

Send me your .phz file(s) for the class project,  
 and any work files (e.g. xls for data entry) in  
a single .zip file ~ e.g.

Bilal - phz - Class - Project . zip  
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 @gmail.com

curtishayswhitson@gmail.com

Use ✓

Ackerman Sample Compositions:

Reliable Lab Data

Sep Oil ... C<sub>7+</sub> & ... C<sub>7</sub> C<sub>8</sub> C<sub>9</sub> C<sub>10</sub> C<sub>11+</sub>  
 ↳ ↳

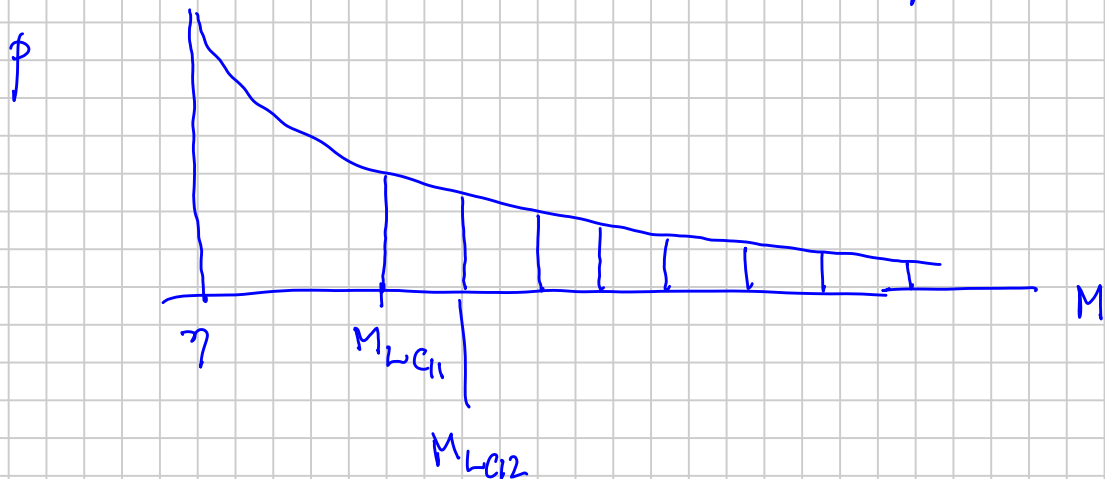
Sep Gas ... C<sub>7+</sub> M<sub>7+</sub> = 102

Exp Split AVG = 102 ⇒

C<sub>7</sub> C<sub>8</sub> C<sub>9</sub> C<sub>10</sub> C<sub>11+</sub>

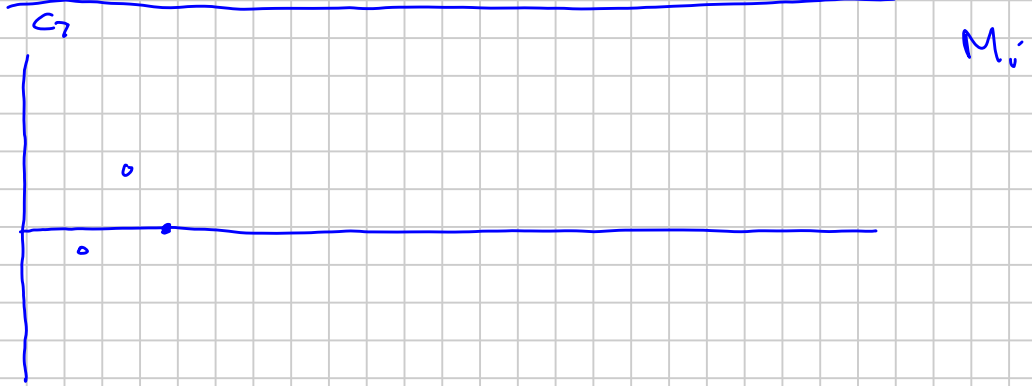
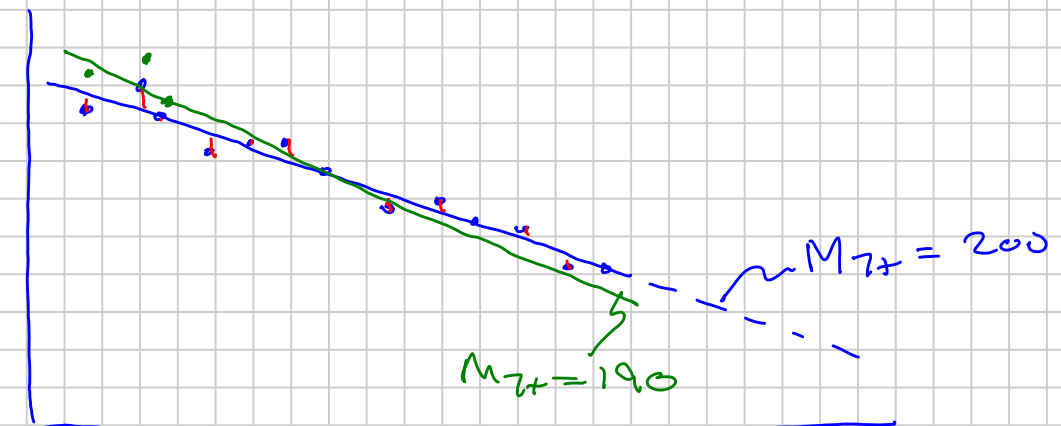
Gamma Fit C<sub>7+</sub>

$\bar{M}_{7+}$   $\gamma$   $\alpha_{7+}$

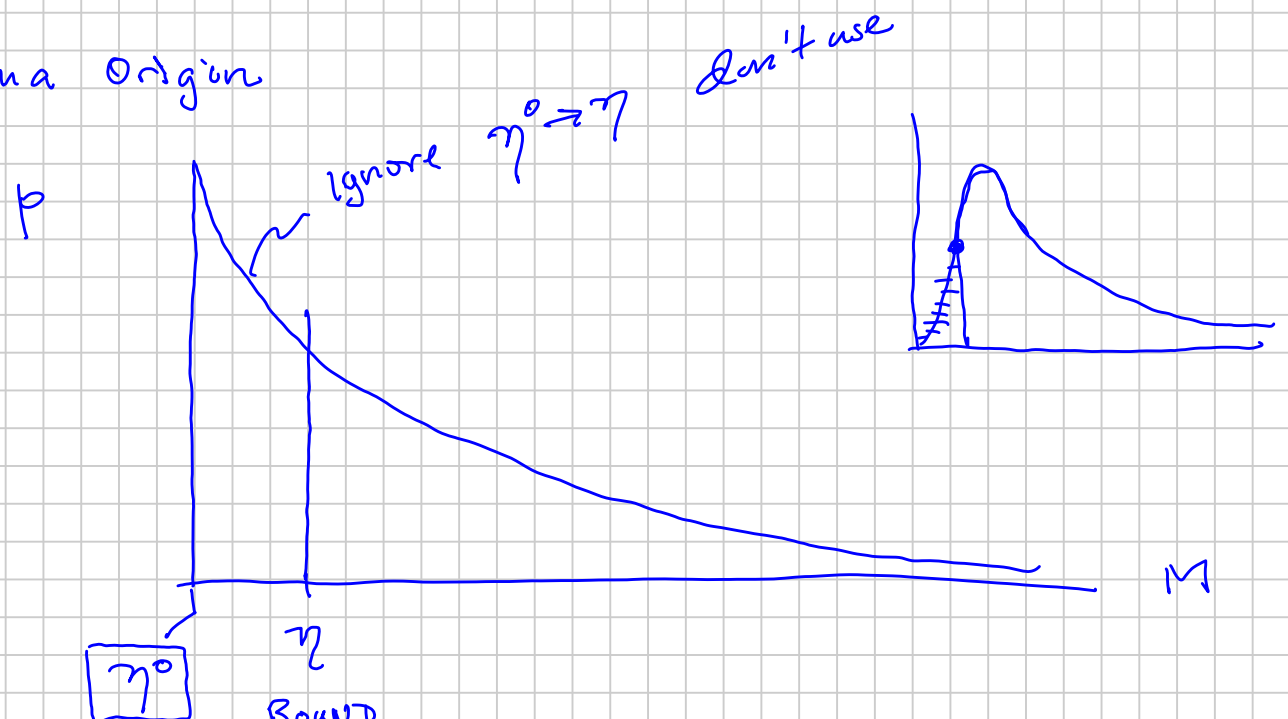


$M$

$Z_i (\log)$



Gamma origin



ORIGIN  
=  $\frac{\eta^0}{\eta}$

Gamma Model

AVERAGE  
SHAPE

$\eta^0 = (\text{ORIGIN} * \text{BOUND})$

# Phase Comp SCN MW "method"

NAME MW

CI I = 6 7 8 ...

Two Correlations:

n-Paraffin Correlations

$$M_p(T_b) \checkmark$$

$$T_{cp}(T_b)$$

$$p_{op}(T_b)$$

Fraction (SCN)  
w/ non-Paraffins

$$M(T_b, \sim \Delta\gamma)$$

$$T_c(T_b, \sim \Delta\gamma)$$

$$p_c(T_b, \sim \Delta\gamma)$$

$$\Delta\gamma = \gamma(T_b) - \gamma_p(T_b)$$

Soreide  $\gamma(\underline{M}, C_f)$   $\gamma(\underline{M}, a_0, a_1, a_2, c_3)$

CI: ①  $T_{bp}(M_{PI} = 14 \cdot I + 2)$

②  $\gamma_p(T_{bp})$

③  $\gamma(M_{PI}, C_f) \leftarrow$  Soreide

④  $\Delta\gamma = \gamma - \gamma_p \quad \left. \vphantom{\Delta\gamma} \right\}$

⑤ Fractional (effective) SCN  $i \leq I$

⊗

$$i = I - 3\Delta\gamma$$

e.g.  $I = 7 \quad i = 6.7$  (default Soreide)

$$\textcircled{6} \quad M_{pi} = 14i + 2 \quad \gamma_{pi}(T_{bi})$$

Phz TB

$$\textcircled{7} \quad T_{bi} = T_{bp}(M_{pi})$$

$T_{wumw} = 1$   
(default)

$$\textcircled{8}_a \quad \left. \begin{array}{l} M_{i \leftarrow} (M_{pi}, \Delta\gamma) \xrightarrow{T_{wu}} \\ \gamma_i \leftarrow (M_i, c_f) \xrightarrow{\text{Solve}} \end{array} \right\} \text{Solve for } M_i$$

$T_{wumw} = 0$   
(often best)

$$\textcircled{8}_b \quad M_i = M_{pi}, \text{ i.e. } \Delta\gamma = 0$$

Any  
 $T_{wumw}$

$$\left\{ \begin{array}{l} \textcircled{9} \quad T_{ci}(T_{bi}, \Delta\gamma) \\ \textcircled{10} \quad p_{ci}(T_{bi}, \Delta\gamma) \end{array} \right.$$

Any experiment defined in Phz (XXX = CCE SEP..)

Mix FEED {what you want the cell filled with initially}  
TEMP 212 F  
{ PRES 6000 PSIA }  
PSAT } Used to define PVT properties called "Relative"

XXX

BASIS {vol, mass, mole}  
Liquid-dependent

$V_{ref} = \text{Flash calc from EOS}$   
w/ ref. amount  
at Pref Tref  
Total Volume

LRVOL

GRVOL

RVOL

↑  
Relative

LRMASS LRMa

GRMASS GRMDL

# Defining Mismatch

Individual (Groups)  
of Weight Factors

<u>i</u>	<u>d<sub>mi</sub></u>	<u>d<sub>ci</sub></u>	<u>Types</u>	<u>Importance</u>	<u>Accuracy</u>
			p <sub>s</sub>	H	H
			q	A	A
			v <sub>r</sub>	L	L
			y <sub>h</sub>		
			x <sub>i</sub>		
			⋮		

N

Difference  $\delta_i = d_{ci} - d_{mi}$  units

$\delta_i^2$  units<sup>2</sup>

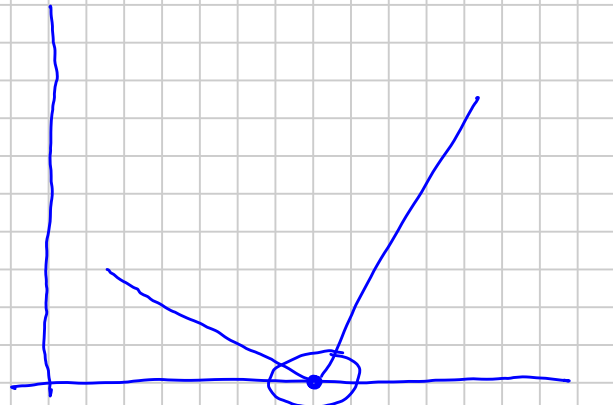
Deviation  $\Delta_i = \frac{d_{ci} - d_{mi}}{d_{mi}}$  ±

$$MM_1 = \frac{1}{N} \sum_{i=1}^N \Delta_i = 0 \quad \text{could still be a bad model}$$

$$MM_2 = \frac{1}{N} \sum_{i=1}^N |\Delta_i|$$

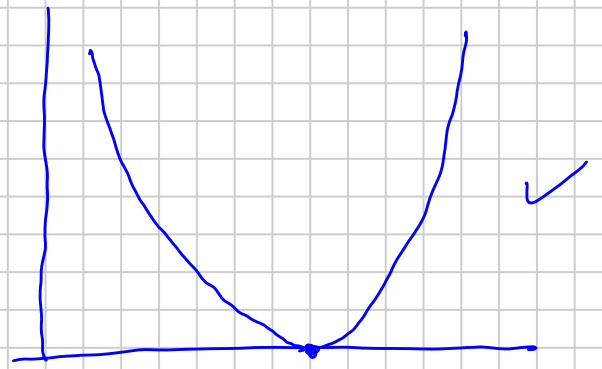
Model  $y = ax$

MM<sub>2</sub>



non-continuous  
derivative

$$MM_3 = \left( \frac{1}{N} \cdot \sum \Delta_i^2 \right)^{1/2}$$



Data Types:

$$\frac{d_{T1 \text{ max}}}{d_{T1 \text{ min}}} = 1.5$$

100

$$\frac{d_{T4 \text{ max}}}{d_{T4 \text{ min}}} = 10^6$$

$P_0$

$P_9$

$K_i$

$y_i$

$x_i$

$V_{r0}$

$\gamma_{C1}$

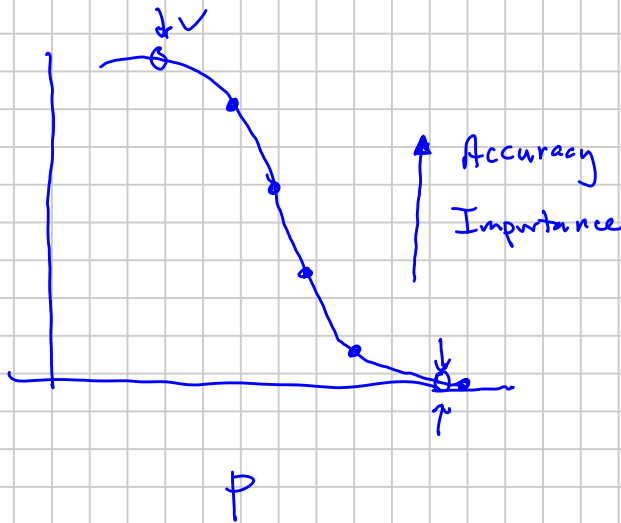
$\gamma_{C30+}$

$10^{-3}$

$\sim 1$

$10^{-6}$

0.5



5 error % in  $C_1$

vs

600

5 error % in  $C_{30+}$

$T4(a) C_1$

$T4(x) C_{30+}$

Residual:  $r_i = \frac{d_{ci} - d_{mi}}{d_{ref,i}}$

built-in weighting  
of "larger" data  
of a given type

$d_{ref,i} = \max \{ d_i \in T4(a) \}$   
e.g.

Each XXX

Each Property Column  
finds  $d_{ref,i}$

MM(FHz) ~

$$\sum_{i=1}^n (r_i \cdot w_i)^2$$

Main Regression

User specified

Program Controlled

@  $P_{min} \sim 1500 \text{ psia}$

Lab  $V_{ro}$

(1)  $\left(\frac{V_o}{V_s}\right)_{\text{Input PVT}}$

(2)  $\left(\frac{V_o}{V_t}\right)_{\text{lab}}$

$\left(\frac{x}{4}\right)$

\* Due Diligence on data entry \*

\* Ranking data: Quality (Accuracy) \*  
• Importance

min RMS

2-5%

$k_{ij} \quad T_c \quad P_c \quad G_{ro} \quad \dots \quad 5-20$

( EOS Model Parameters )

Individual Sample Compositions

$V_{c1} \quad V_{c2} \quad V_{c3}$

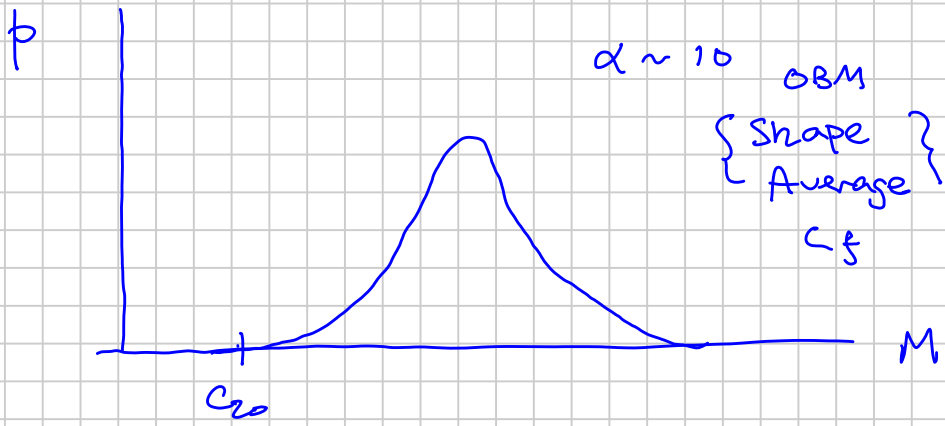
$M_{7+} \quad GOR \quad Z_{30+}$

Pick Variables (EOS | Sample-comp):

$V_i$  : starting value, min bound, max bound



OBM :



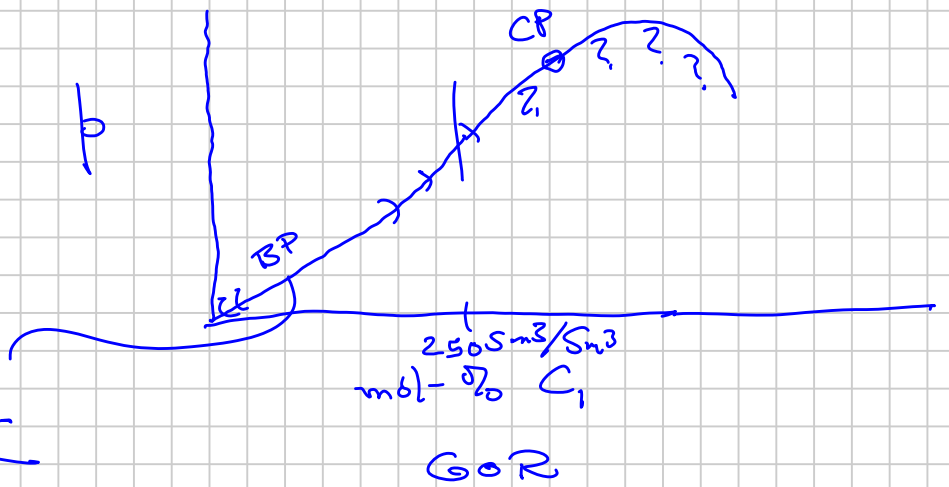
$[OBM + c_1]$

$(T, p)$

$[g(p, T)]$

$\underline{p}_s(c_1, T) :$

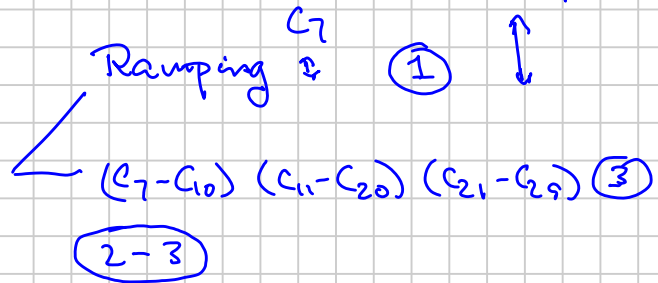
SRK :



# Regression Variables

① EOS Properties 3-7 variables

- $C_{7+}$   $T_c$   $P_c$   $\{T_b | w\}$
- $C_{N+}$   $T_c$   $P_c$   $\{T_b | w\}$



- {Soreide Model parameter ( $C_f$ )}  $\Rightarrow S_i (C_{7+})$  ①

- BIPs :  $C_1 - C_{7+}$  ①
- $C_1 - C_{N+}$  ①

{HC-HC} including  $C_{7+} - C_{7+}$  BEWARE!  
 {nm-HC - HC} only if "gas inj" PVT w/ nm-HCs

\* Affect all samples in a 'SIMILAR' way

② Sample Compositions (1-3) sample

- GOR recombination

Sep Gas  $\cdot V$  + Sep Oil  $(1-V)$

(a) Test sampler

(b) Flashed from wellstream

-  $M_{N+} \updownarrow \pm 5-10\%$

mass-to-mole conversion  $\updownarrow$   
 amount of  $C_{N+}$

-  $Z_{N+} \updownarrow$

\* Impact only the particular sample quality match

- Any composition change should improve most/all PVT predictions

- max note % ( $C_1, C_{7+}$ )  $\updownarrow \pm 0.5-1.5$  (2)

