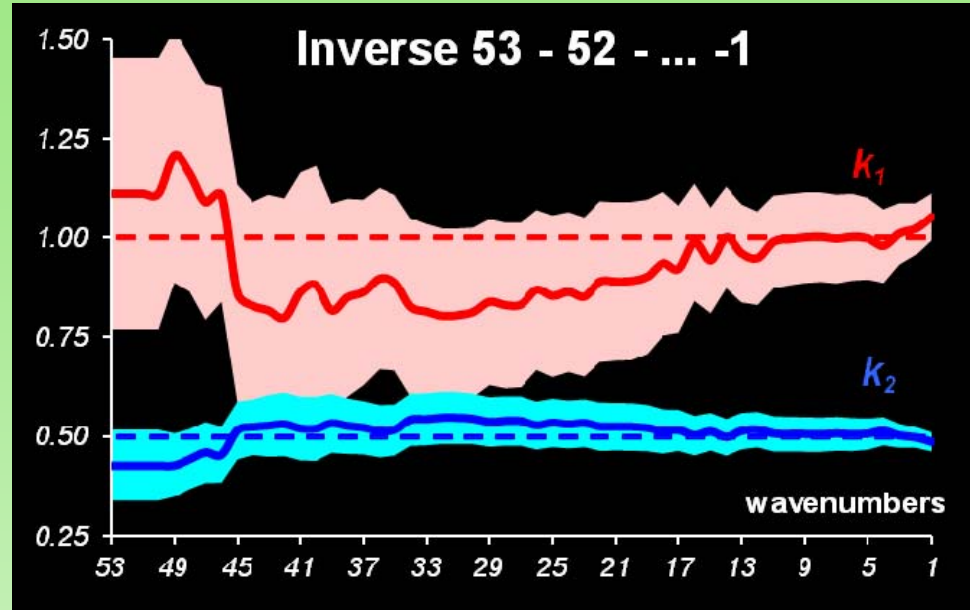
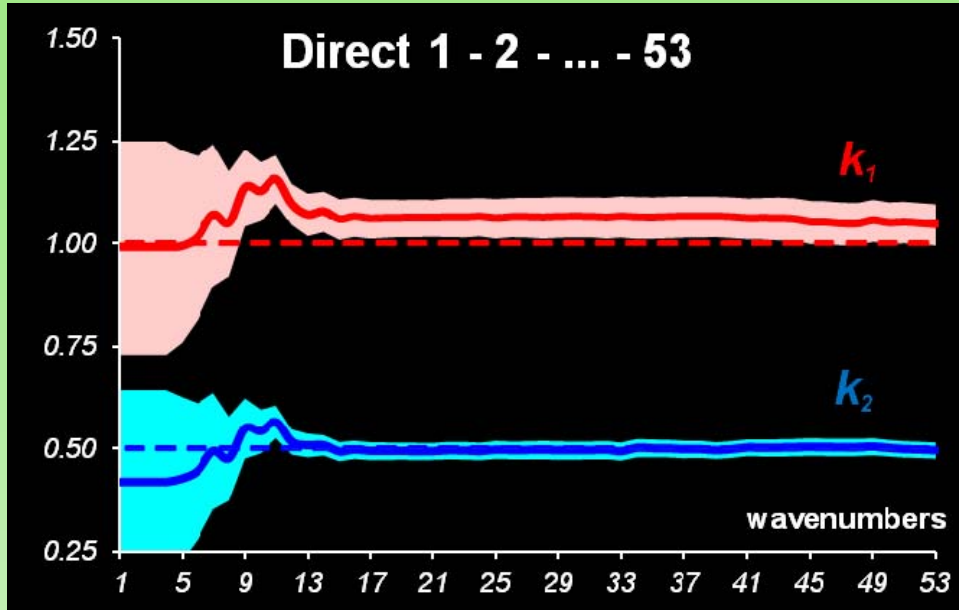
$$\hat{\mathbf{a}}_{\text{SBE}}^{12} = \hat{\mathbf{a}}_{\text{SBE}}^{21} = \hat{\mathbf{a}}_{\text{OLS}}$$

2) Non linear regression

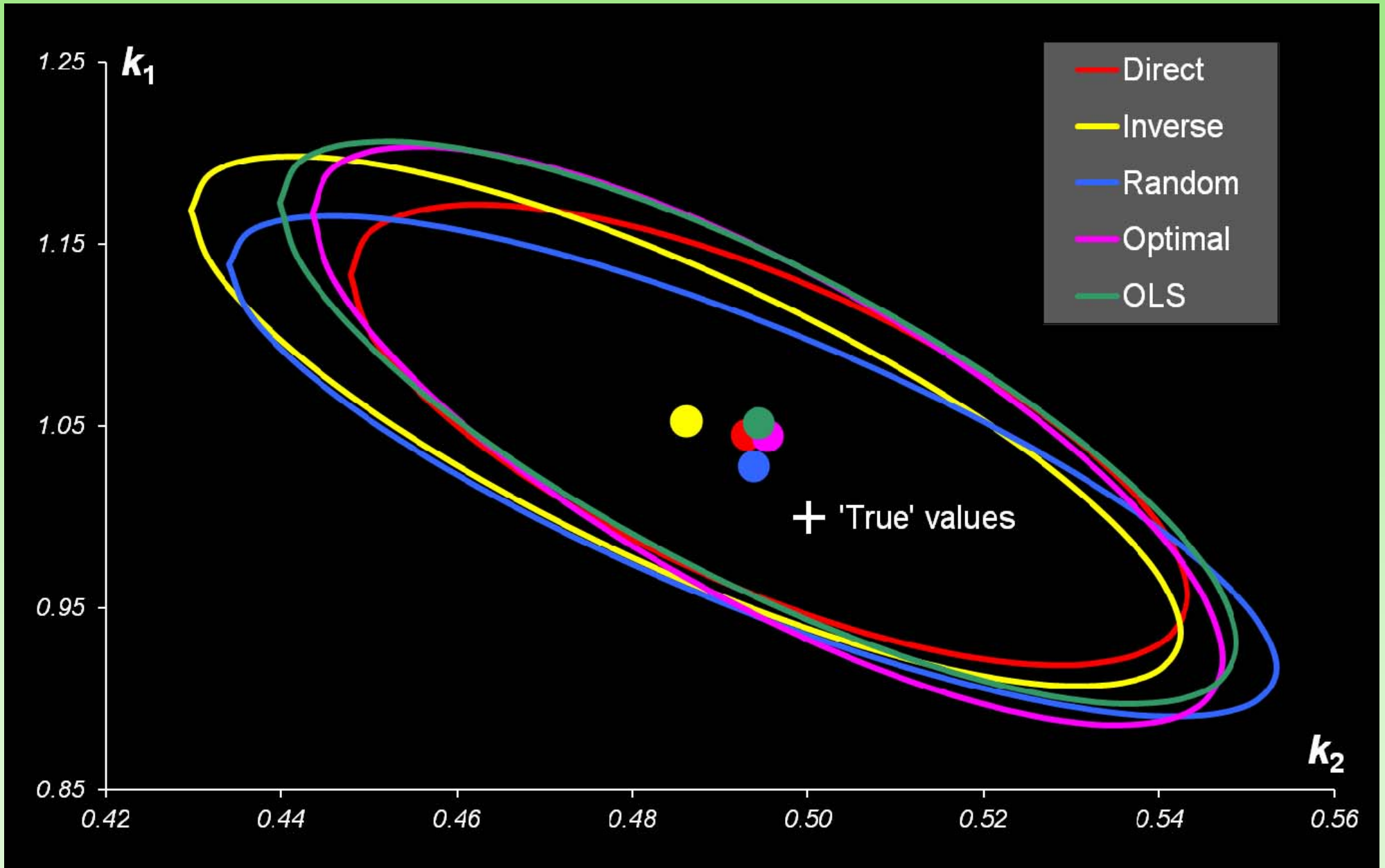
$$\hat{\mathbf{a}}_{\text{SBE}}^{12} \rightarrow \hat{\mathbf{a}}_{\text{OLS}} \leftarrow \hat{\mathbf{a}}_{\text{SBE}}^{21} \quad \text{at } N \rightarrow \infty$$

Example: a hard MCR

$A \rightarrow B \rightarrow C$

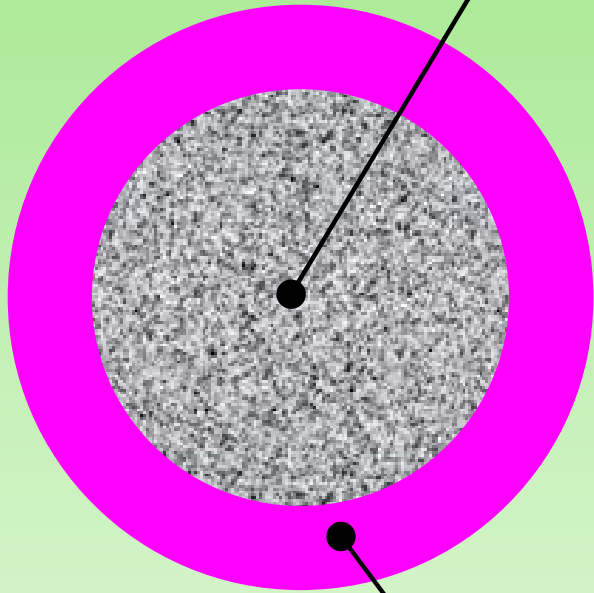


Example: a hard MCR



Case study: Pellets' release

Sugar +API



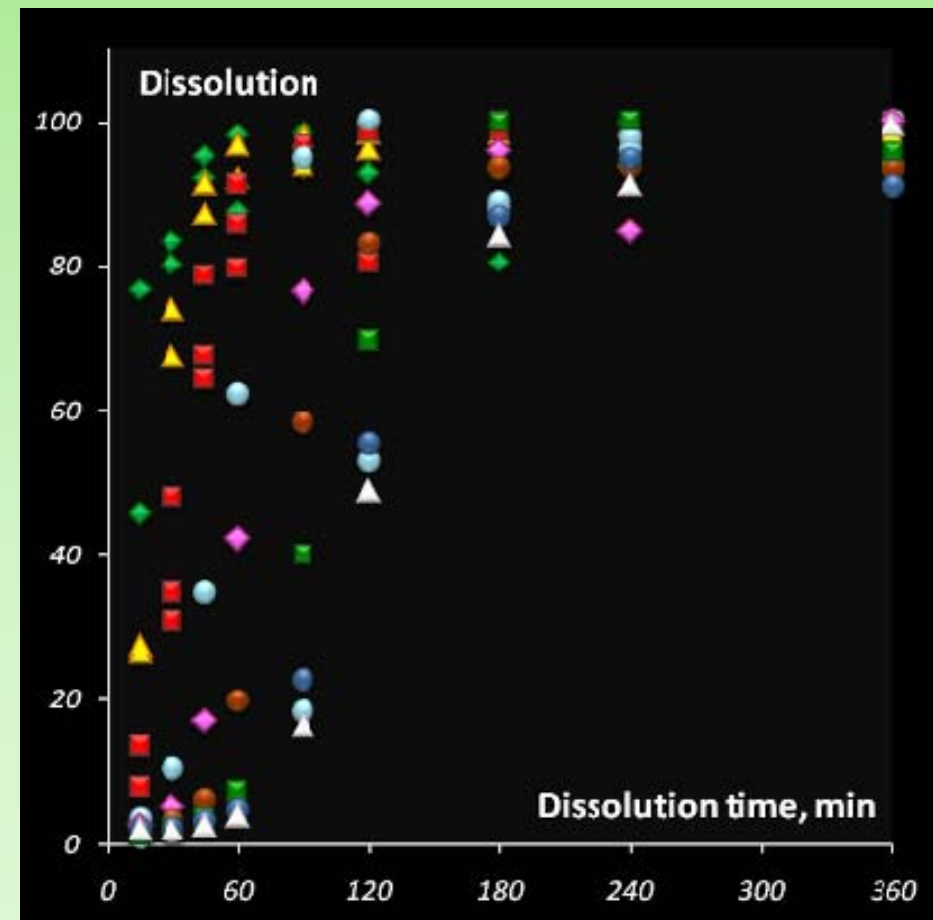
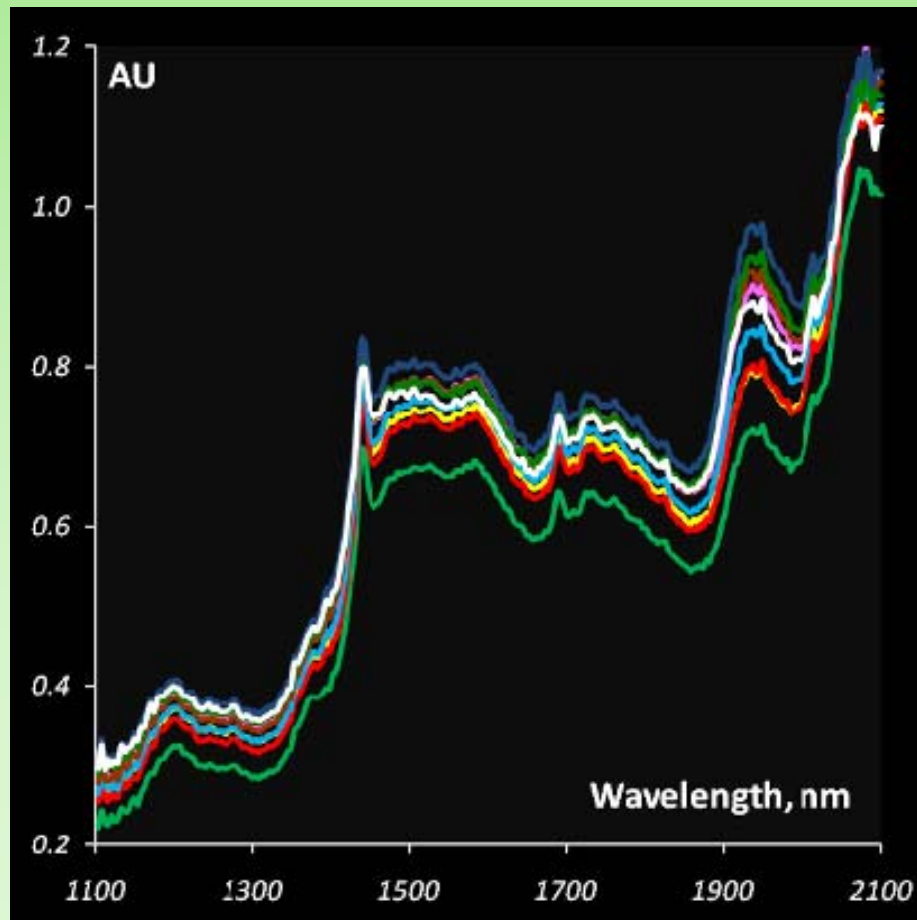
Coating: Acryl EZE

Experiment

NIR Spectra

Dissolution Profiles

$t = 105$



Dissolution data

White dyer

Yellow dyer

Batch W1

Batch Y1

Batch W2

Batch Y2

Batch W3

Batch Y3

Batch W4

Batch Y4

Batch W5

Batch Y5

Batch W6

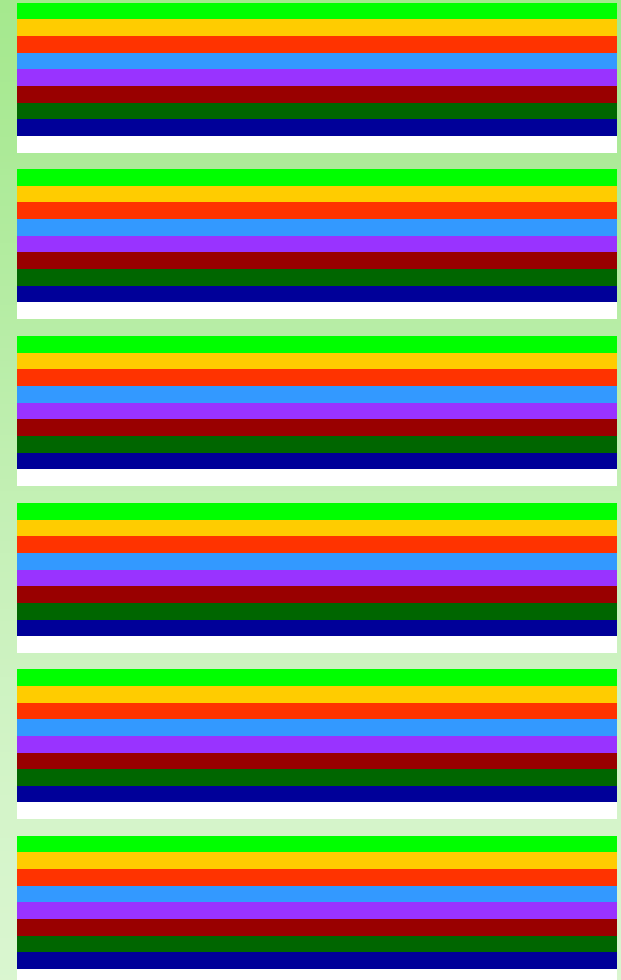
Batch Y6

Batch W7

28.02.12

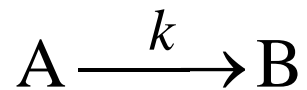
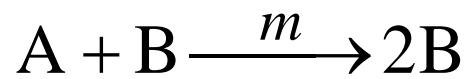
WSC-8

18



Autocatalysis

$$\varphi(t, m, k) = 100k \frac{\exp[(m+k)t] - 1}{m + k \exp[(m+k)t]}$$

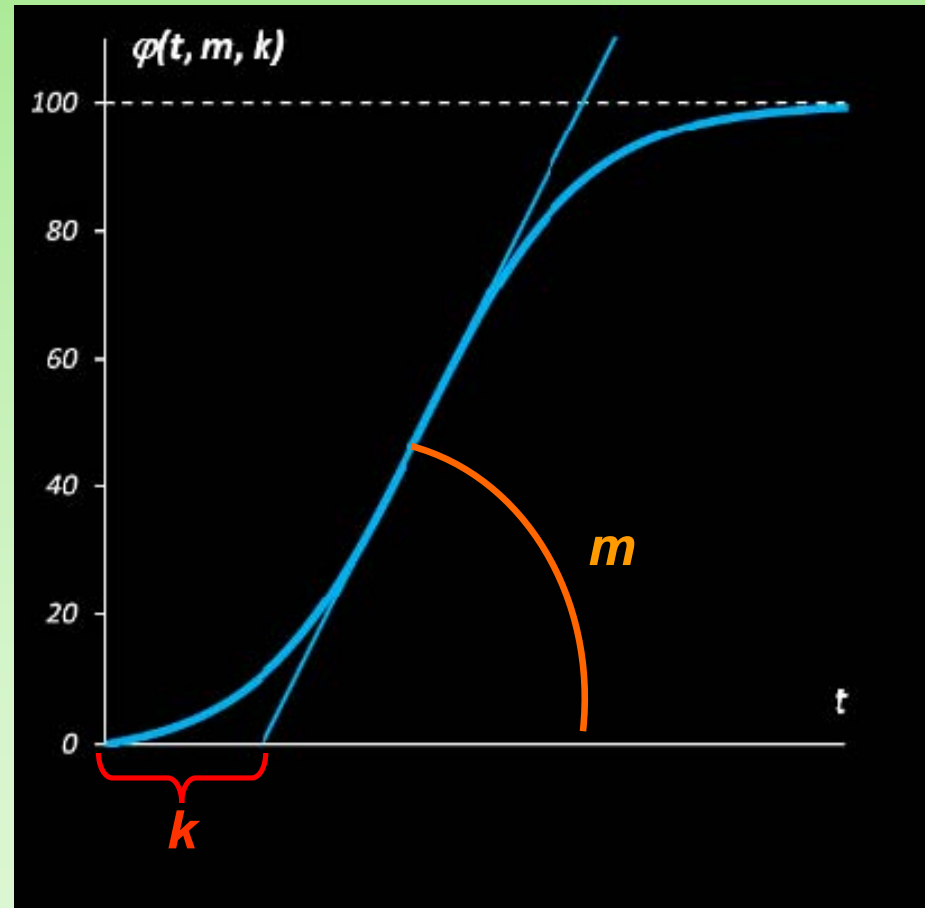


$$[A] + [B] = 100$$

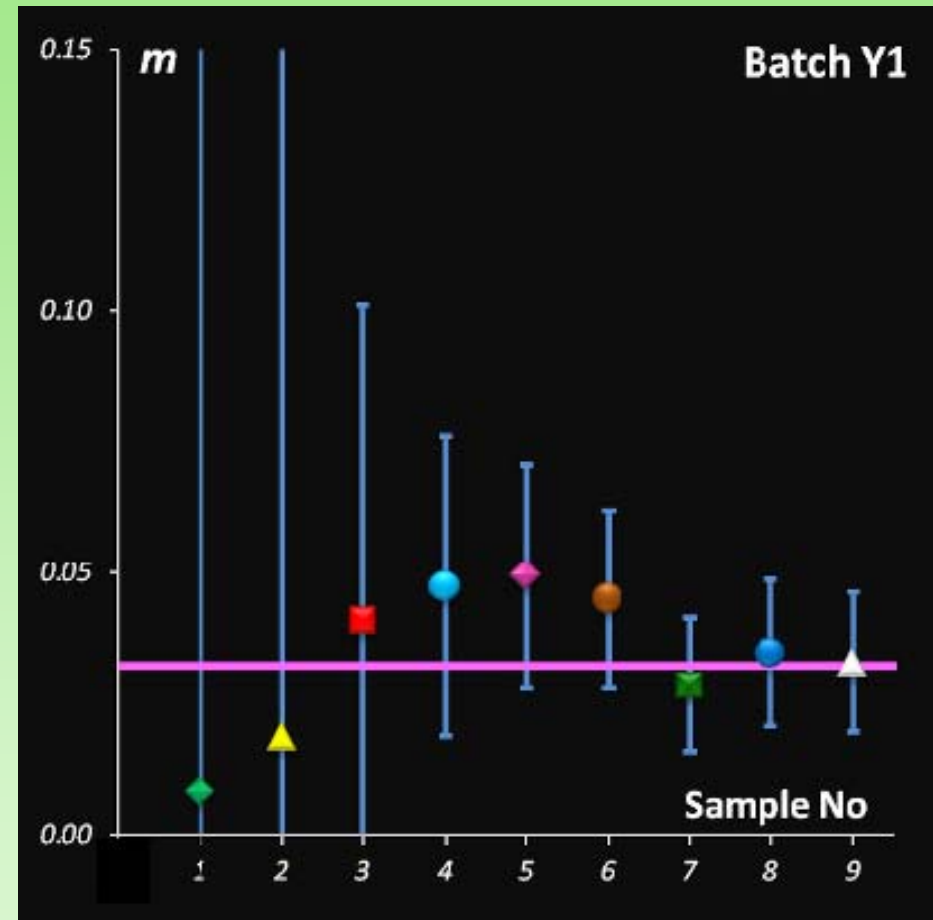
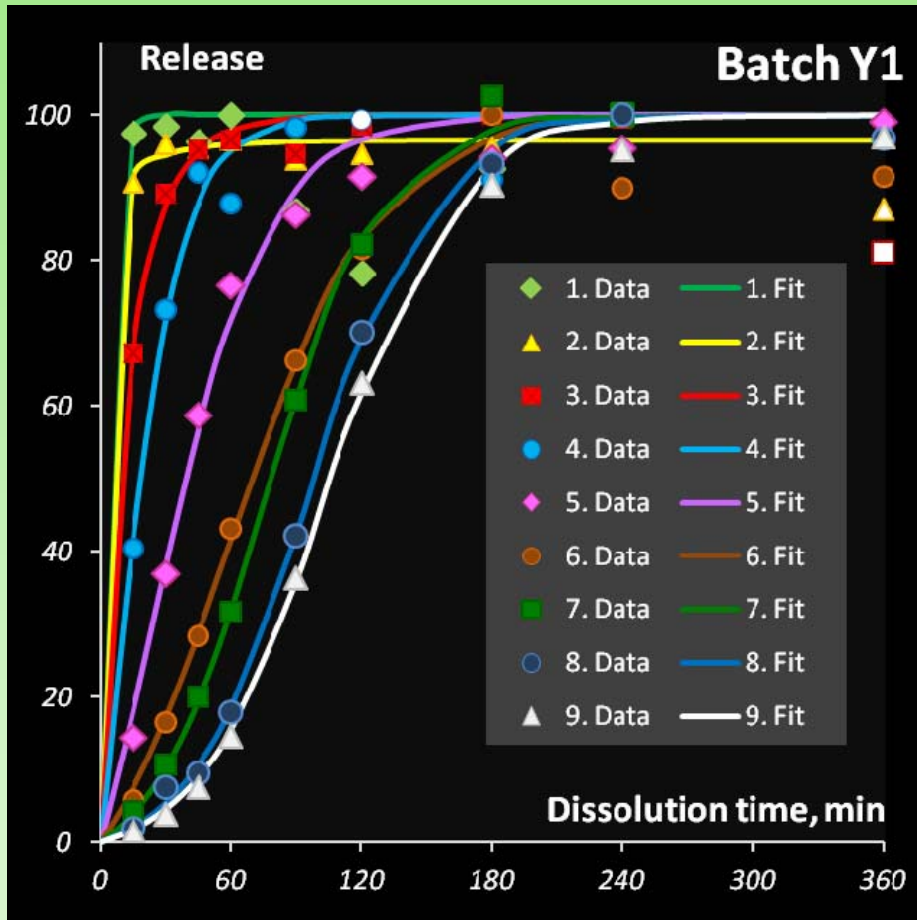
$$[B](0) = 0$$

$$\varphi = [B]$$

**(7+6)*9 k 's + (7+6)*9 m 's =
234 unknown parameters**

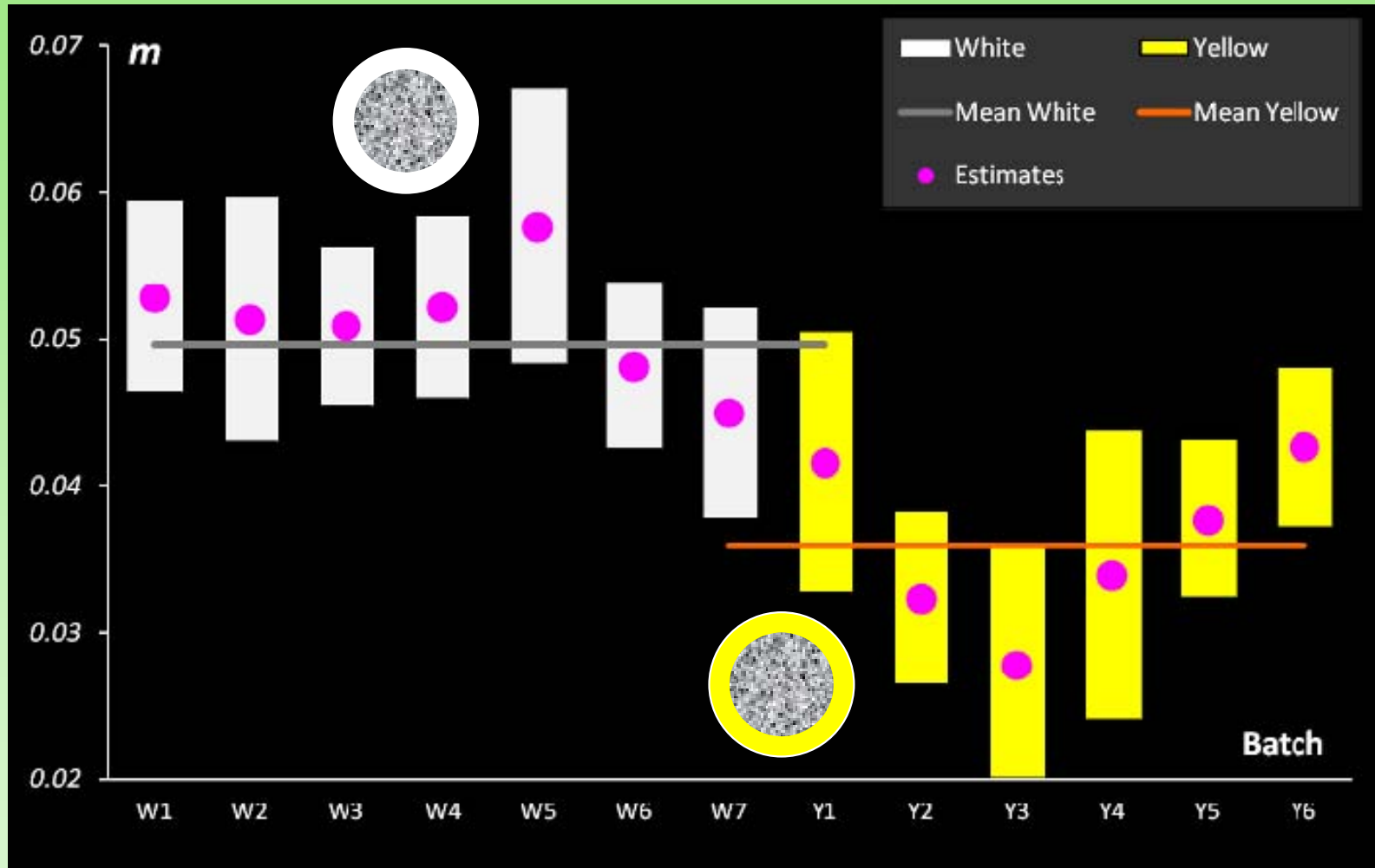


Parameter m is common within a batch



$(7+6)*9$ k 's + $(7+6)$ m 's = 130 unknown parameters

Parameter m and the layer grade



$(7+6)*9$ k 's + $(1+1)$ m 's = 119 unknown parameters

White dyer batches

$$\varphi(t, m, k) = 100k \frac{\exp[(m+k)t] - 1}{m+k \exp[(m+k)t]}$$

Batch W1



$k_{11}, k_{12}, \dots, k_{19};$ m

Batch W2



$k_{21}, k_{22}, \dots, k_{29};$ m

Batch W3



$k_{31}, k_{32}, \dots, k_{39};$ m

Batch W4



$k_{41}, k_{42}, \dots, k_{49};$ m

Batch W5



$k_{51}, k_{52}, \dots, k_{59};$ m

Batch W6



$k_{61}, k_{62}, \dots, k_{69};$ m

Batch W7



$k_{71}, k_{72}, \dots, k_{79};$ m

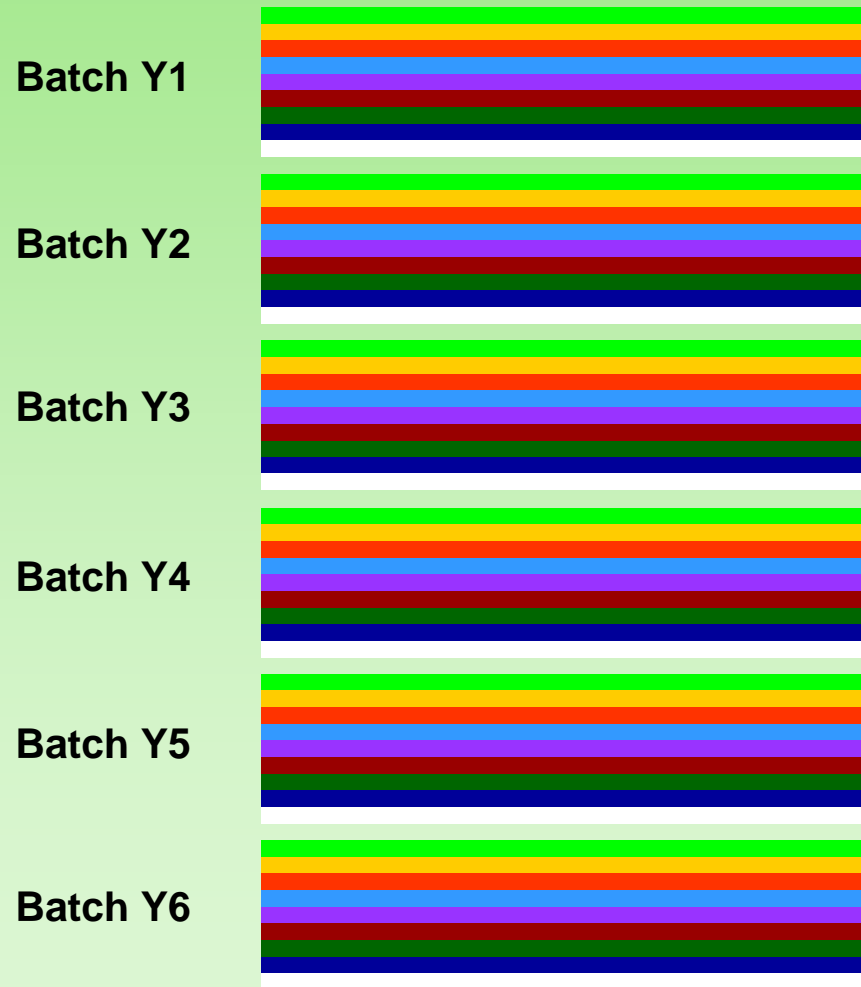
28.02.12

WSC-8

22

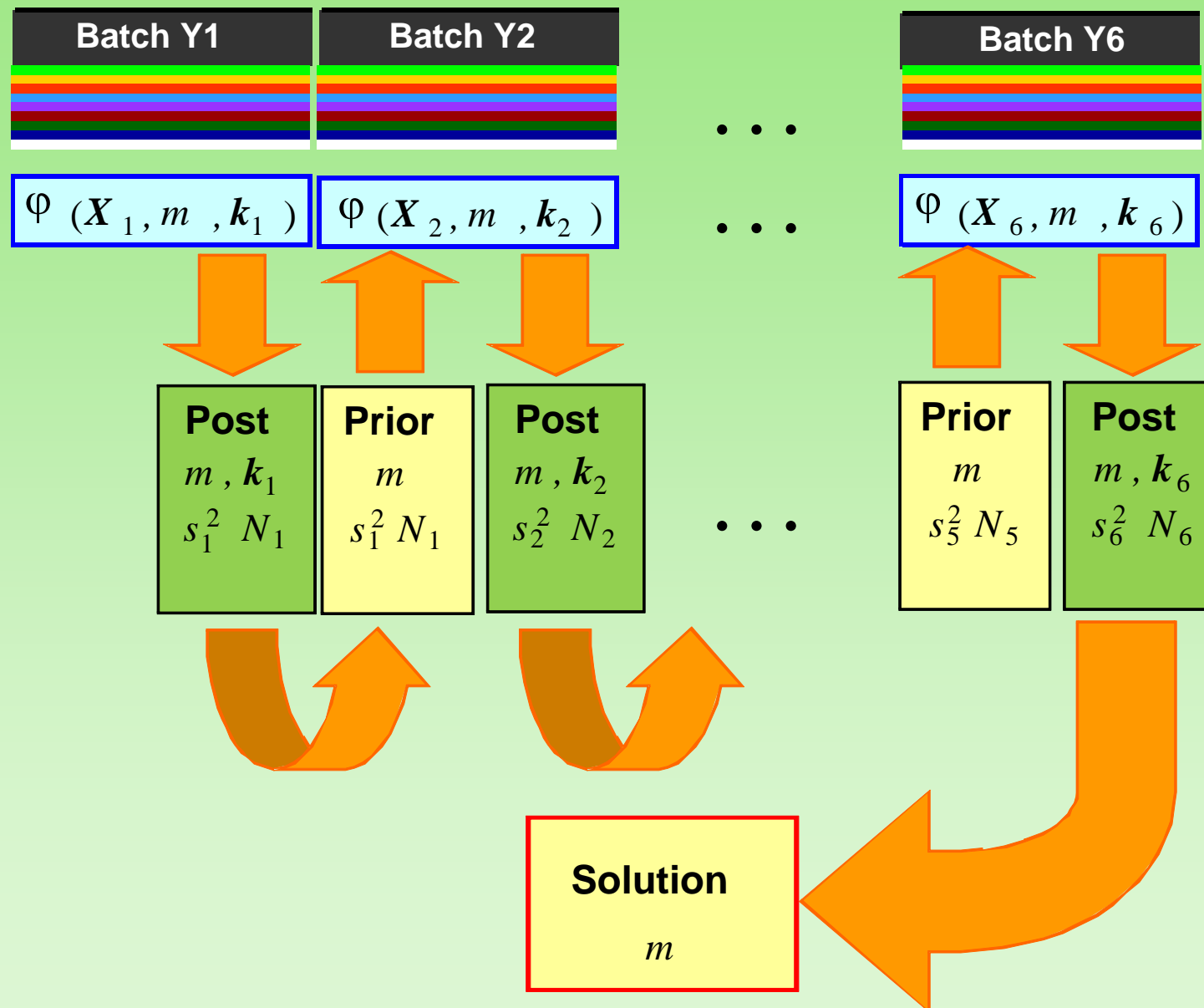
Yellow dyer batches

$$\varphi(t, m, k) = 100k \frac{\exp[(m+k)t] - 1}{m+k \exp[(m+k)t]}$$

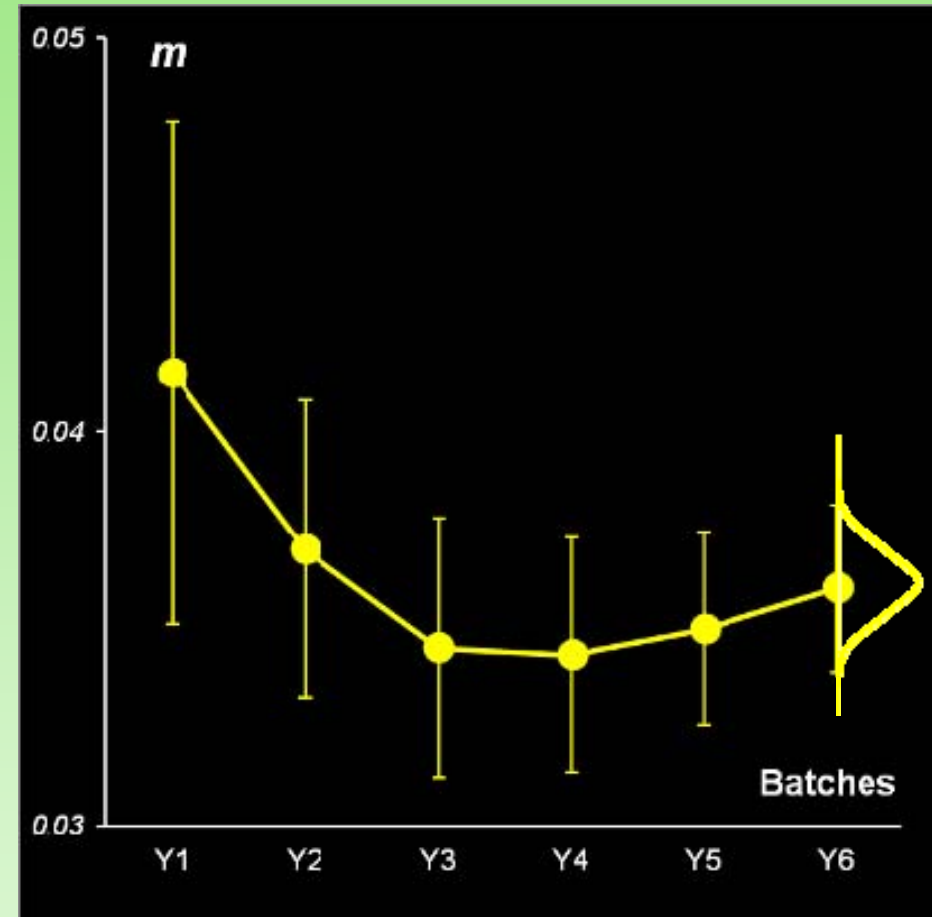
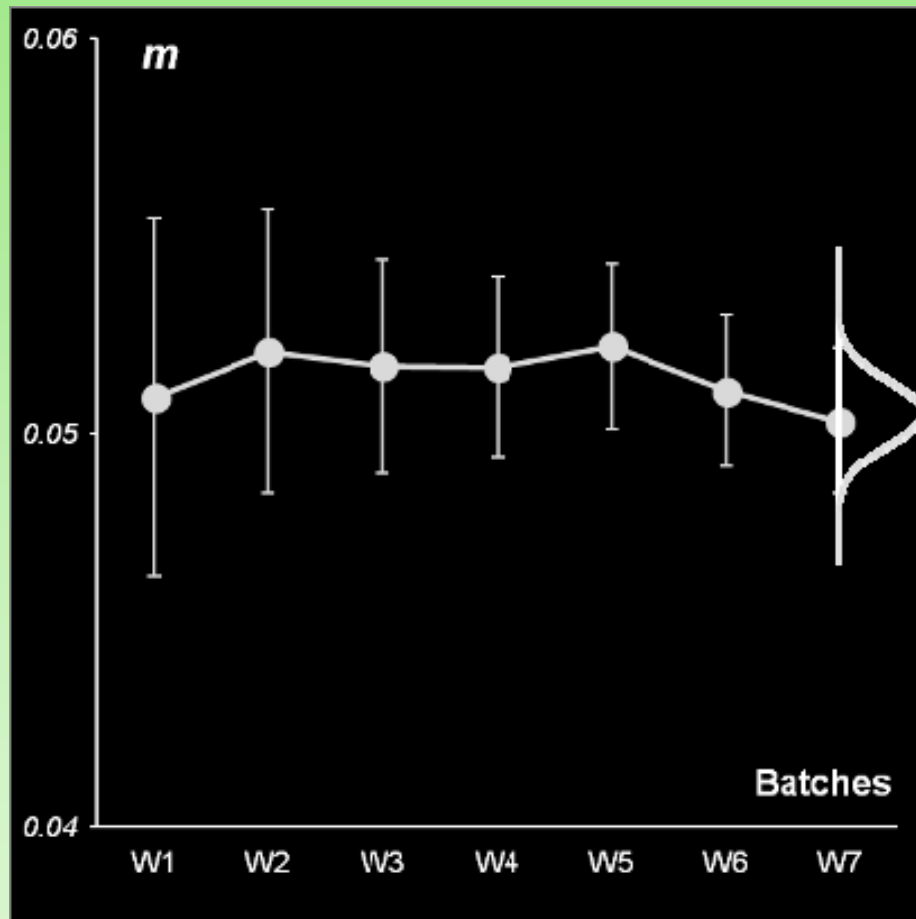


$k_{11}, k_{12}, \dots, k_{19};$	m
$k_{21}, k_{22}, \dots, k_{29};$	m
$k_{31}, k_{32}, \dots, k_{39};$	m
$k_{41}, k_{42}, \dots, k_{49};$	m
$k_{51}, k_{52}, \dots, k_{59};$	m
$k_{61}, k_{62}, \dots, k_{69};$	m

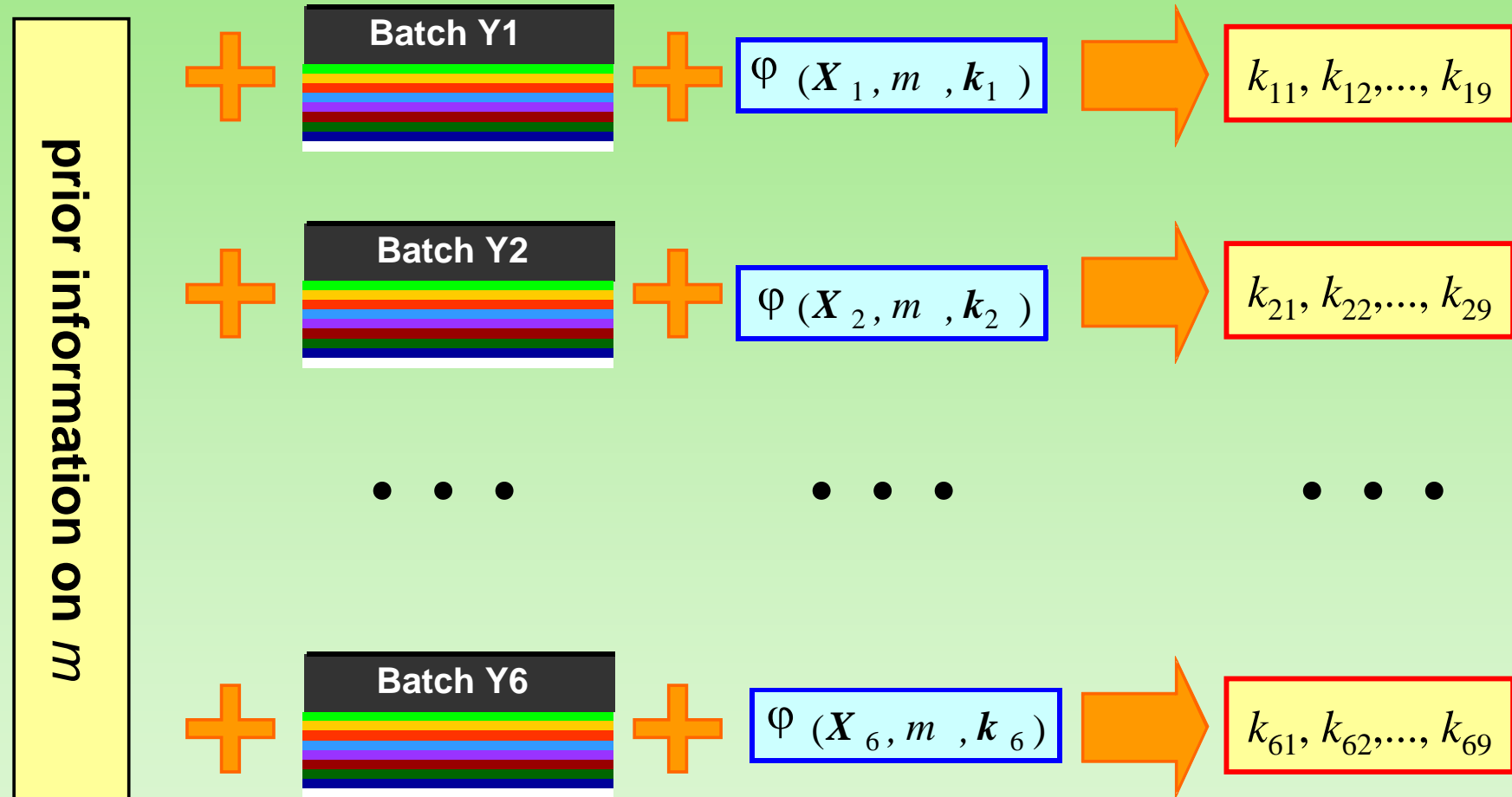
Forward SBE procedure



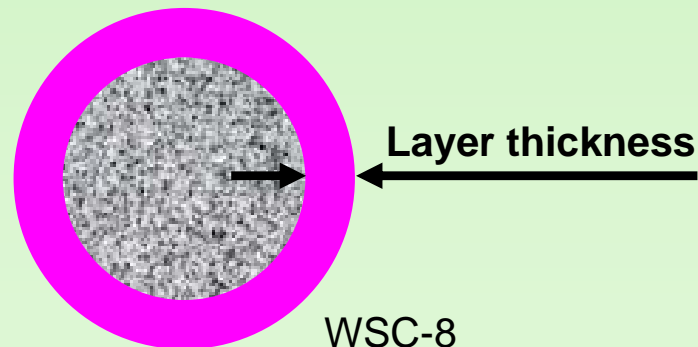
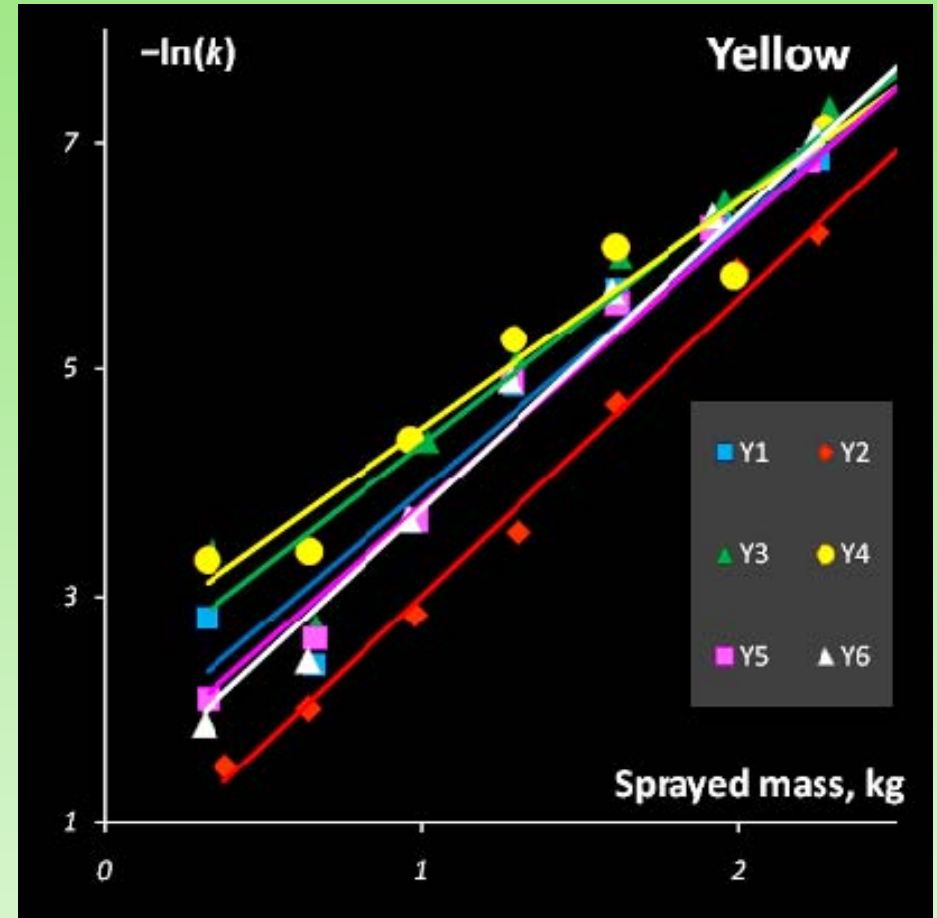
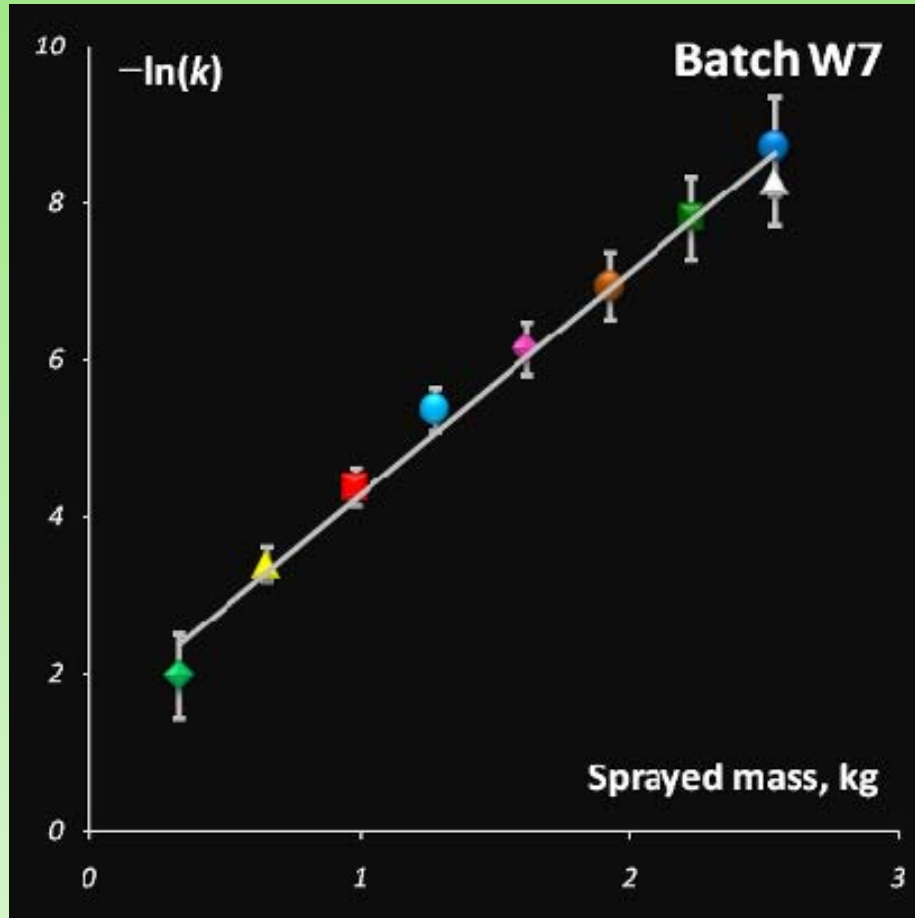
Successive estimates of m



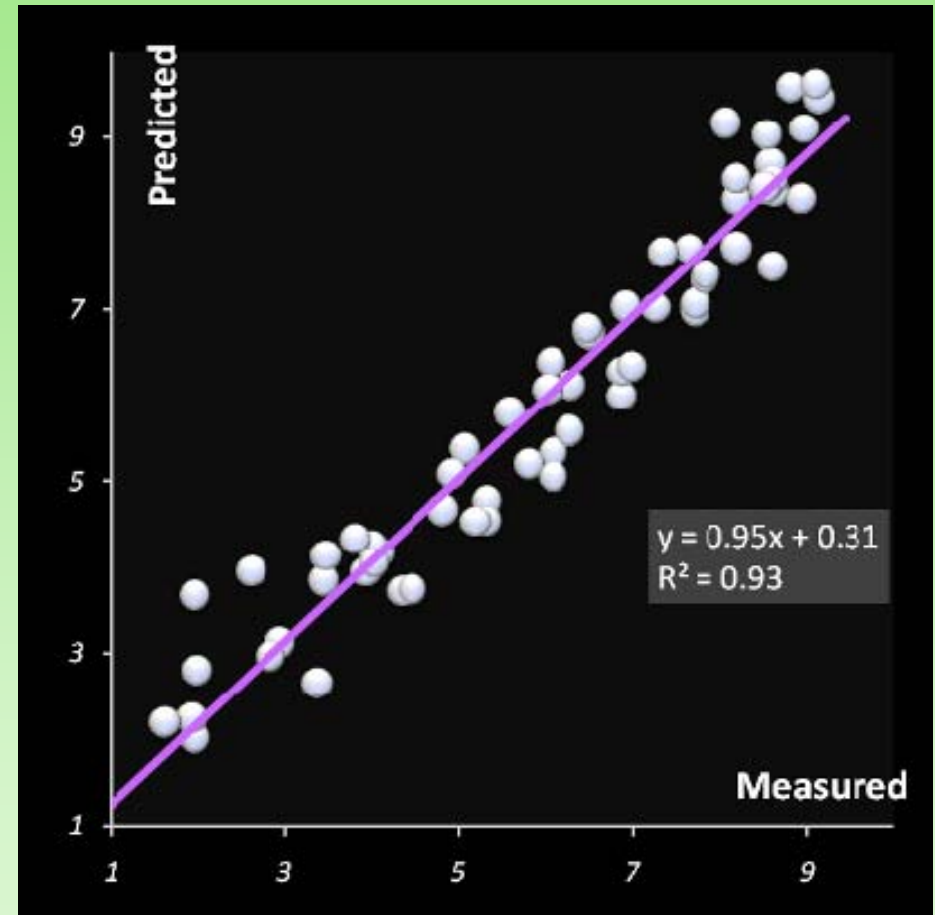
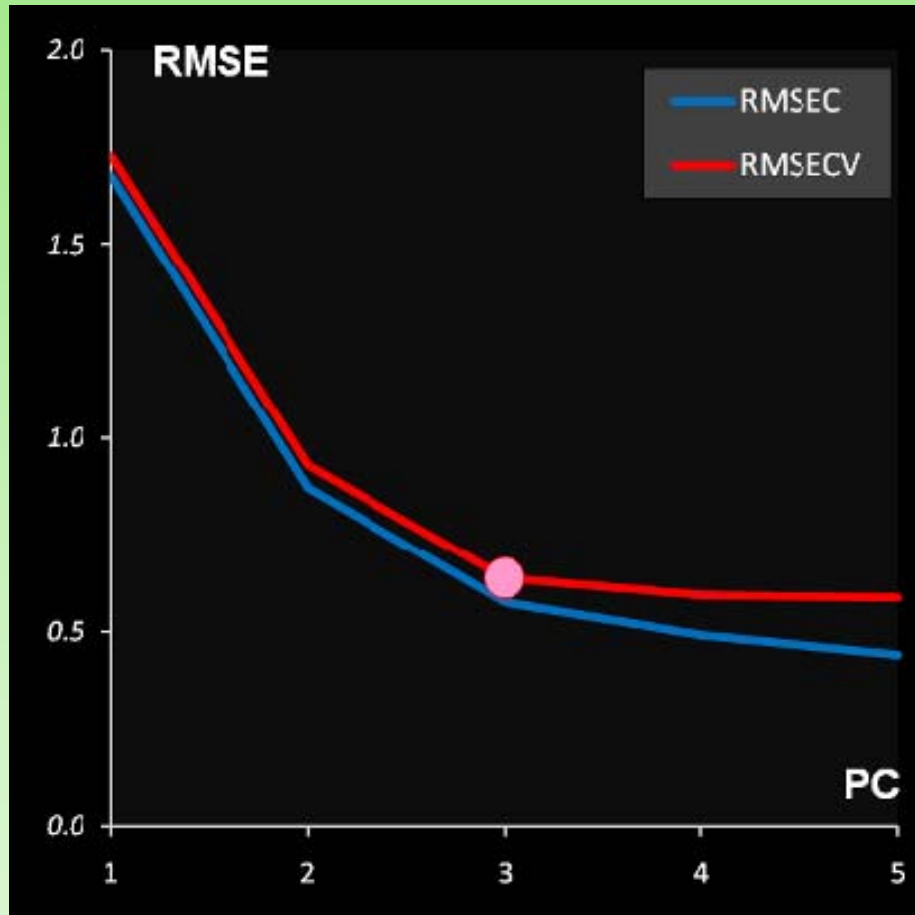
Backward SBE procedure



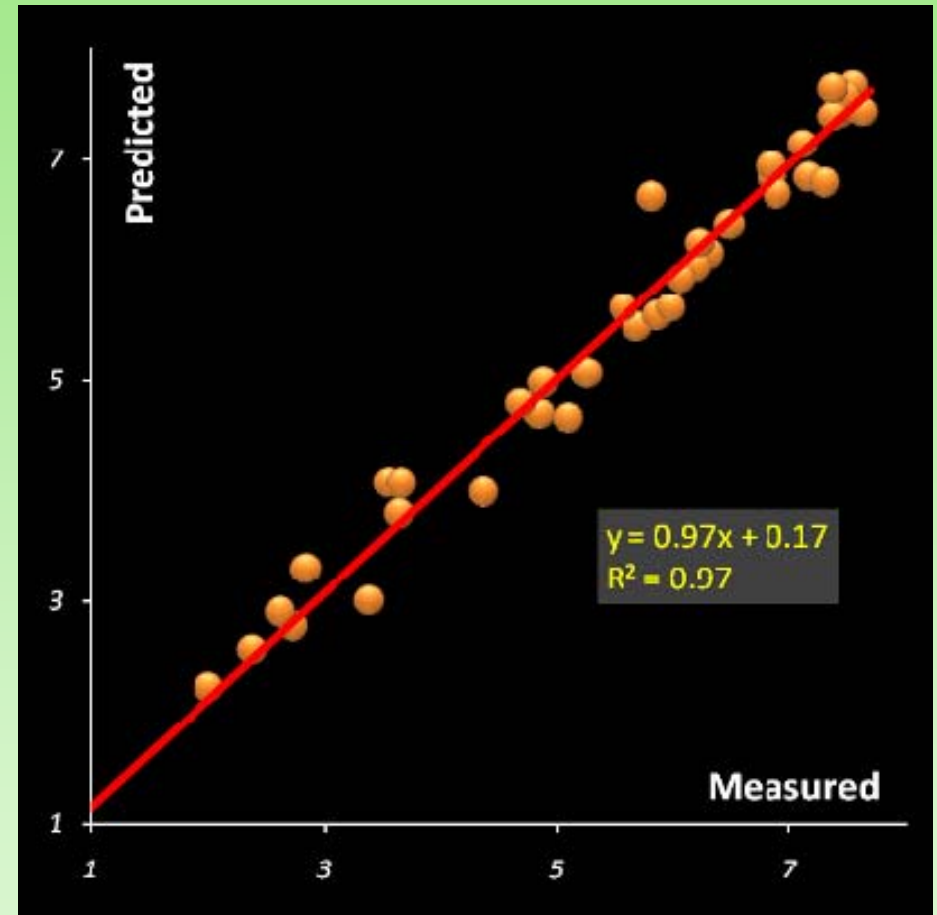
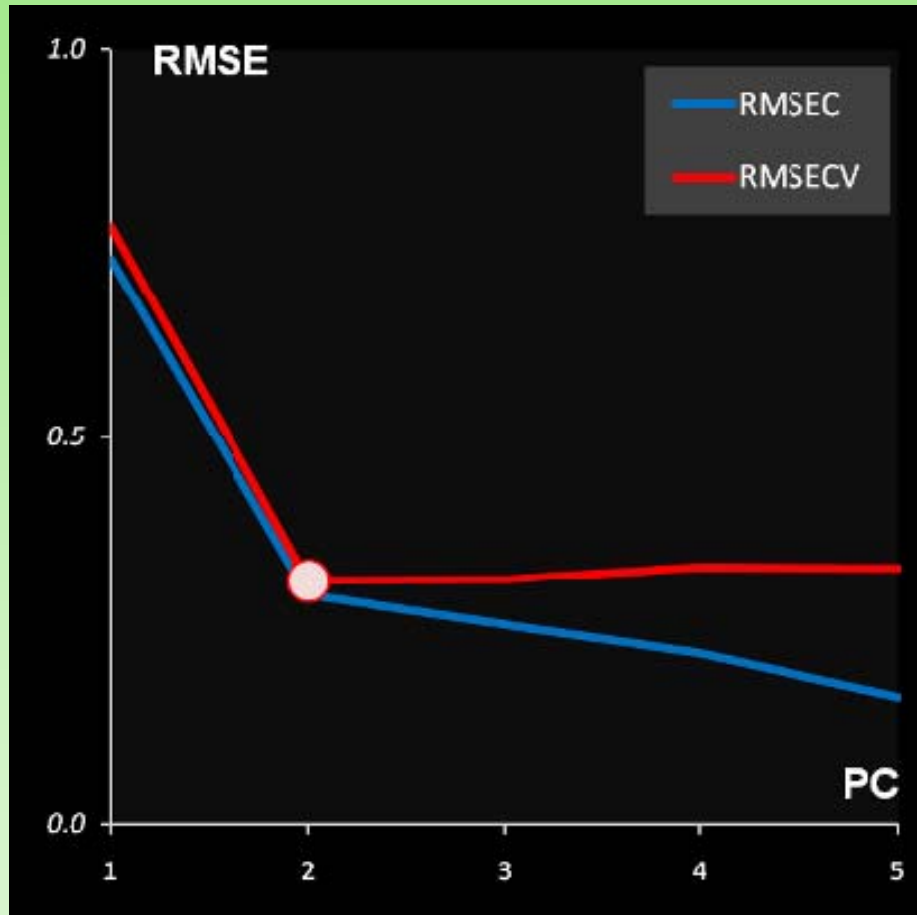
Parameter k and the layer thickness



NIR prediction of k : White subset



NIR prediction of *k*: Yellow subset



Conclusions

White models: formal kinetics, mechanism, rate constants, etc

Semen Spivak

Grey models: curve resolution, pure components, etc

Cyril Ruckebusch & Veli-Matti Taavitsainen

Black models: features extraction, 'pre-processing' for MVA

me