

WEDNESDAY REVIEW

PVT Lab Tests (Conventional)

	<u>CCE</u>	<u>SEP</u>	<u>DLE</u>	<u>CVD</u>
⊙ Lab Procedures	*			
⊙ Measured data				
⊙ Reported quantities	*			
• Uncertainties	*			
⊙ QC	*			
⊙ Applications (direct)				

Ch. 6

- ⊙ Covered
- ⊙ Partly Covered
- Not discussed

CCE: Constant Composition (Mass)

SEP: Multi-Stage Separator Test

DLE: Differential Liberation

CVD: Constant Volume Depletion

* VERY important to tuning EOS model

* EOS MODEL TUNING

- STRATEGY
- MECHANICS
- EXAMPLE (CHOI)

* VISCOSITY MODEL TUNING

* BLACK-OIL PVT MODEL

- What
- How to Build from EOS
- Challenges & Problems
- When good enough, when not

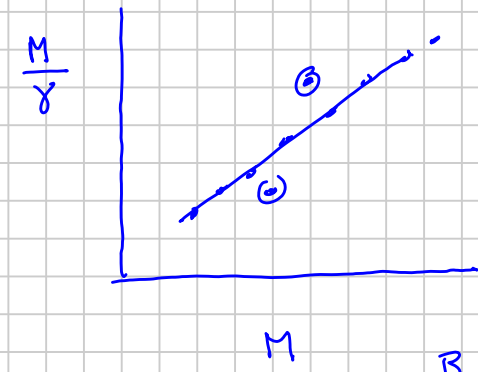
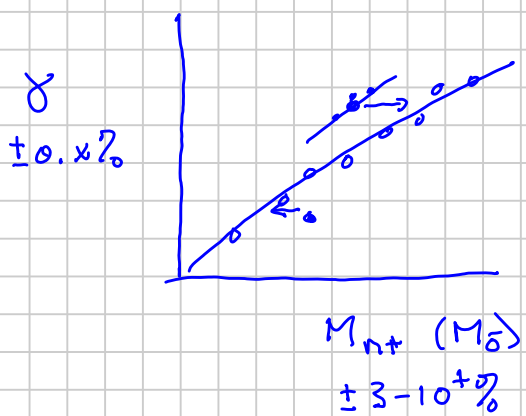
EOS MODEL TUNING STRATEGY

* Compositional Data QC (Excel)

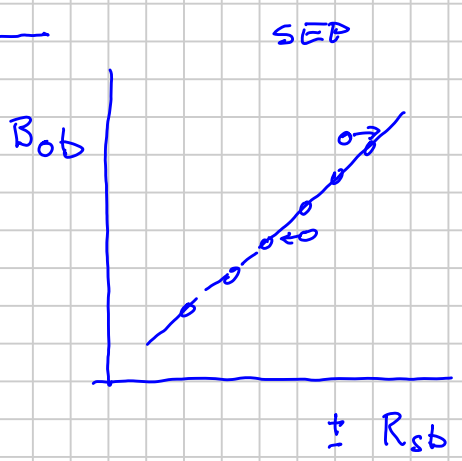
* — " — Enter to PVT Program (QC)

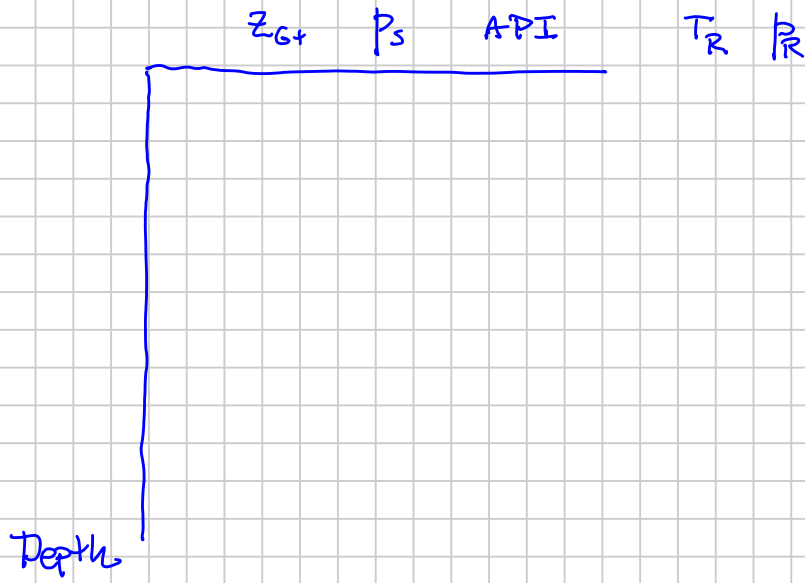
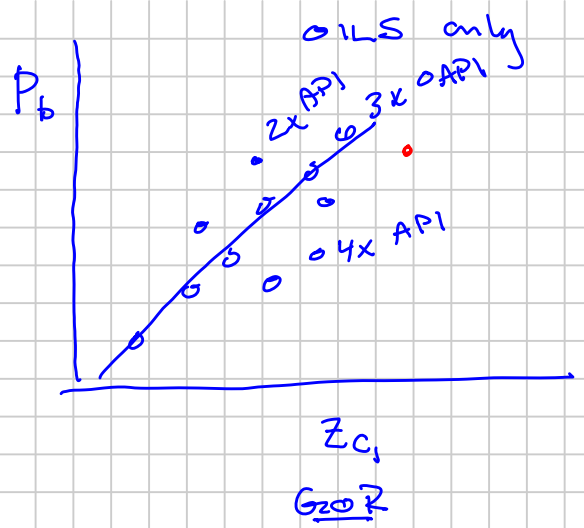
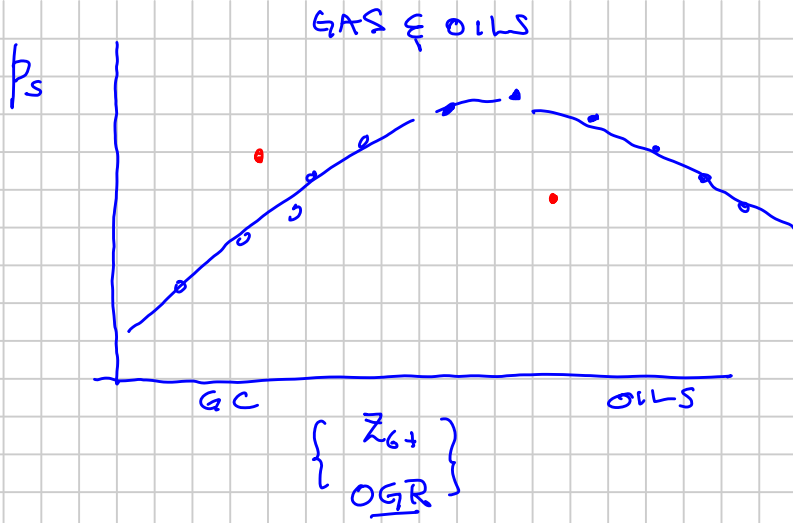
* PVT Data QC (Excel) -

- Individual Samples : mat. balance, plotting
- Trend plots (define trends, outliers) Multiple PVT Studies



TBP & Cnt





* PUT Data Entry to PUT Program (QC)
 ↳ Simple plotting

* Prediction EOS Calculation (using Default Characterization)

- Expect "Reasonable" match w/ PUT data

⇒ • $p_s \pm 2-15\%$ $= a + b \ln(i)$

⇒ • $\left. \begin{matrix} \rho_o \pm 1-3\% \\ Z_g \end{matrix} \right\} > 2-3\% \left(\gamma_i = f(\mu_i, C_{\underline{f}}) \right)$

Fix this now

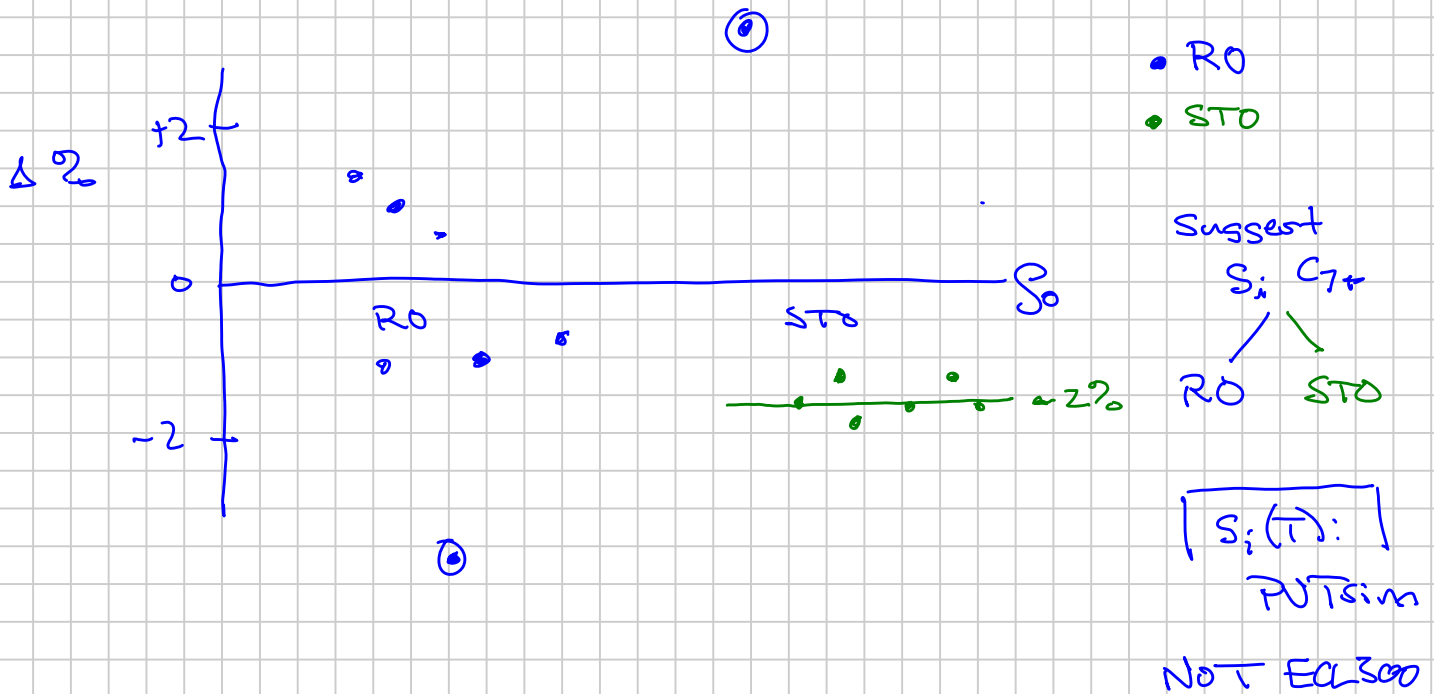
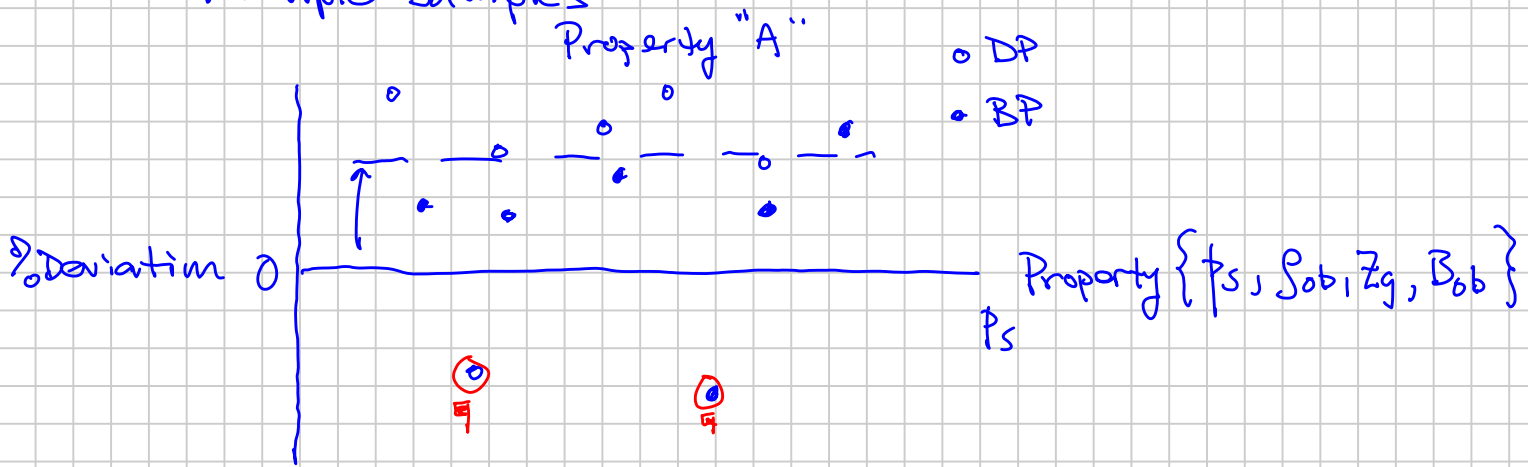
⇒ • SEP: $B_{ob}(\text{oils}) \text{GOR/ORR} \pm 2-5 (B_o) \pm 5-1x\% \text{GOR}$

• V_{ro}^{CVD} ZLE
 GC. $B_{od} \sim V_o$ oil shrinkage
Volatile oils

• $y_{j,t}^{CVD}$ (p) ! Gas Cond.

* Assessment of the Mismatch

- Multiple Samples



$$c_i = a_0 + a_1 T$$

$$s_i = \frac{c_i}{b_i}$$

$$c_i(60^\circ F) \quad 22 \quad b_i = R_0 \frac{R^2 T_0^2}{P_0^2}$$

$$c_i(280^\circ F) \quad 22$$

cc/g-mole

* Act on the Assessment (P_s, P_0 become unbraced)

- P_0 braced : fix the $\gamma_i (M_i, C_i)$

- p_s biased: Change (manually)

unbiased



$k_{C_1-C_7+}$ default
BIP
HC-HC

C_1-C_7 0.02
 C_1-C_{30+} 0.08 } PR

Casep / PVTsim
Dewpoint

$k_{HC-HC}^{SER} = 0$
default

$$\left\{ \begin{array}{l} p_d = f(K_{n+}) \\ p_b = f(K_{C_1}) \end{array} \right. \updownarrow$$

$$\frac{k_{C_1-C_7+} \neq 0}{1} < 0.2$$

$K_{C_1} \uparrow \uparrow p_b$
 $K_{C_7+} \downarrow \uparrow p_d$

Prausnitz-Chueh

$$k_{ij} = A \left[1 - \left(\frac{2v_{ci}^{1/6} v_{cj}^{1/6}}{v_{ci}^{1/3} + v_{cj}^{1/3}} \right)^B \right], \dots (5.79)$$

$A \sim 0.18$ $B = 6$
 $A = 1$ $B = 1$

$$v_{ci} = \frac{RT_{ci} z_{ci}}{p_{ci}} \quad 0.000x \quad 0.000x \quad \rightarrow \quad 0.0x$$

Many PVT Reported Quantities are Relative to p_s

Reference State

* OK slightly modified default EOS $p_0 \checkmark p_s$ unbiased

* Choosing Variables to Tune With (Pool of Variables)

(1) - C_{7+} (Table 1) Properties $\{T_{ci} \quad p_{ci} \quad \omega_i\} = \Theta_i$

- More uncertainty in Θ_i as SCN increases

NOT S_i : $S_i \Leftrightarrow \delta_i(M_i, C_i)$
unchanged

- Best if changes were "similar" for all the C_{7+} but more change for heavier

(2) - C_{N+} (25+ 30+ 35+) could be quite different in chemical makeup and most uncertain Θ_{N+}

(3) - k_{ij} all HC-HCs

- $C_1 - C_{7+}$ biggest effect on f_s
- Entire matrix Chueh-Prausnitzz (single variable) "A"

- $C_{7+} - C_{7+}$ huge impact on high p $K_i(p)$ "nose" impact retrograde condensation, CP, near-critical

$\sqrt{(1)-(3)}$ 10s - 100s will impact all samples in a similar, consistent way | Max. 5-10 EOS Variables

Sample-Dependent Variables \Rightarrow mol-% change in C_1 | $C_{7+} \lesssim 2$ mol-%

- M_{N+} : used to convert $w_i \rightarrow z_i$

$$z_i = \frac{w_i / M_i}{\left(\sum_{j=1}^{n-1} w_j / M_j \right) + \frac{w_{N+}}{M_{N+}}}$$

(2) should \sim all PVT calculations

- w_{N+} z_{N+} always uncertain

Could be very different sample to sample

- Flash-GC-Recomb
"GOR"

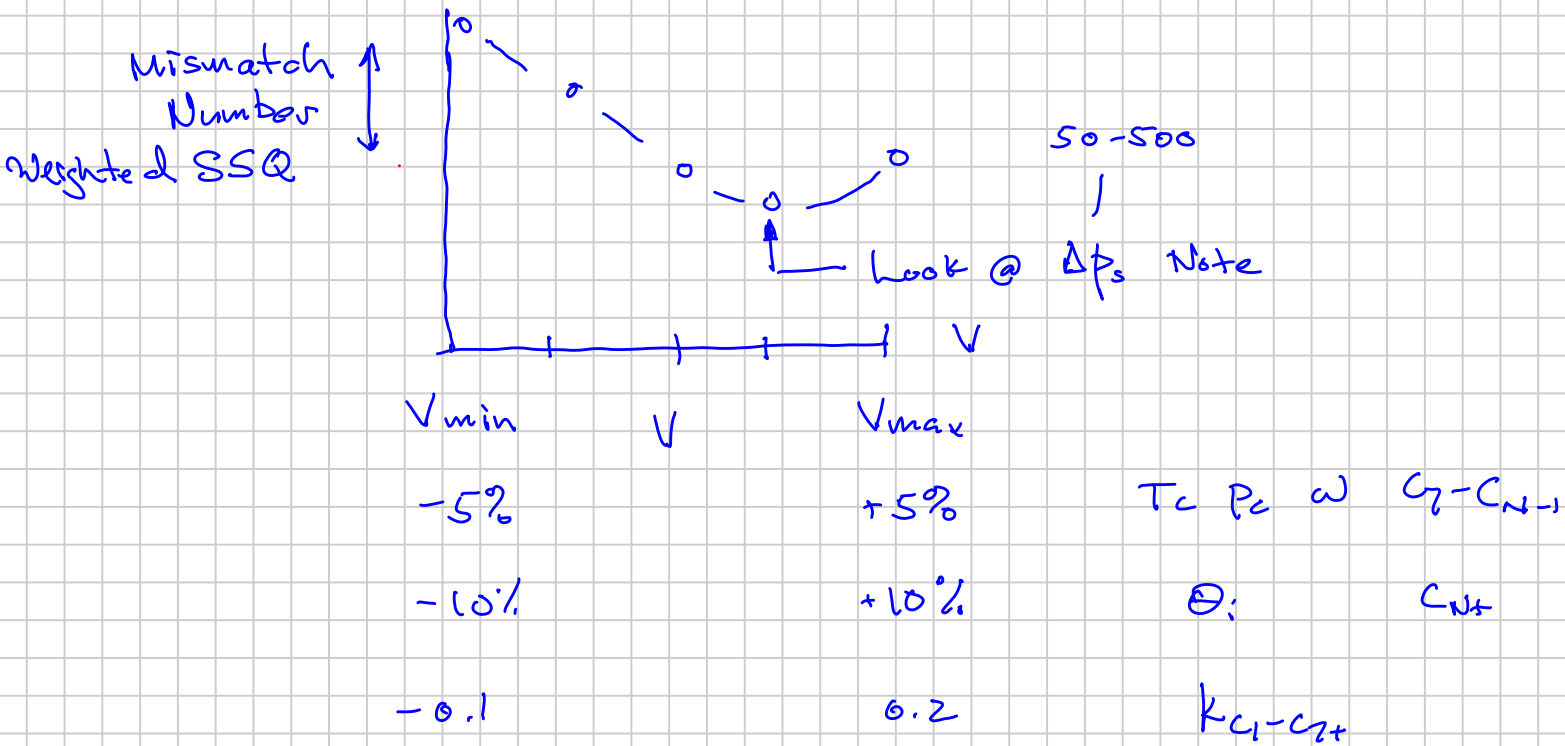
(Which PVT programs allow)
PVTx | PhaseComp

* Select to Weigh some data $\uparrow \downarrow$

- Importance of data w/ how many of that kind of data
- Experimental uncertainty $\Delta p_s \pm 1-5 \text{ psi}$

* Start systematic Tuning

- Try '1D' (single variable) sensitivities



- Adjust roots on key-data weight factors (p_s)

- Multi-variable automated regressions

- Look at all calculations
Plots of Tables
each one
- Trying different combinations of V_j
 - Adjusting a bit some weight factors
 - keep a Journal
 - Quit
- ~ days
~ weeks

Residual Weighting Within PhazeComp

Aaron A. Zick, Ph.D.
Zick Technologies

January 22, 2007

$$\chi^2 = \sum_{i=1}^N (w_i r_i)^2,$$

↙ you control

w_i = weight factor

r_i = residual

min \hat{r} (circled) = $\sqrt{\frac{\chi^2}{\sum_{i=1}^N w_i^2}}$

RMS

0

1-3%

Reference: $r_i = \Delta_i$ use $y^r = y^e$

EOS LAB

$$r_i = 100(y_i^c - y_i^e) / y_i^r,$$

$\frac{0.01}{1} = 1\%$

= max of data of a given type in a given experiment



$$\frac{0.01}{0.03} = 0.33$$

33%

min WSSQ by changing V_j (boxed) EOS Sample-Specific

① Which data are poorly predicted?

x - $R_{sd,b}$

x - $B_{od}(p)$

Plot $R_{sd}(p)$

$B_{od}(p)$

$B_{od}(R_{sd})$

② Which data are most important?

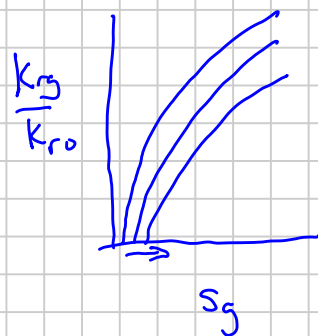
- Depletion SGD: Low vs high R_{Fo} ?

$$\frac{V_o(p) \Delta R_s}{B_{od}(p): R_{sd}(p)}$$

$$\frac{k_{rg}}{k_{ro}}(S_g)$$

③ Any missing data?

$$a = \rho_a \frac{R^2 T_c^2}{p_c} \cdot \alpha$$

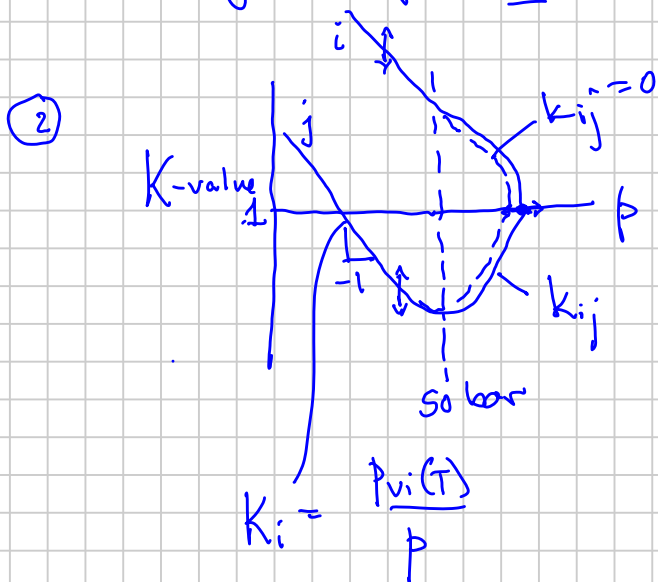


$$\alpha_{RK} = 1/\sqrt{T_r}$$

$$\alpha_{SRK} = f(T_r, \omega)$$

BIPs $[k_{ij}]$ Are correction terms

① Mainly change $\underline{k_i}$ and $\underline{k_j}$ in the $p \approx 50$ bar



All components ^{k-values} are affected also, because joined @ P_k

TABLE 6.7—SEPARATOR TESTS (RESERVOIR-FLUID) OF GOOD OIL CO. WELL 4 OIL SAMPLE

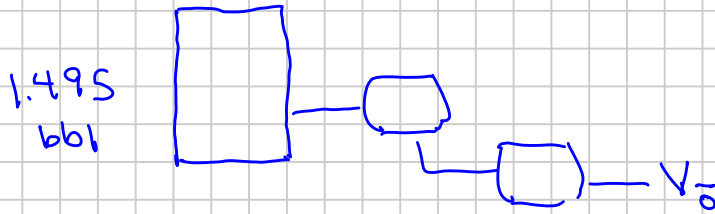
Separator Pressure (psia)	Separator Temperature (°F)	GOR ^b (ft ³ /bbl)	GOR ^c (ft ³ /bbl)	Stock-Tank Gravity (°API)	FVF ^d (bbl/bbl)	Separator Volume Factor ^e (bbl/bbl)	Flashed-Gas Specific Gravity
50 to 0	75	715	737			1.031	0.840
	75	41	41	40.5	1.481	1.007	1.338
100 to 0	75	637	676			1.062	0.786
	75	91	92	40.7	1.474	1.007	1.363
200 to 0	75	542	602			1.112	0.732
	75	177	178	40.4	1.483	1.007	1.329
300 to 0	75	478	549			1.148	0.704
	75	245	246	40.1	1.495	1.007	1.286

^aGauge.
^bIn cubic feet of gas at 60°F and 14.65 psi absolute per barrel of oil at indicated pressure and temperature.
^cIn cubic feet of gas at 60°F and 14.65 psi absolute per barrel of stock-tank oil at 60°F.
^dIn barrels of saturated oil at 2,620 psi gauge and 220°F per barrel of stock-tank oil at 60°F.
^eIn barrels of oil at indicated pressure and temperature per barrel of stock-tank oil at 60°F.

MIX FEED OIL
 PRES 2620 PSIG
 TEMP 220 F
 DLE ID "Oil separator test: Psp1=300 psig."
 BASIS 1.495 BBL

TEMP F	PRES PSIG	MREM SCF	LVOL BBL	LDEN G/CC	GSG	LAPI
220	2620	0	1.495	0.6562		
75	300	549	1.148		0.704	
75	0	795	1.007		1.286	
60	0		1			40.1

End; DLE (separator test)



$$B_{ob}^{Eos} = \frac{1.495}{0.9981} \quad \text{vs}$$

$$B_{ob}^{Lab} = \frac{1.495}{1}$$

Pool of Variables:

$$* k_{C_1-C_7+} \text{ or } k_{\text{H}_2\text{O}} \quad [1]$$

$$* C_{30+} \quad T_{C_i} \quad P_{C_i} \quad \omega_i \quad [3]$$

$$* M_{7+} \quad (\omega_i \rightarrow z_i) \quad [1]$$

$$* z_{30+} \quad [1]$$

$$* GOR_{FGCR} \quad [1]$$

} Composition Variables

$$\left\{ * C_{7-29} \quad T_{C_i} \quad P_{C_i} \quad \omega_i \quad [3] \right\}$$

Ph2: Introduce a Variable

VARIABLE v-name v-start { v-min v-max }

Apply the variable v-names

MULTIPLY
DIVIDE
INCREASE
DECREASE
REPLACE



BIPS-CI-C7+-MULTIPLIER

0
0.5
1
> 1.01
1.5
2

RMS %
11.62
7.585
5.075
8.332
15.42

Pb [2620] Lab
psig
2241
2414
2615
2620
2852
3139

Next Try: D = default

k_{rec}	k_{HC-HC}	T_{30+}	p_{30+}	w_{30+}	M_{7+}	"Flash-66- Recomb"	GOR	Final RMS %	Pb psig [2620]
1	(1.01)							5.075	2620 [767]
2	(0.866)							5.864	2591
3	1	-	(1)					5.075	2620
4	1	-	D	(1)				5.075	2620
5	1	-	D	D	(1)			5.075	2620
<hr/>									
6	(1.47)	(1.036)						5.03	2620
7	(1.41)	D	(0.9)					5.04	2620
8	(1)	D	D	(1.004)				5.075	2620

9 (1.55) (1) (0.9) (1) 4-variable 5.036 2618

10 (2) (1.041) (1.05) (1) w/ C7+ not C30+ 4.405 2612
749

Replace BIPS of N2 with i-C4 to C30+ by 0.08
Replace BIPS of CO2 with i-C4 to C30+ by 0.15

CORRELATE

```
; VARIABLE TC-C30+ 1 0.9 1.1
; MULTIPLY TC OF C30+ BY TC-C30+
; VARIABLE PC-C30+ 1 0.9 1.1
; MULTIPLY PC OF C30+ BY PC-C30+
; VARIABLE AF-C30+ 1 0.9 1.1
; MULTIPLY AF OF C30+ BY AF-C30+
```

```
VARIABLE BIP-C1-C7+-MULTIPLIER 1 0 2
CHUEH C1 with C7 to C30+ by 1
MULTIPLY BIPS C1 WITH C7 TO C30+ BY BIP-C1-C7+-MULTIPLIER
```

```
; REG2
; VARIABLE BIP-HC-HC-MULTIPLIER 0 0 2
; CHUEH C1 TO C30+ with C1 to C30+ by 1
; MULTIPLY BIPS C1 TO C30+ WITH C1 TO C30+ BY BIP-HC-HC-MULTIPLIER
```

```
VARIABLE TC-C7+ 1 0.95 1.05
MULTIPLY TC OF C7 TO C30+ BY TC-C7+
VARIABLE PC-C7+ 1 0.95 1.05
MULTIPLY PC OF C7 TO C30+ BY PC-C7+
VARIABLE AF-C7+ 1 0.95 1.05
MULTIPLY AF OF C7 TO C30+ BY AF-C7+
```

```
VARIABLE MWAvg-C7+ 1 0.9 1.1
MULTIPLY GAMMA AVERAGE BY MWAvg-C7+
```

$$M_{7+} \times 0.9$$

$$K_{ij} \times 1.29$$

$$C_f \quad \underline{0.296}$$

$$WT \quad API \quad 10$$

$$RMS \quad \underline{3.794}$$

$$P_b = 2618 \checkmark$$

$$P = 159$$

$$R_{sd} = \underline{762}$$

$$B_{od} = \underline{1.1965}$$

FGCR GOR \equiv EQV

EQV

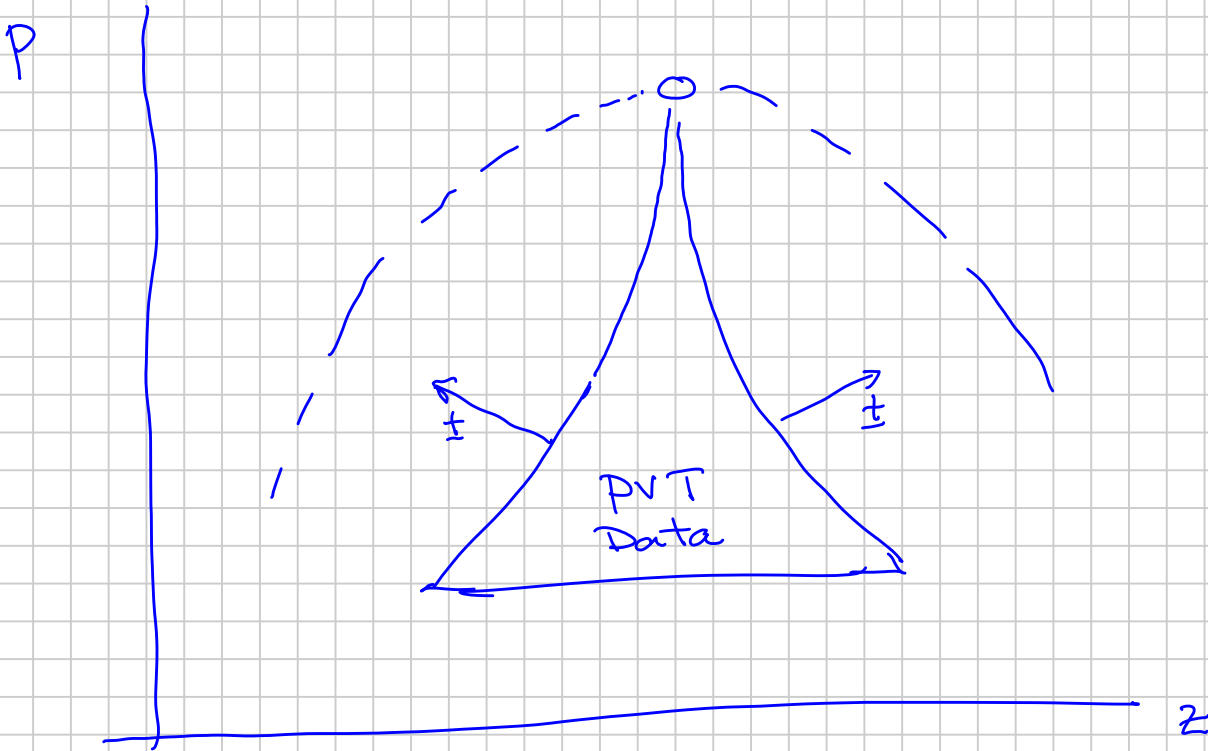
$$Z_i = \underbrace{f_g y_i}_{\text{FLASHED-GAS}} + (1-f_g) x_i$$

Var Num	Reg Var	Variable	Current	Initial	Lower Bound	Upper Bound
1		CF	✓ 2.90000e-01			
2	1	TC-C30+	✓ 9.95453e-01	1.00000e+00	9.00000e-01	1.10000e+00
3	2	BIP-C1-C7+-MULTIPLIER	✓ 1.12834e+00	1.00000e+00	0.00000e+00	2.00000e+00
4	3	FG	5.80748e-01	5.00000e-01	0.00000e+00	1.00000e+00

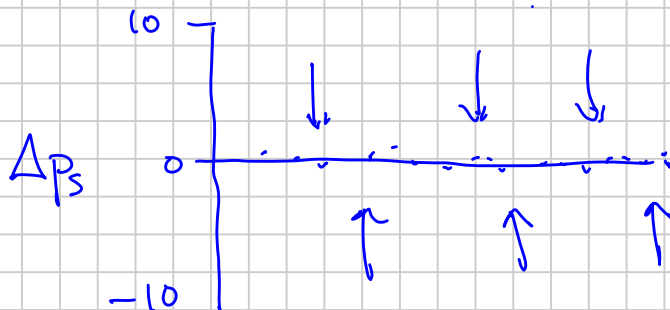
$RMS = 4.05$

$R_{sd} = 736$ $B_{od} = 1.21$

ρ_o ✓



{ Smaller EOS Property Changes
 +
 ⊗ Smaller by-sample composition changes



OIL VISCOSITY MODELING

Alani-Kenedy

① LBC (Lorentz Bray Clark)

"Jossi-Thodos"

oils $[v_{ci}(T) (M_{7+}, X_{7+})]$

Pure Compounds

② Pedersen

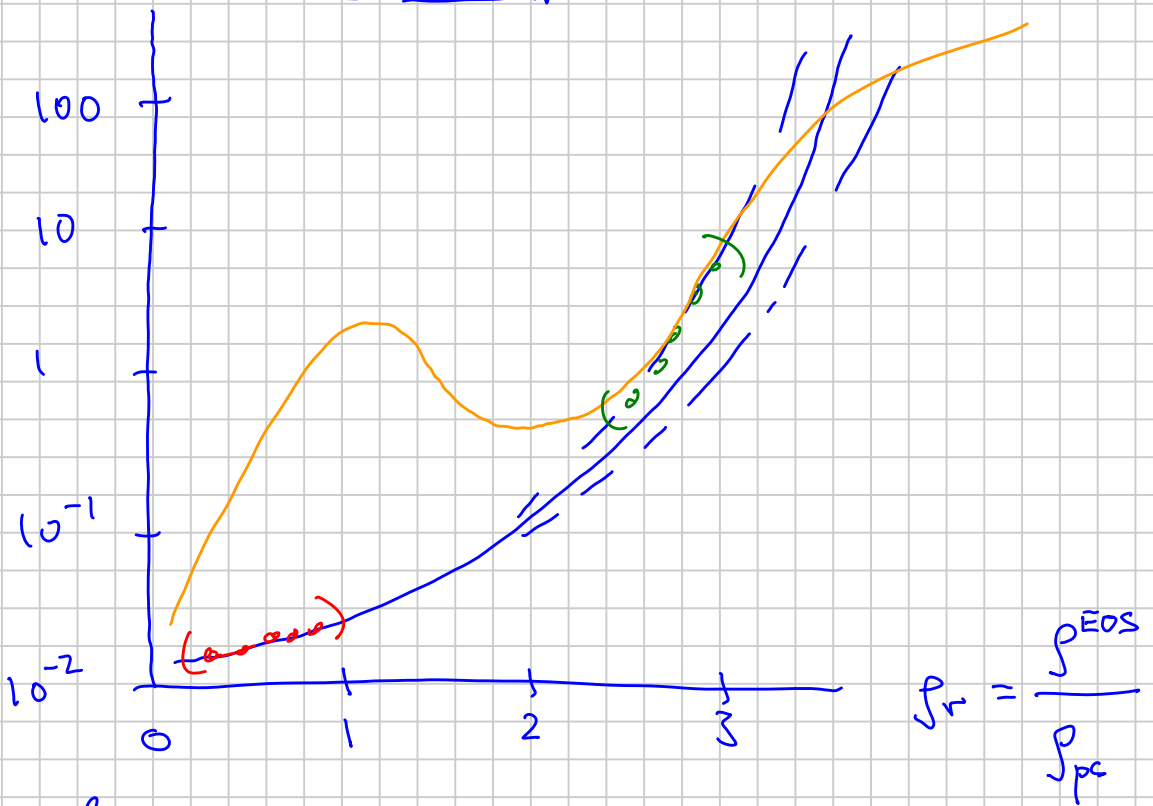
- Tuning: don't know, ask Calsep

$$p_r = \frac{p}{p_{pc}} \quad p_{pc} = \frac{M}{v_c}$$

LBC:

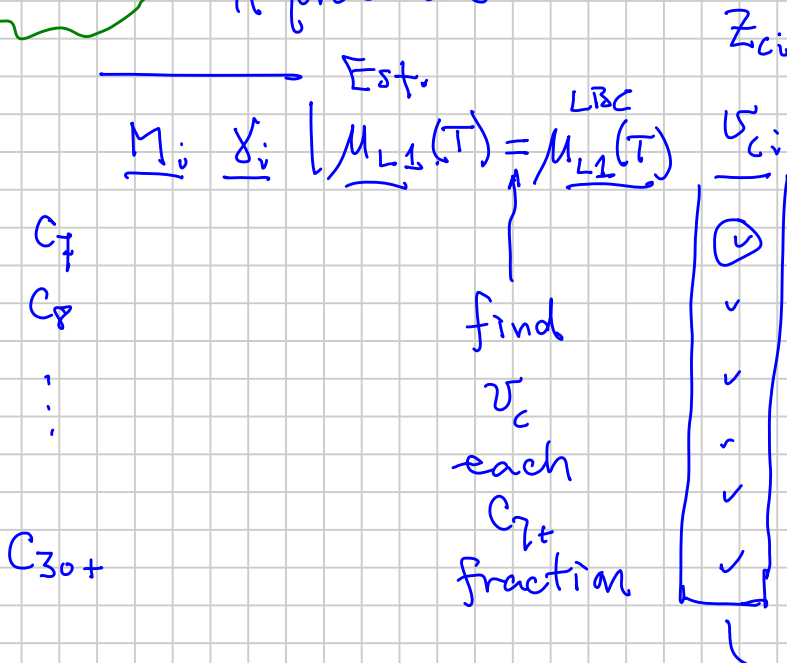
$$\mu - \mu^o = \left[a_0 + a_1 p_r + a_2 p_r^2 + a_3 p_r^3 + a_4 p_r^4 \right]^4$$

↑ ↑ ↑



Yang Fevang

A procedure

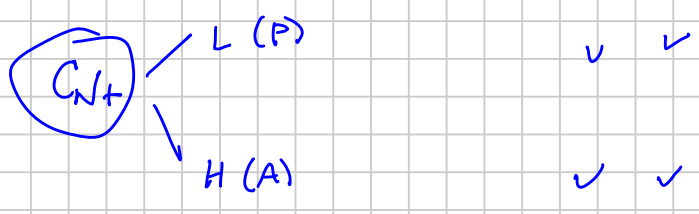
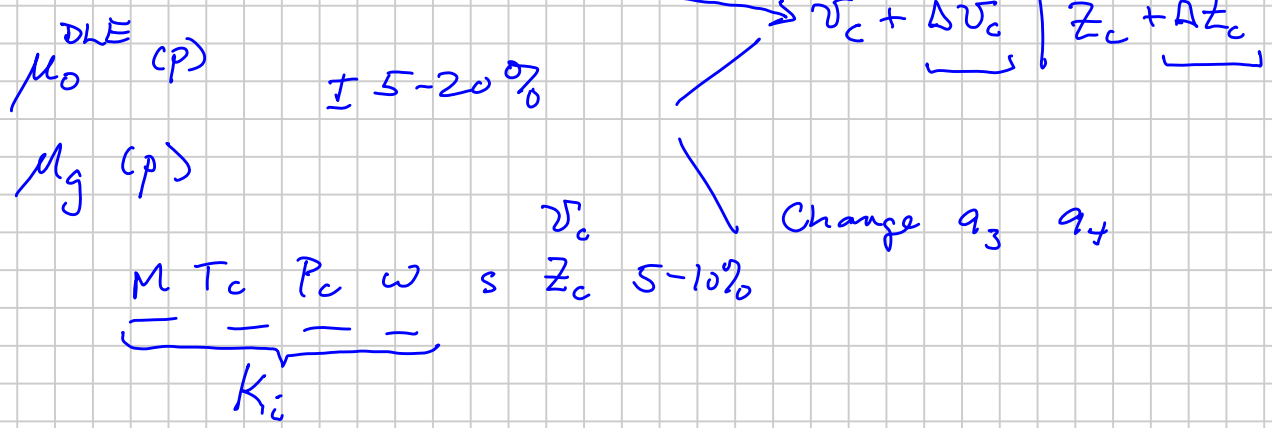


$$p_{pc} = \frac{p_{pc} M}{R T_{pc} Z_{pc}} = \frac{M}{v_{pc}}$$

$$v_{pc} = \sum u_i v_{ci}$$

↑
 $y_i = x_i$

KEY: v_{ci} C_{7+} fractions



API variations
 Troubles fitting
all oil μ_0

Black-Oil PVT

optimize

to minimize error

* BO Properties:



in the important
 $S_g(p,T)$ $S_o(p,T)$

OIL PHASE

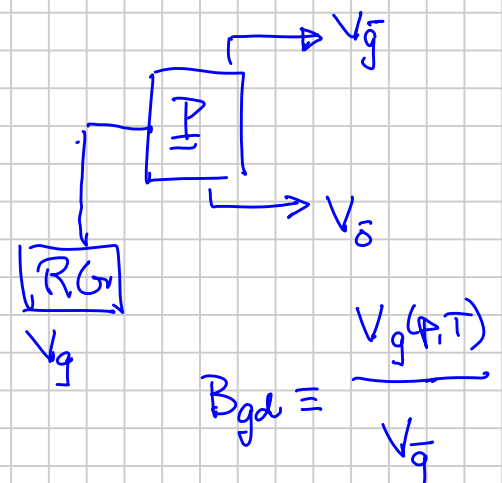
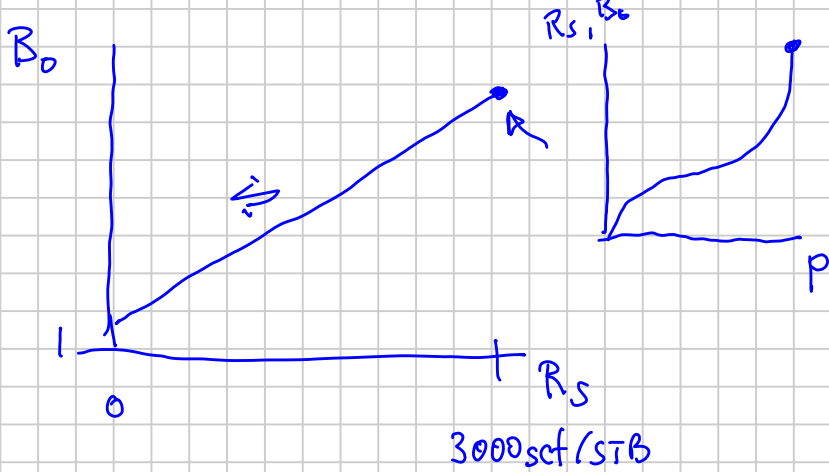
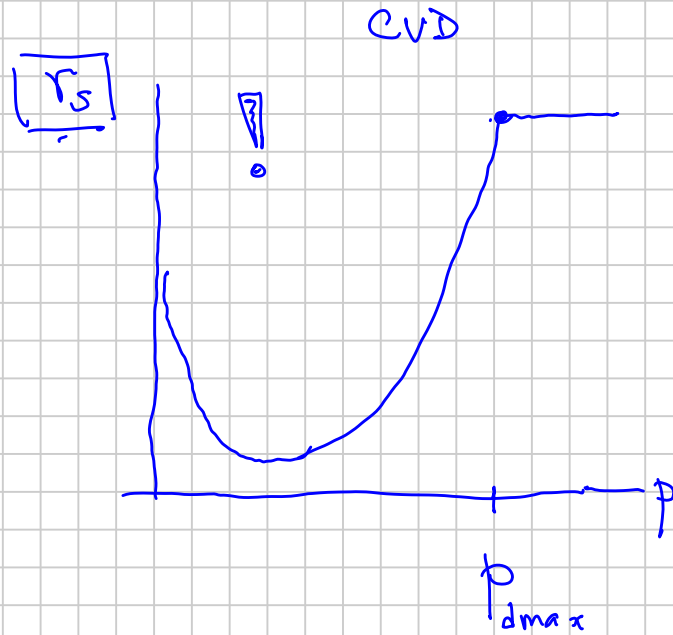
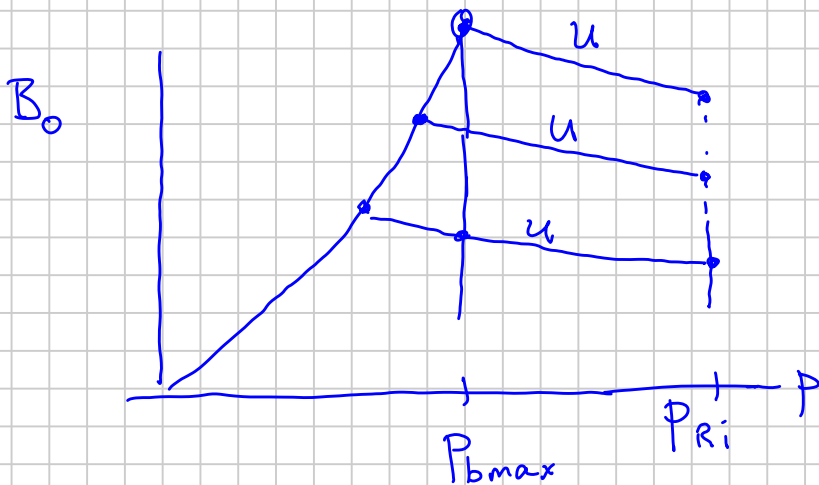
GAS PHASE

$B_o(p)$ $R_s(p)$ $\mu_o(p)$

$\frac{B_g(p)}{Z_g(p)}$ $\frac{\Gamma_s(p)}{R_v}$ $\frac{\mu_g(p)}{p}$

$S \sim \mu @ p$

$S \neq \mu @ p$



$$\rho_g(p,T) = \frac{\rho_{gg} + \Gamma_s(p,T) \rho_{og}}{B_{gd}(p,T)}$$

$$B_{gw} = \frac{P_{sc}}{T_{sc}} \cdot \frac{T_r Z_g^{sk}}{p}$$

$$\rho_o(p,T) = \frac{\rho_{oo} + R_s(p,T) \rho_{go}}{B_o(p,T)}$$

$$B_{gd} = B_{gw} \frac{1}{(1-y_{nt})}$$

$$C \cdot \frac{y_{nt}}{1-y_{nt}} = r_s$$

How to Build BOPT from an EOS (Ch. 7)

Required Input:

1. EOS (EOS-R | EOS-S | $s_i(T)$)

2. Z_{Boi} "pick the sample with the highest p_s "

SPE 63087

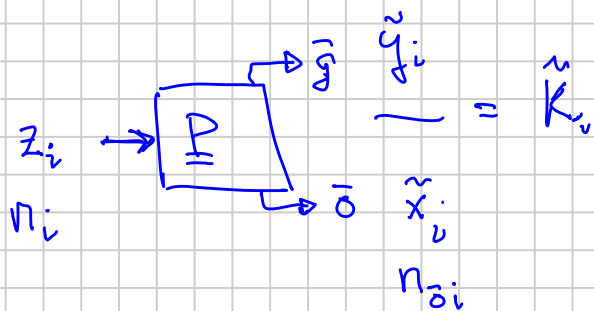
GOC (S/U) @ GOC if ands $\frac{1}{2}$ butts

3. Surface Process ($\rightarrow \bar{g} \bar{o}$)

✓ - EOS-based

{ - \tilde{K}_i (Statoil)

{ - RF_{oi} (Z_{nt})



$$RF_{oi} = GP_i = \frac{n_{oi}}{n_i}$$

"GP Table" = $f(Z_{nt})$

4. Method of Depletion (p_{smax} of $Z_{Boi} \rightarrow$ lower p 's)

CCE | CRD | DLE

5. Saturated List of $p < p_{smax}$

6. Undersaturated List of $p > p_{smax}$

{ 7. } Method saturated table extrapolation

$$p_s > p_s(\underline{Z_{Boi}})$$

- Gas injected
-
-

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pg

The surface process of Z_{Boi}

pg