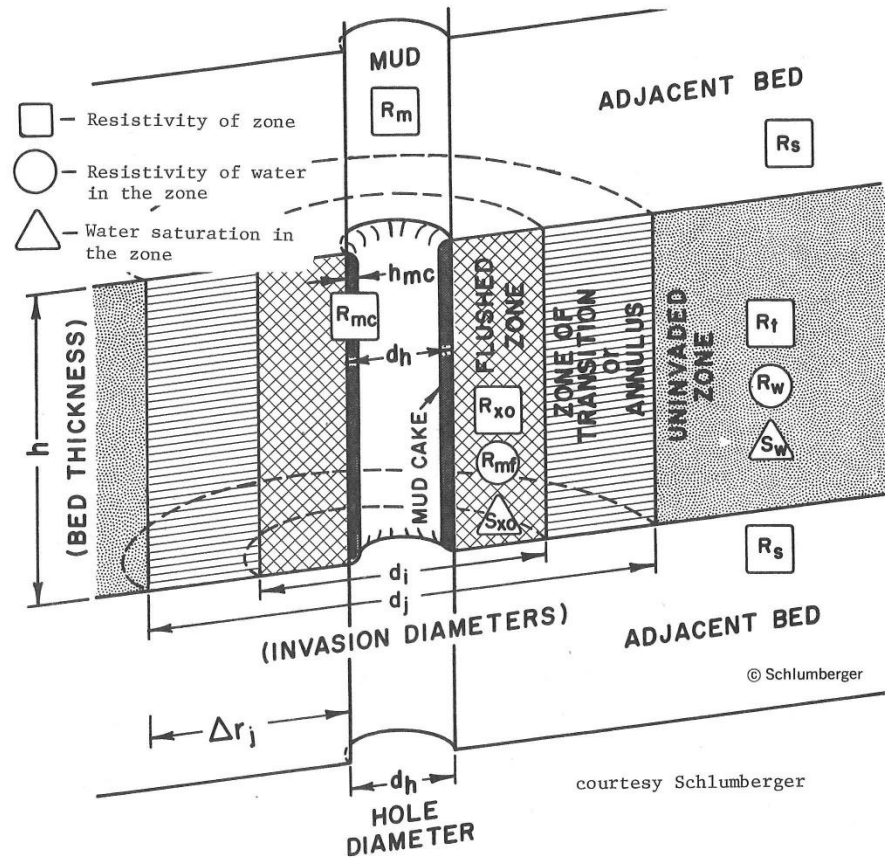


OPENHOLE LOG EVALUATION

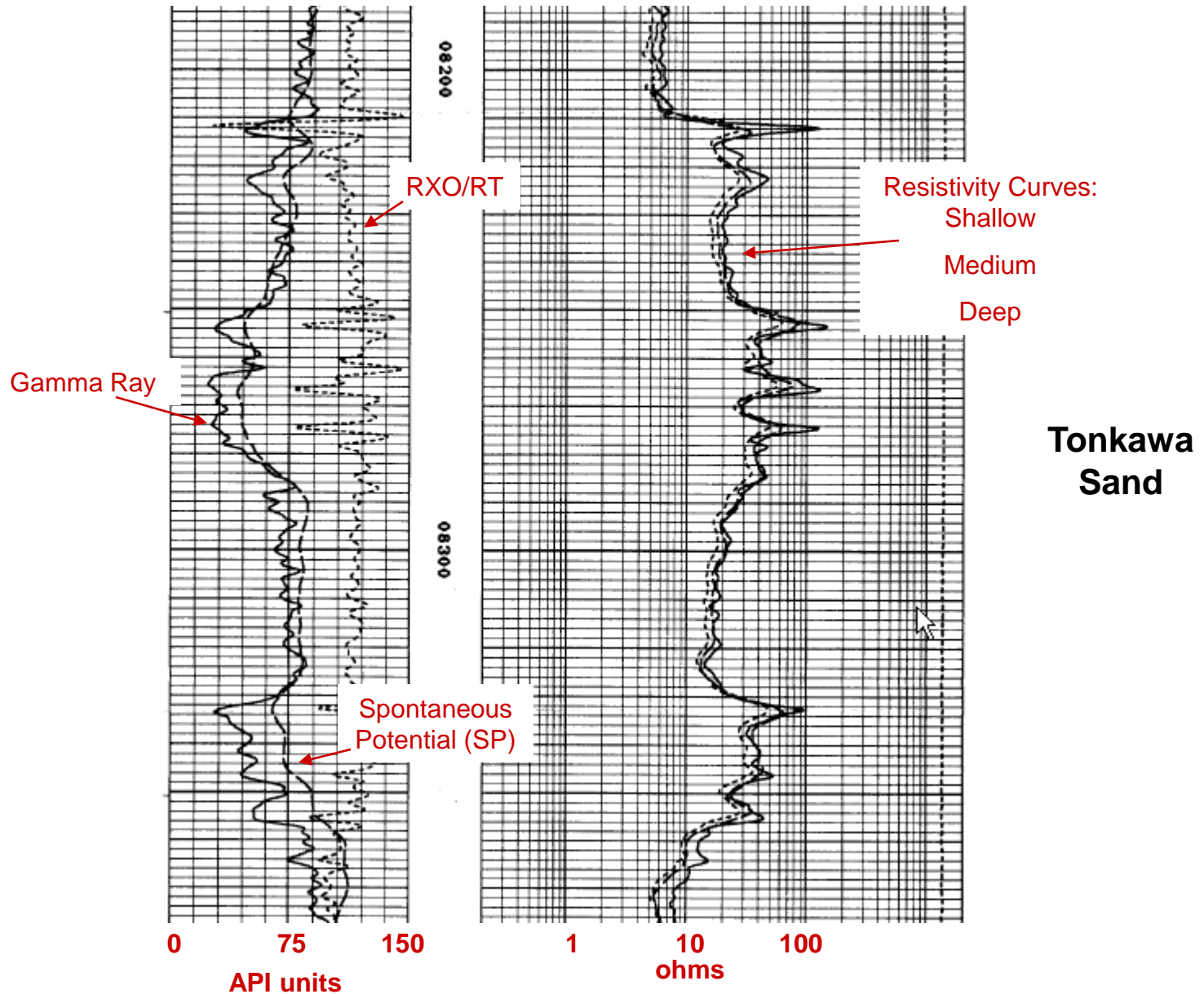
Prepared By
Bob Gaddis

Figure 2-8 Schematic Cross Section of a Well Bore



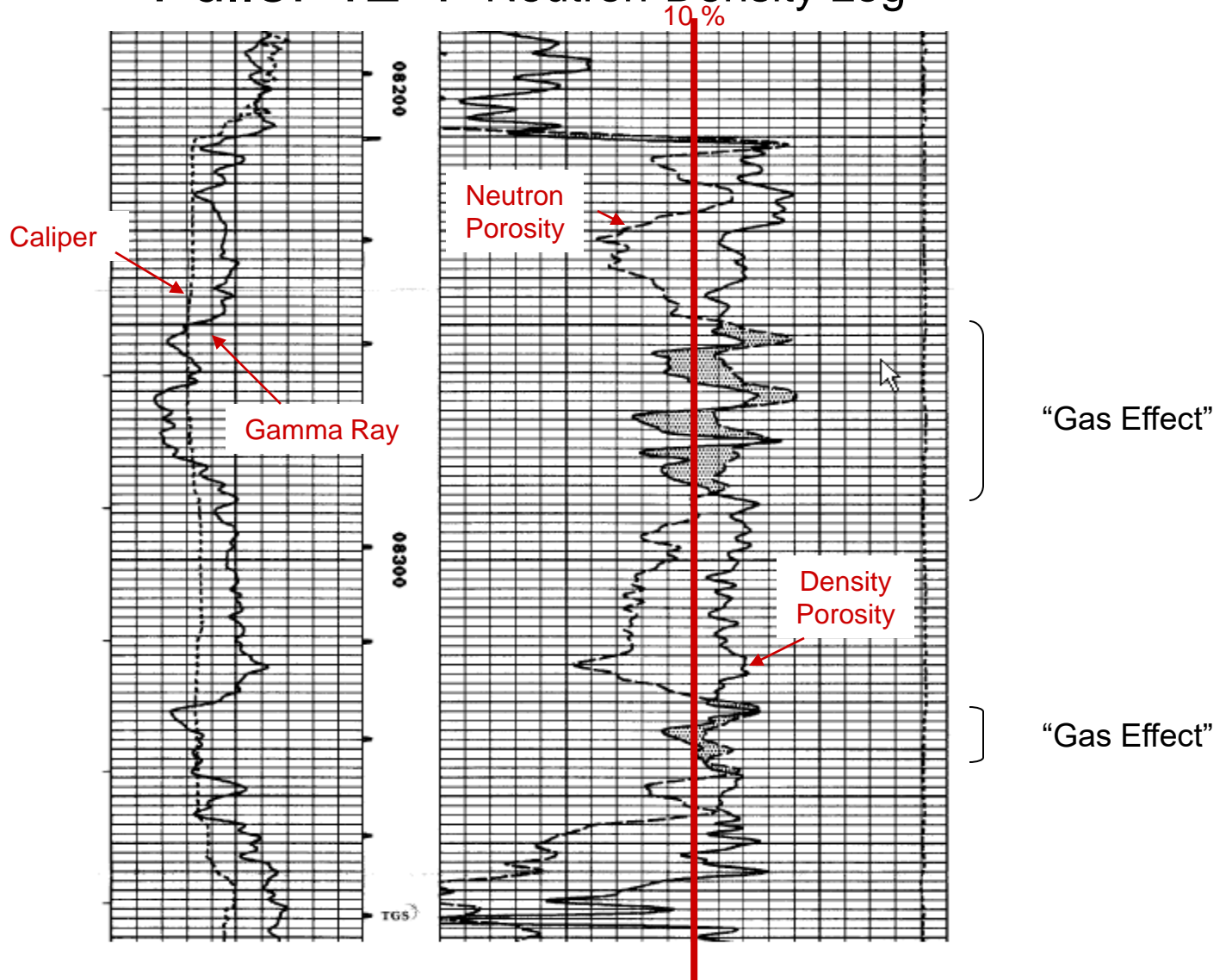
Parameter	Flushed Zone	Transition zone	Virgin Zone
resistivity	R_{xo}	R_i	R_t
porosity	ϕ	ϕ	ϕ
water saturation	S_{xo}	S_i	S_w
water	R_{mf}	R_z	R_w
equations	$S_{xo} = \sqrt{\frac{F_R R_{mf}}{R_{xo}}}$	$S_i = \sqrt{\frac{F_R R_z}{R_i}}$	$S_w = \sqrt{\frac{F_R R_w}{R_t}}$
note: with invasion	$S_{xo} > S_i > S_w$	unless all are 100%	
no invasion	$S_{xo} = S_i = S_w$ and R_w in all pores.		

Fuller 12-1 Dual Induction Log

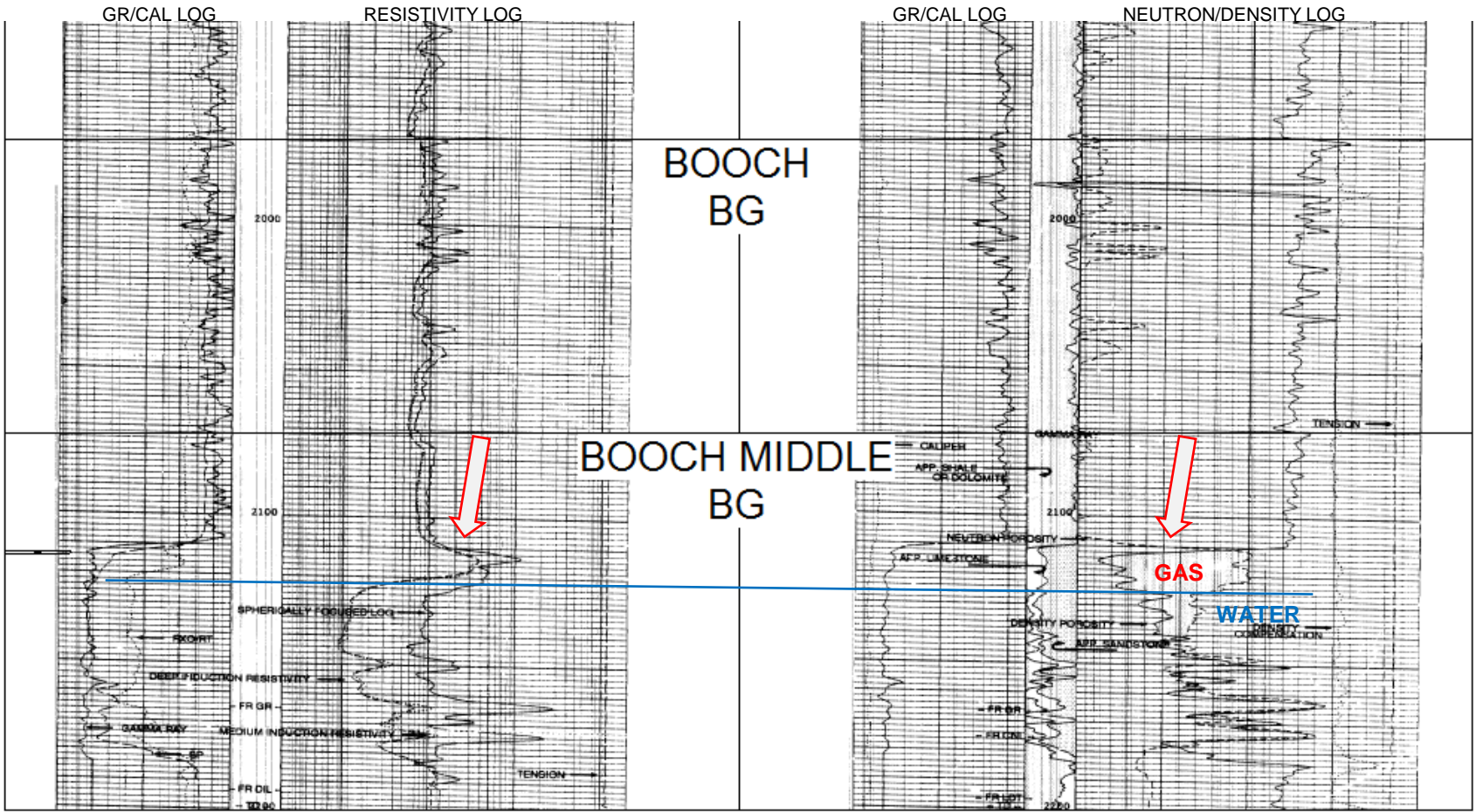


- Section 12-13N-19W, Custer Co., Okla.

Fuller 12-1 Neutron-Density Log

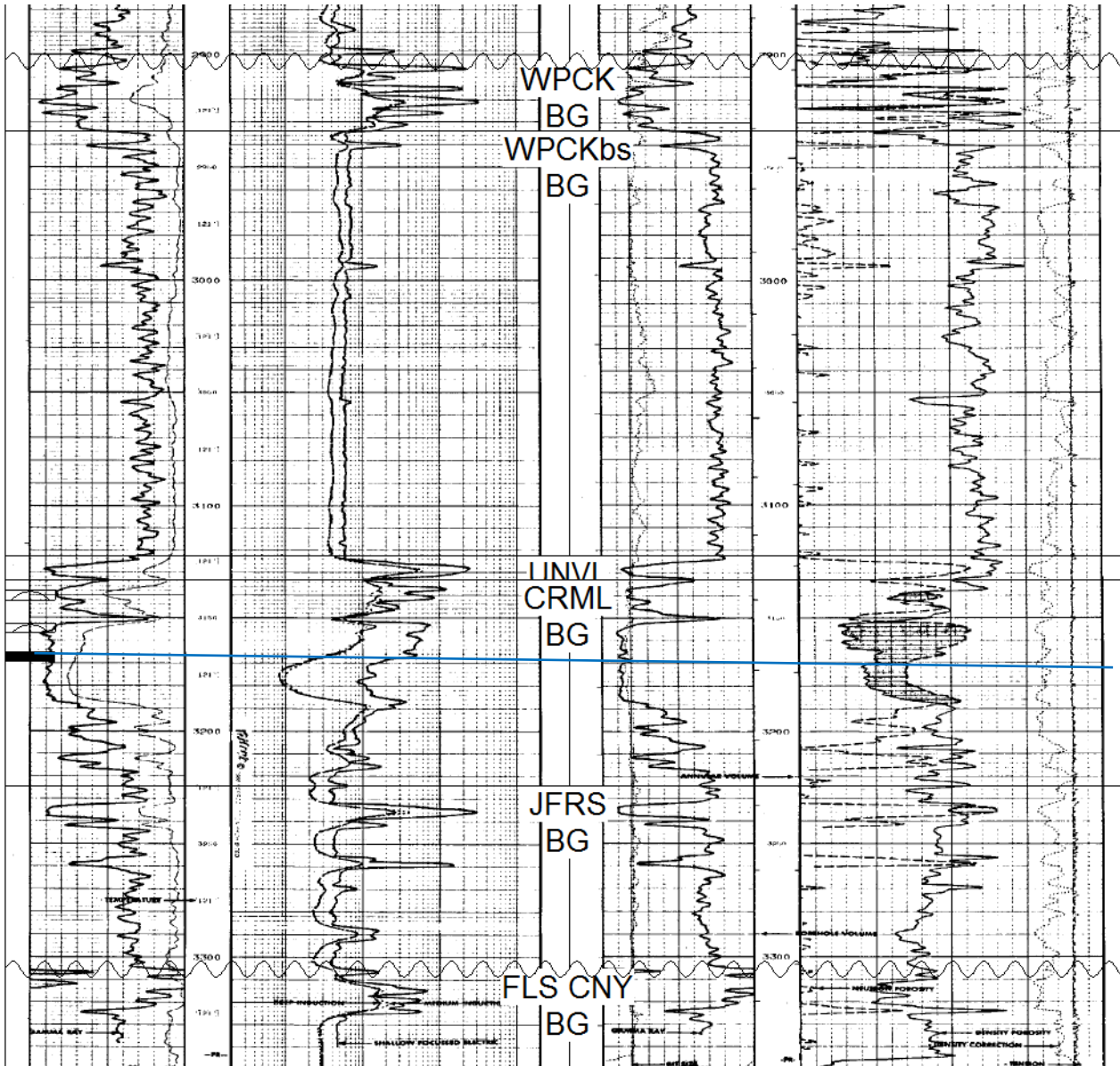


- Section 12-13N-19W, Custer Co., Okla.

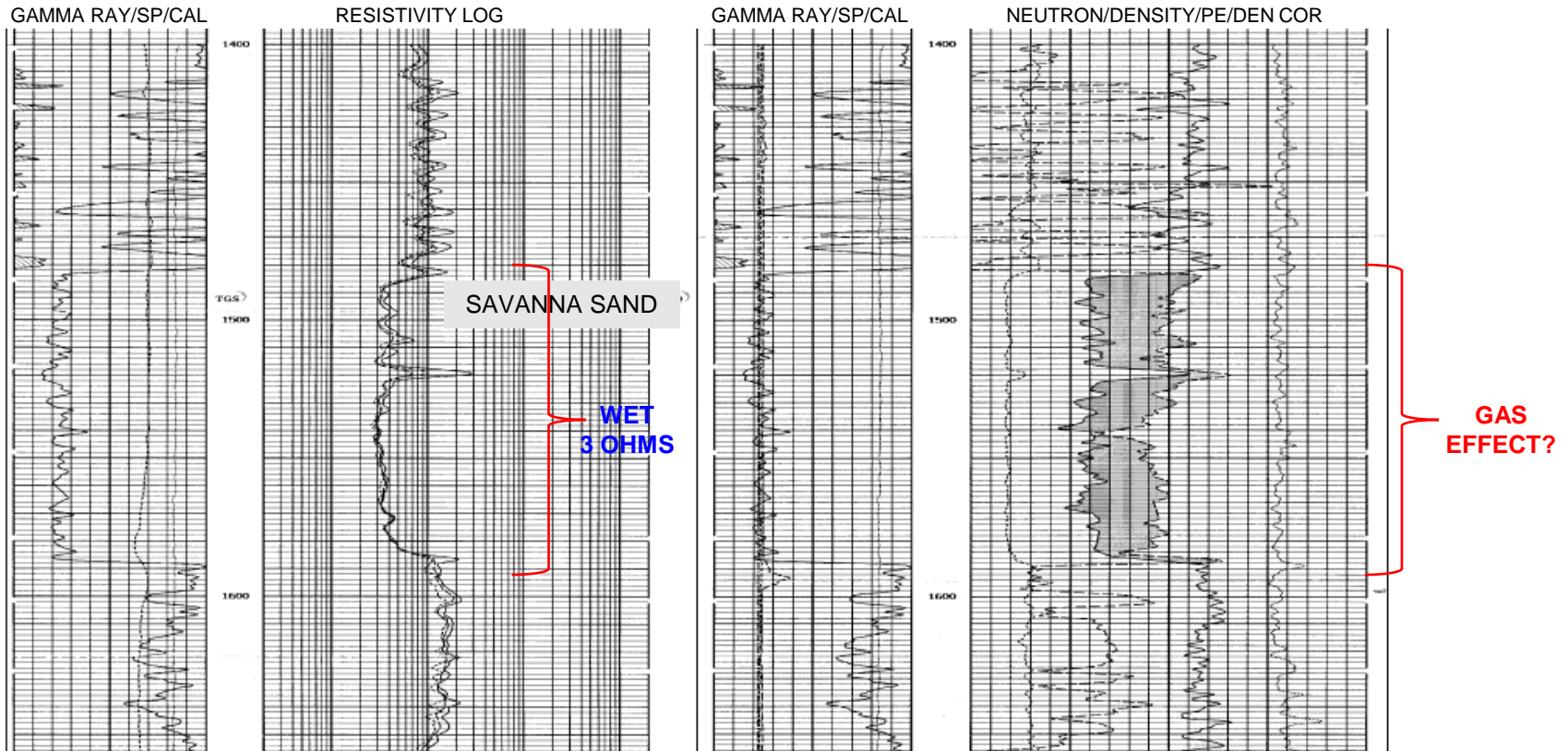


TD=2200.00

JORDAN O&G
ROSA 3-2
SE SE NE SEC. 2-9N-10E
HUGHES CO.
35063226050000



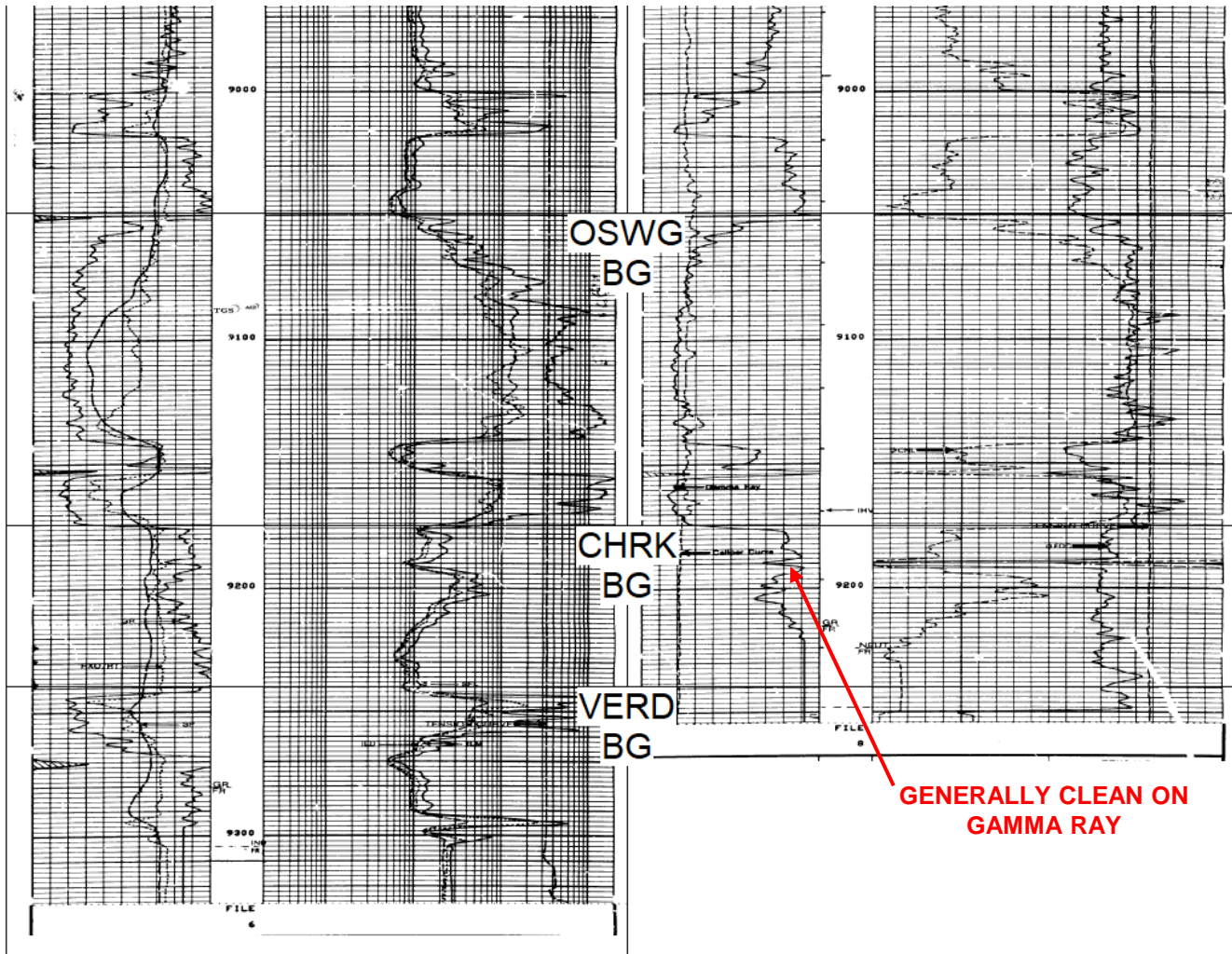
TD=3350.00



NO, NOT "GAS EFFECT"
BUT A CLEAN WATER SAND

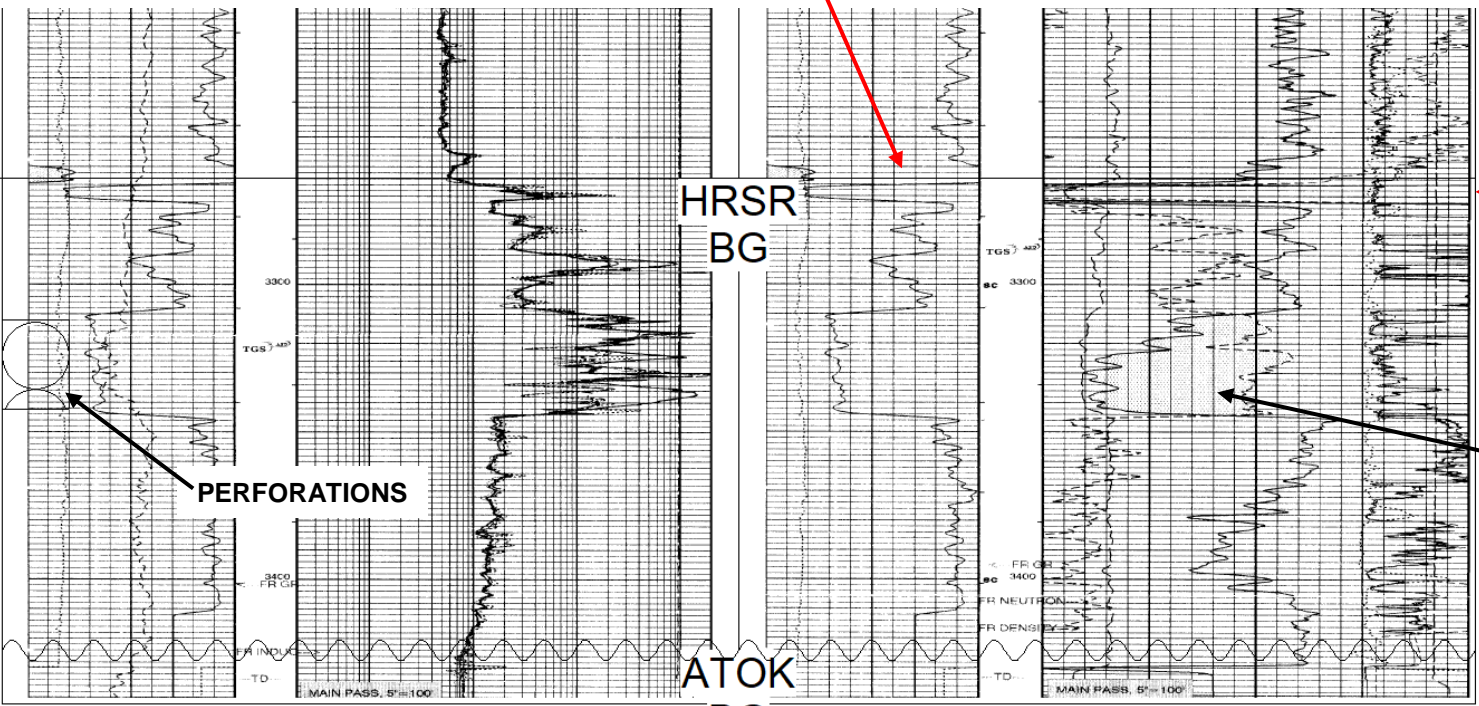
SO, THINK "CROSS-OVER", NOT "GAS EFFECT"

HPC
CLARK 1
SE NW SEC 29-13N-8W
CANADIAN CO.
3501722325000

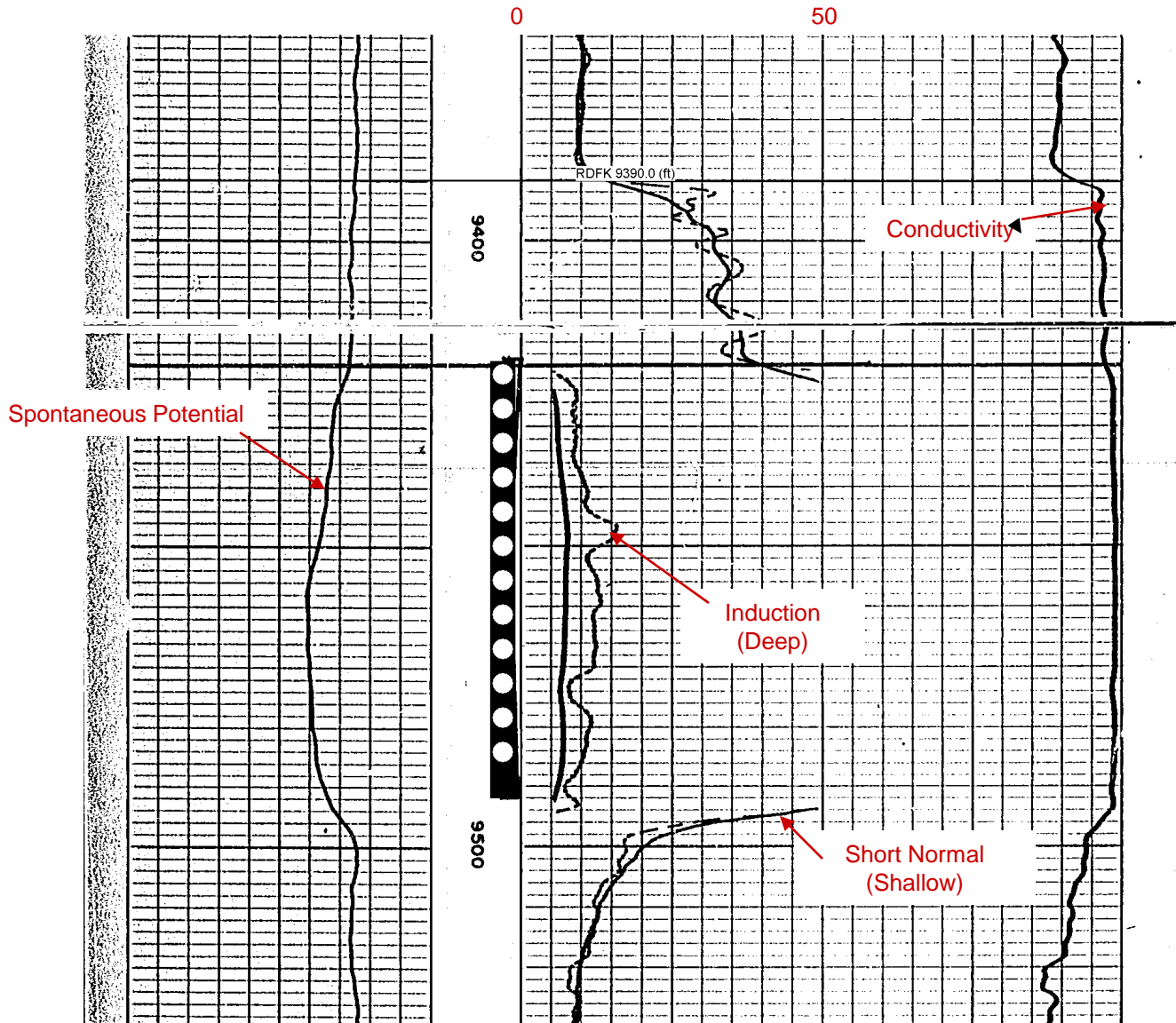


DAVIS OPER
DONNA 1
NE SEC 16-5N-12E
PITTSBURG CO.
35121224330000

GENERALLY CLEAN ON
GAMMA RAY

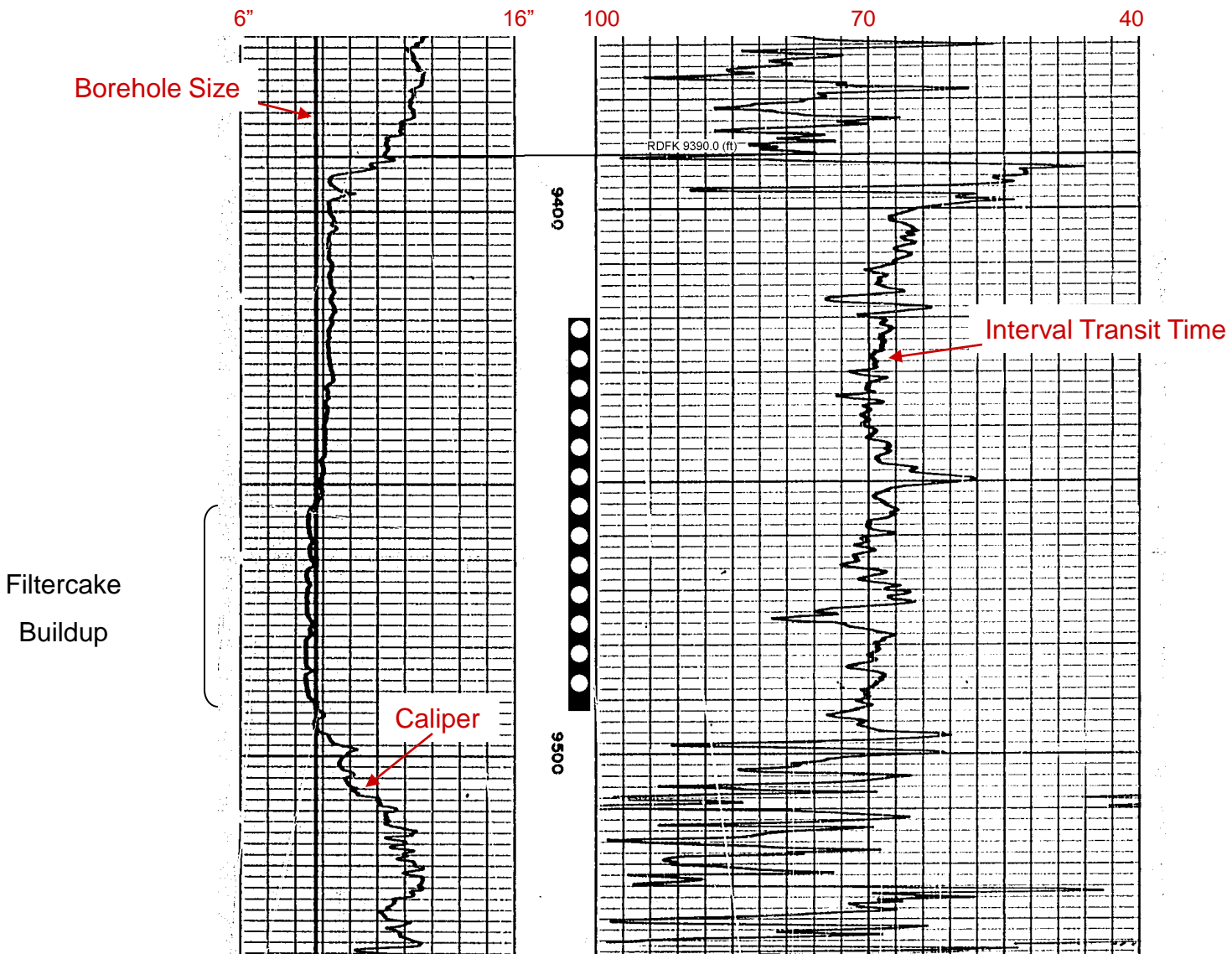


Fluitt 1 Induction Electric (IE) Log



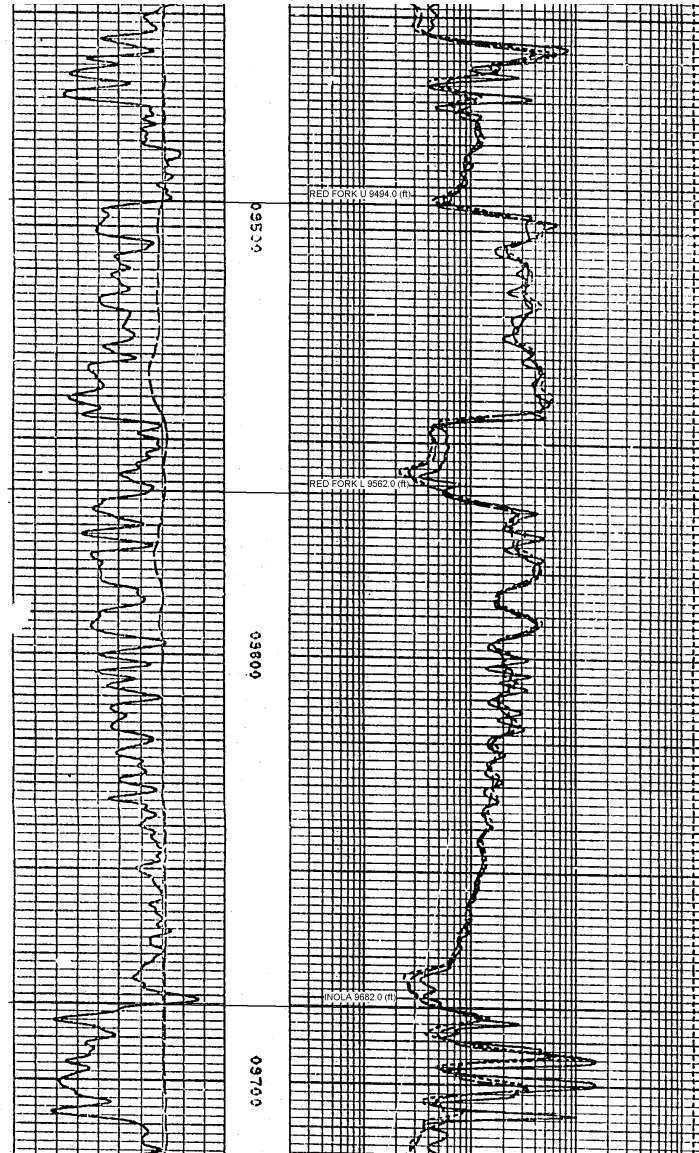
- Section 14-7N-4W, McClain Co., Okla.

Fluitt 1 Sonic Log



• Section 14-7N-4W, McClain Co., Okla.

Gambrel 1 Dual Induction Log



Pink Lime

Red Fork

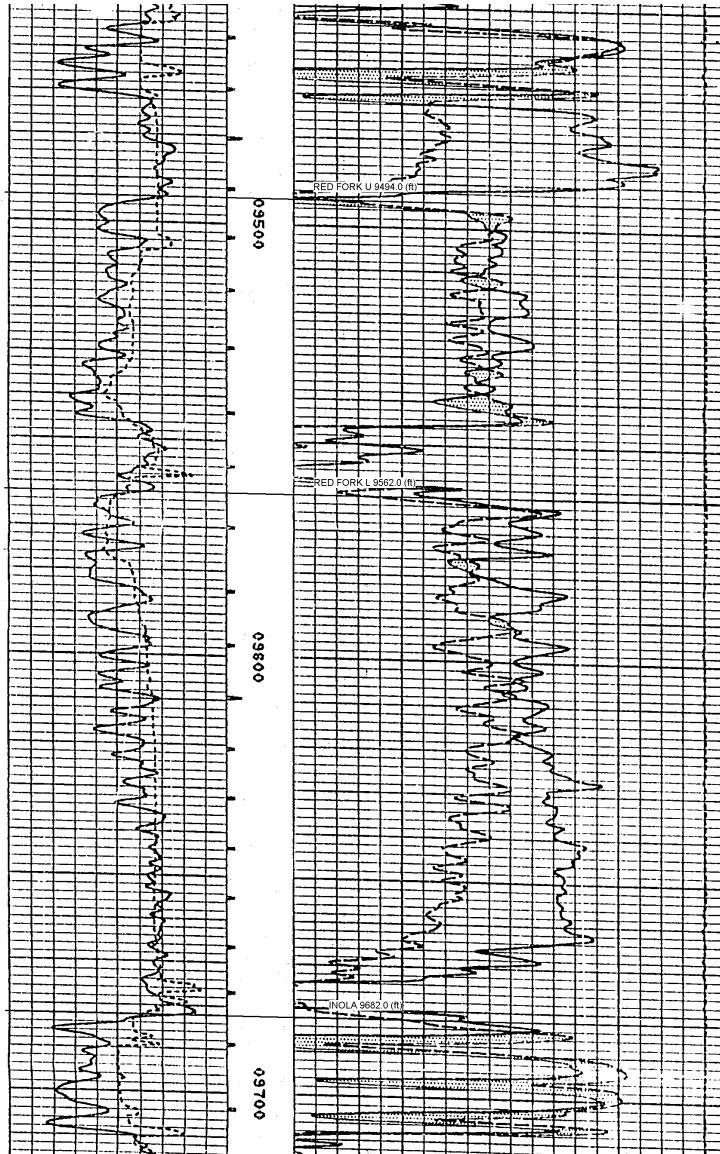
Upper

Lower

Inola Lime

- Section 13-17N-17W, Dewey Co., Okla.

Grambrel 1 Neutron/Density Log



- Section 13-17N-17W, Dewey Co., Okla.

Albers 1-34 Dual Induction Log

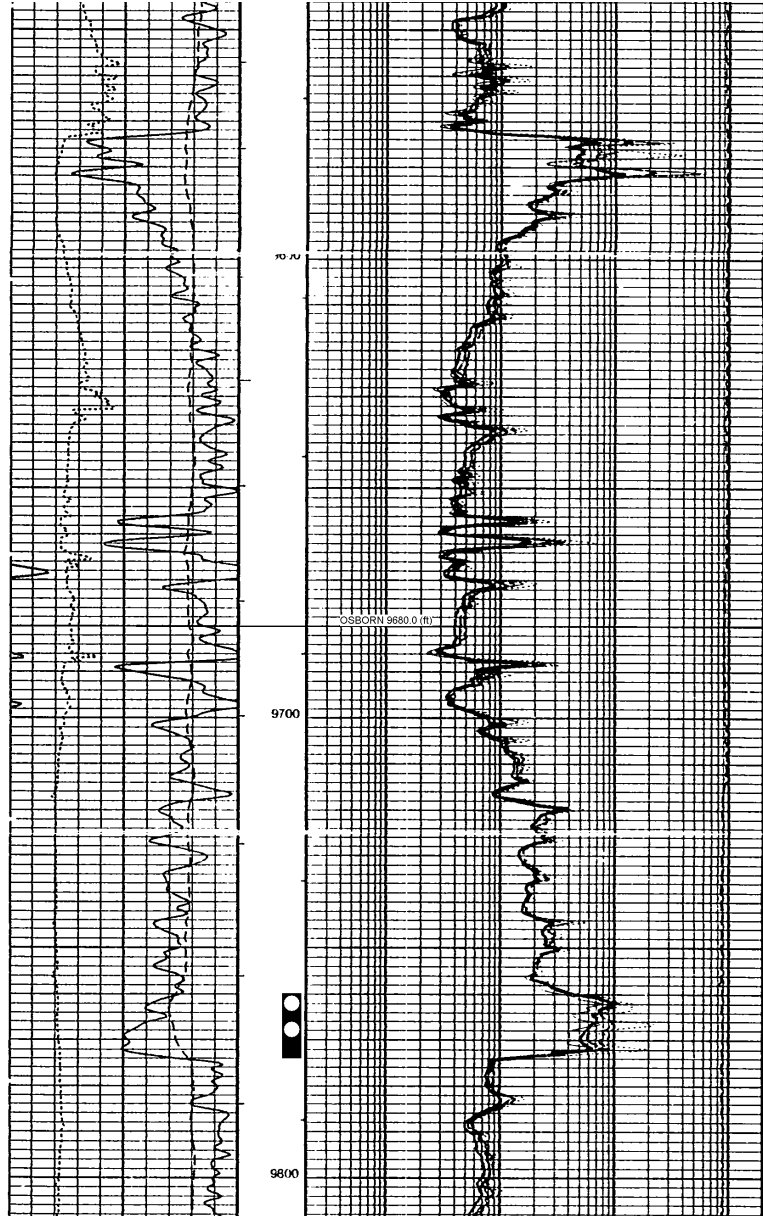
Near Shore
Marine

7

NOTE THE "SEVENS"

"Osborn Mbr"

Fluvial
Channel



Lower Skinner Mbr.

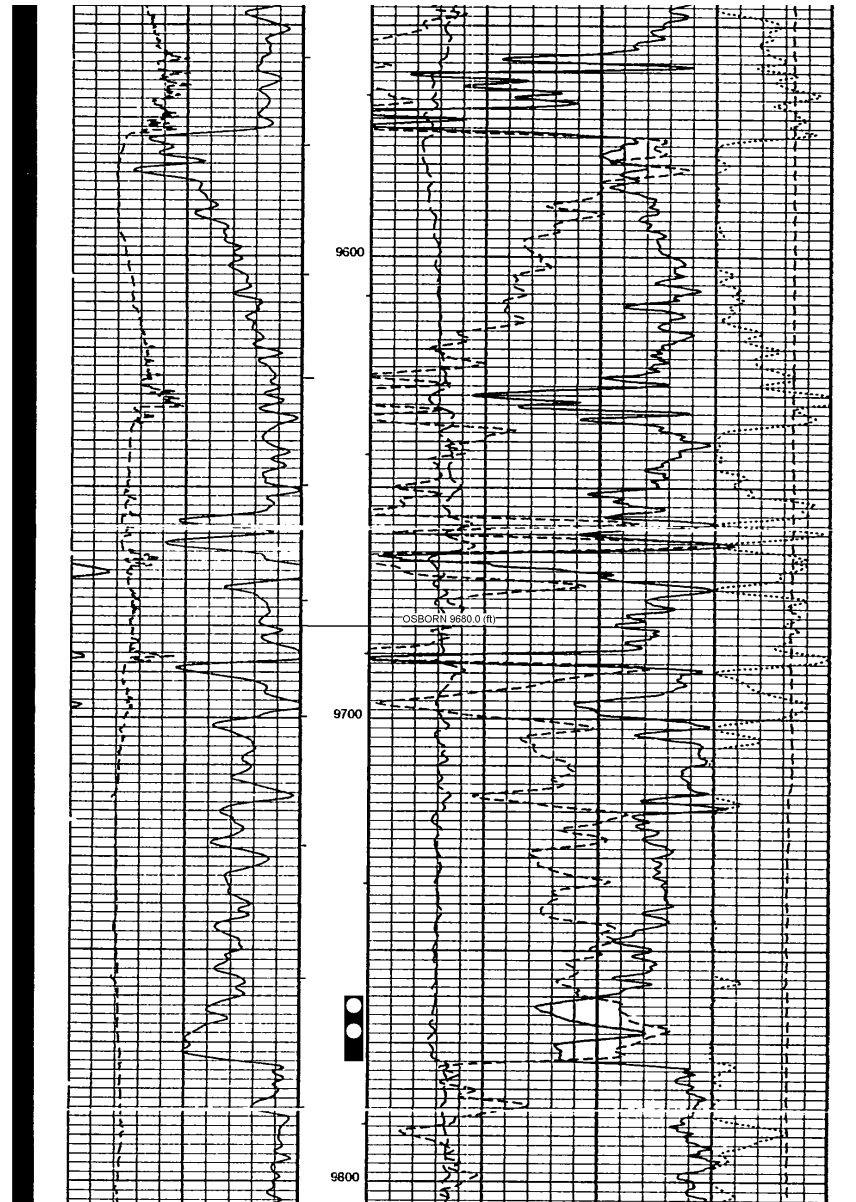
Pink Lime

Red Fork

- Section 34-11N-7W, Canadian Co., Okla.

Albers 1-34 Neutron/Density Log

10%

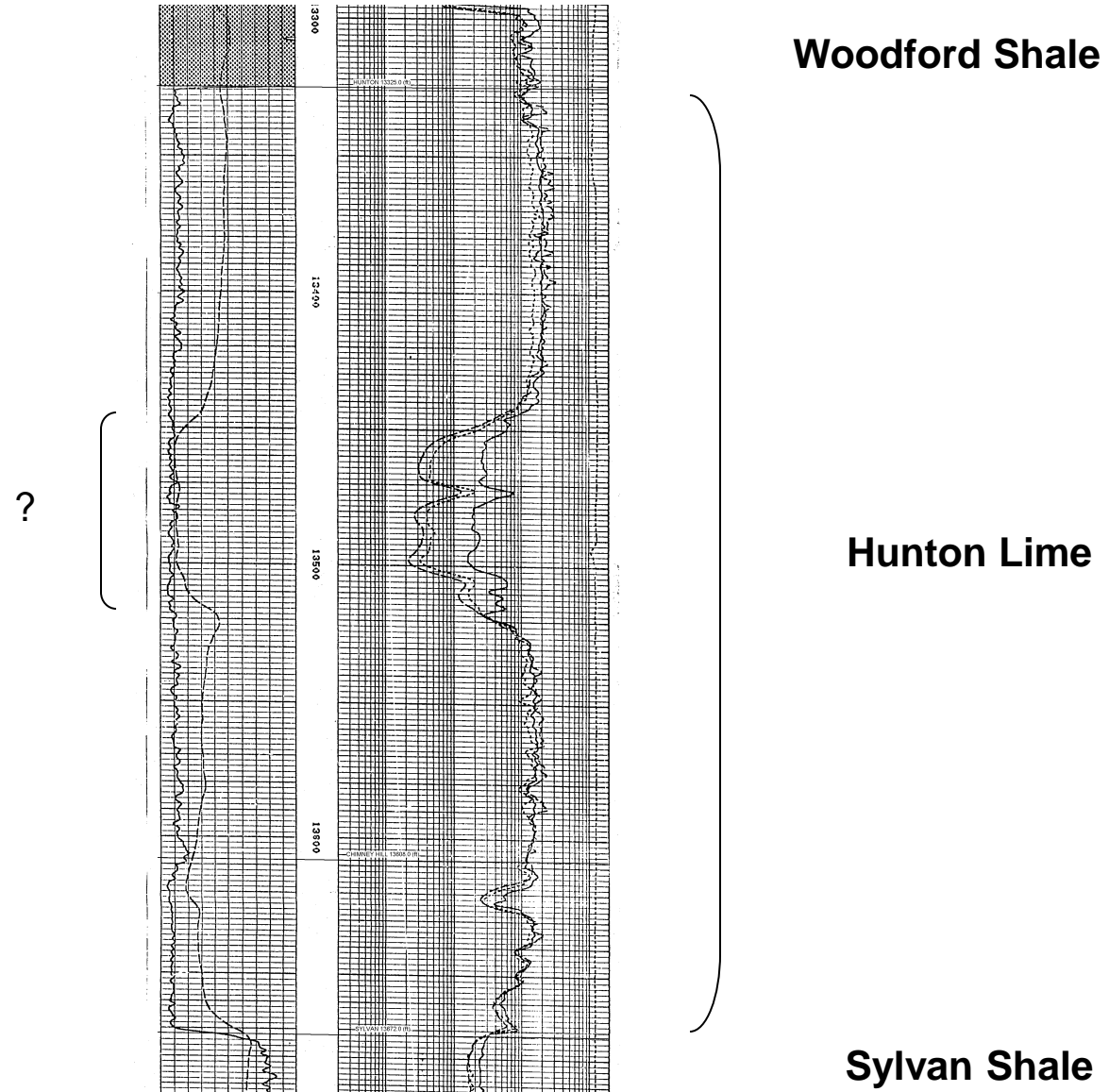


Lower Skinner Mbr.

Osborn Sand

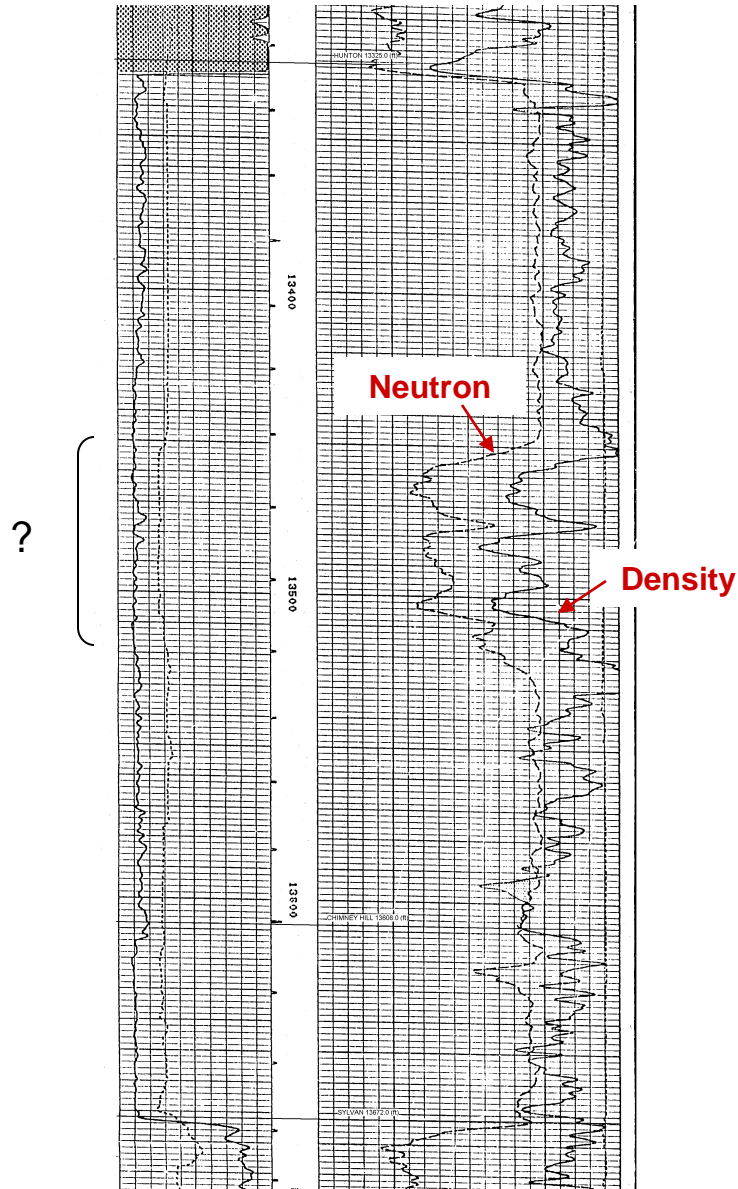
- Section 34-11N-7W, Canadian Co., Okla.

Gambrel 1 Dual Induction Log



- Section 13-17N-17W, Dewey Co., Okla.

Gambrel 1 Neutron/Density Log



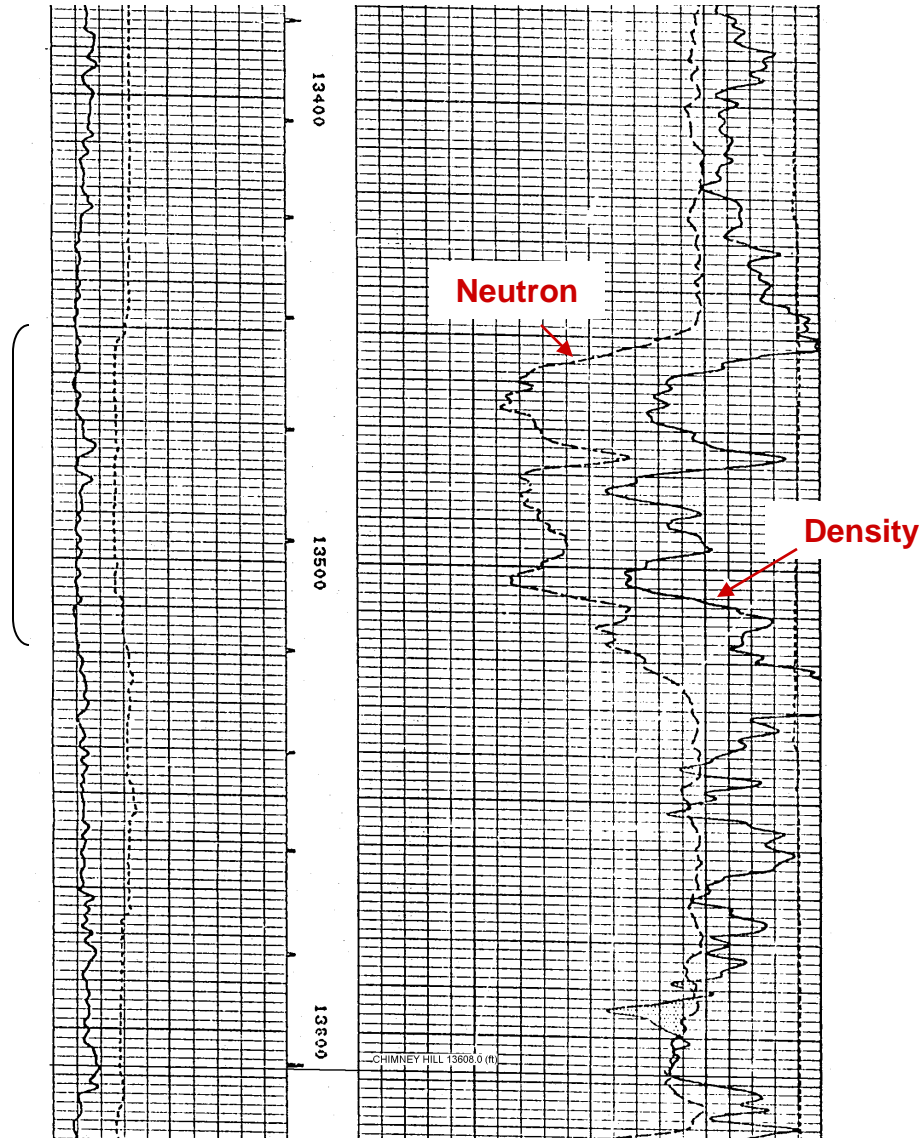
- Section 13-17N-17W, Dewey Co., Okla.

Gambrel 1 Neutron/Density Log

Limestone

Dolomite

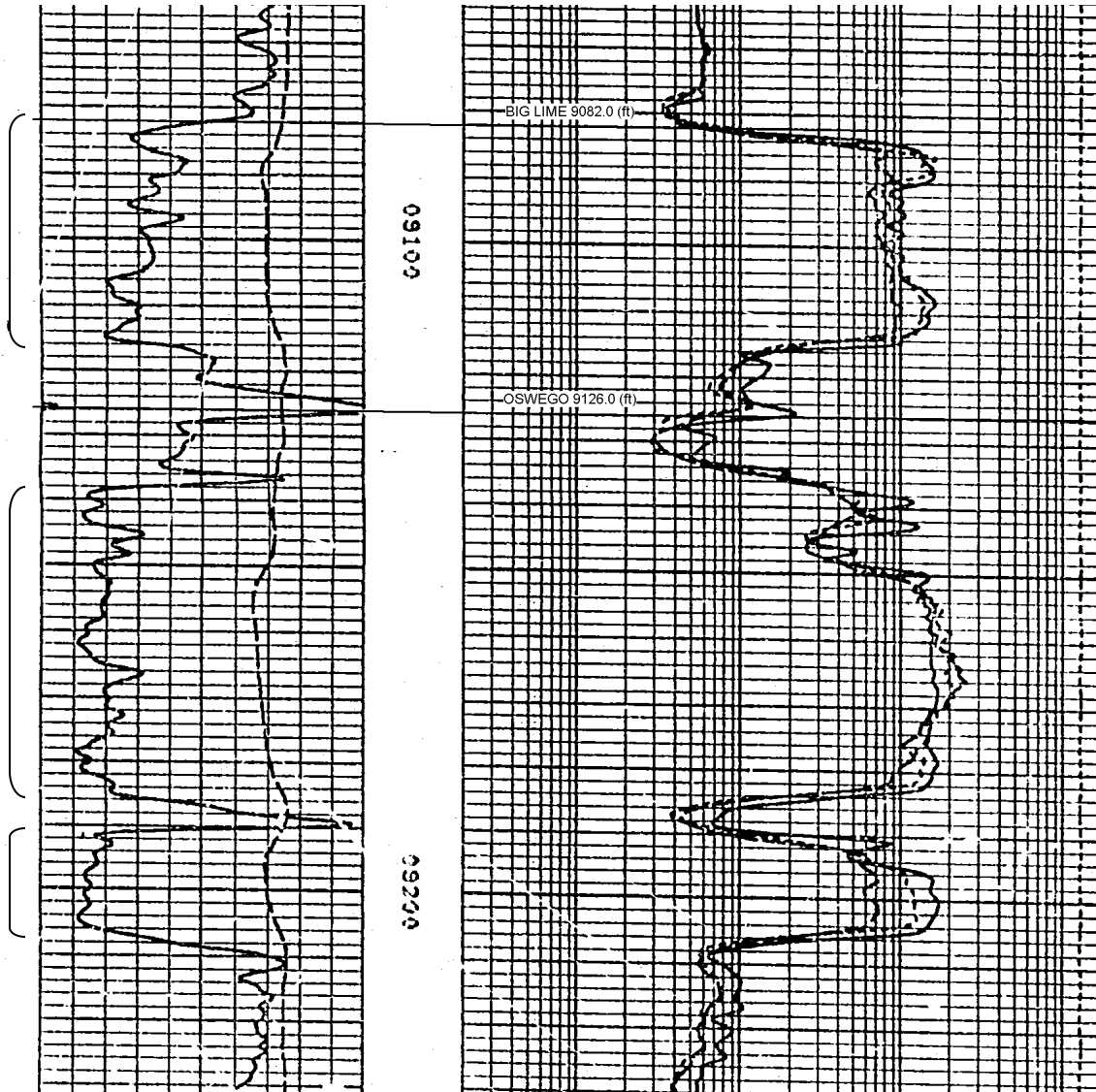
Limestone



- Section 13-17N-17W, Dewey Co., Okla.

Gambrel 1 Dual Induction Log

100



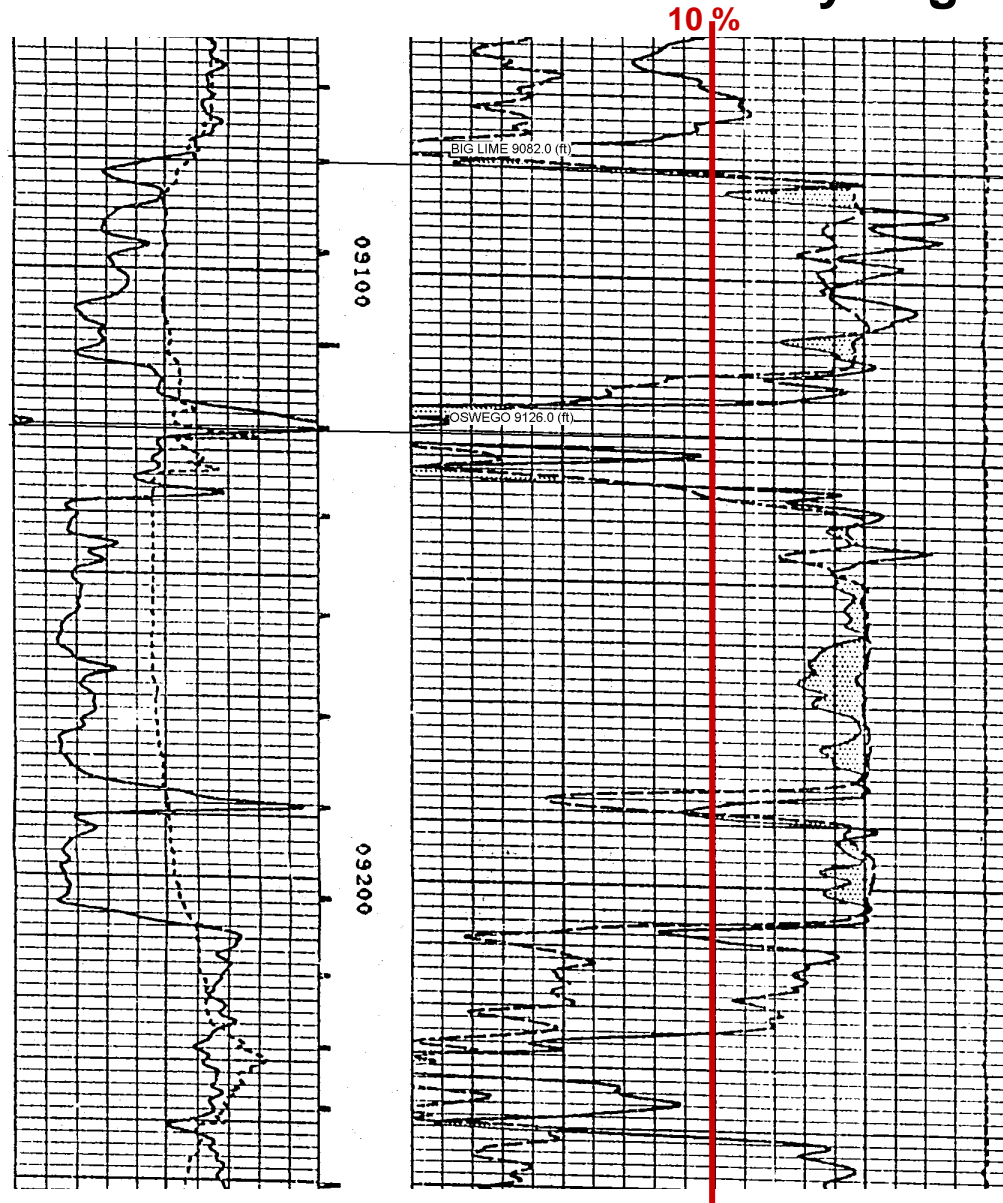
Lime Banks

Big Lime

Oswego Lime

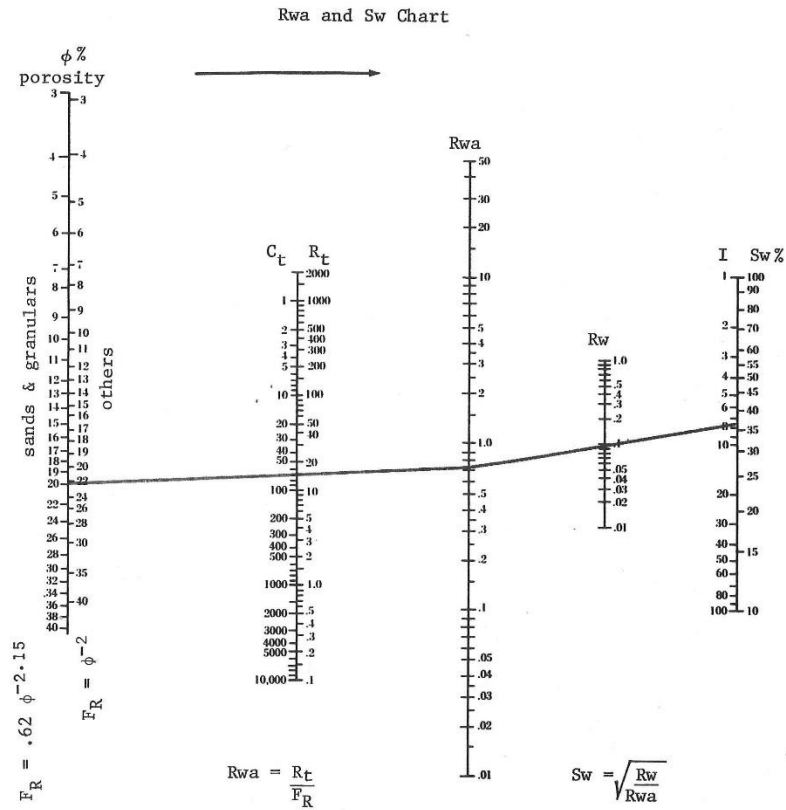
- Section 13-17N-17W, Dewey Co., Okla.

Gambrel 1 Neutron/Density Log



- Section 13-17N-17W, Dewey Co., Okla.

Figure 2-6 Solution of Water Saturation Equation (2-10)



Example: Porosity is 20% in a sandstone. R_t is 15. These give you an R_{wa} of 0.72. If R_w is .095 then $Sw = 36\%$

Density log:

$$\rho_b = \Phi \cdot \rho_f + (1 - \Phi) \cdot \rho_{ma}$$

$$\Phi = \frac{(\rho_{ma} - \rho_b)}{(\rho_{ma} - \rho_f)}$$

ρ_{ma} :	Quartz	2.65 gm/cc	ρ_f :	1.00
	Calcite	2.71		
	Dolomite	2.87		

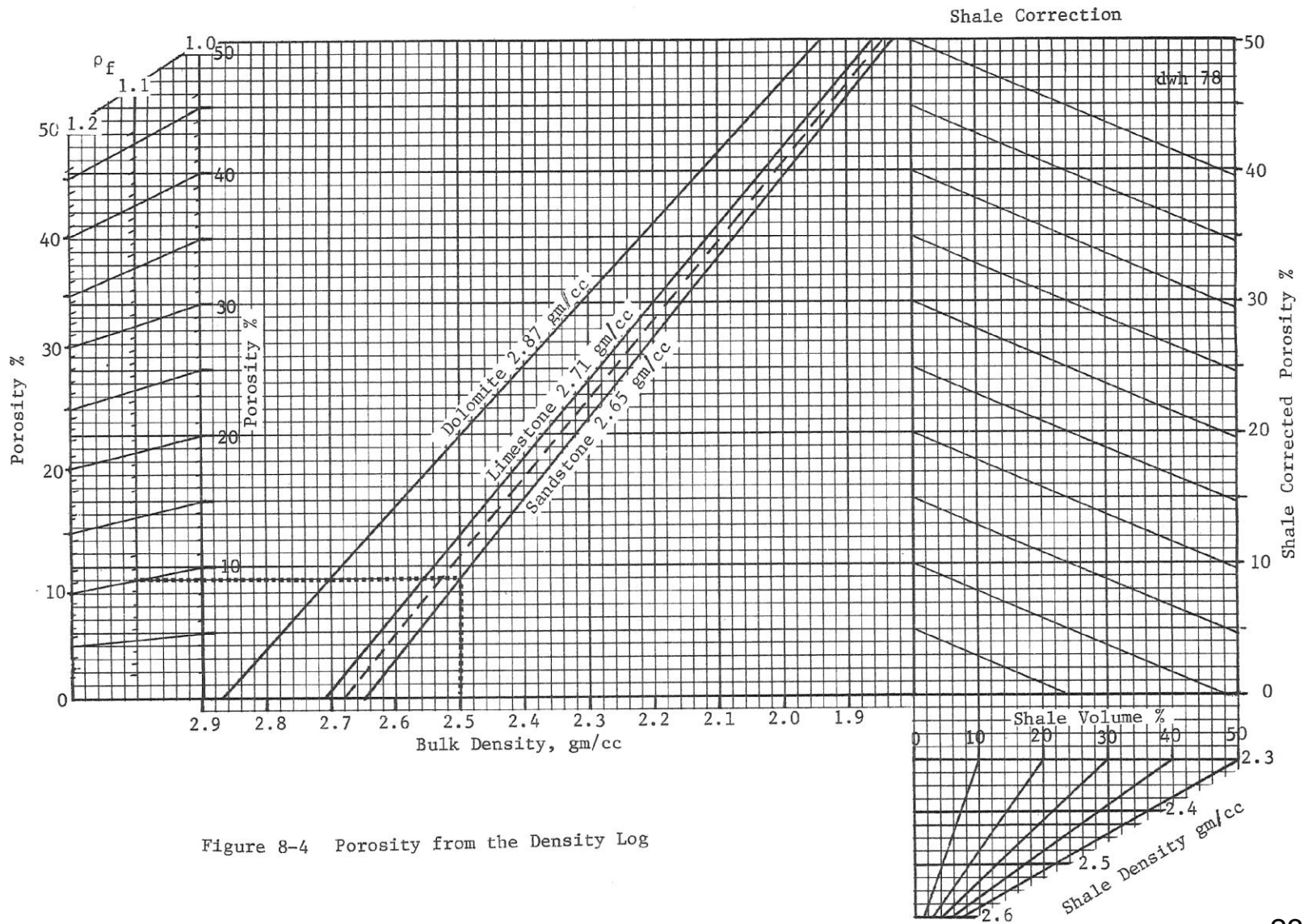


Figure 8-4 Porosity from the Density Log

13-7

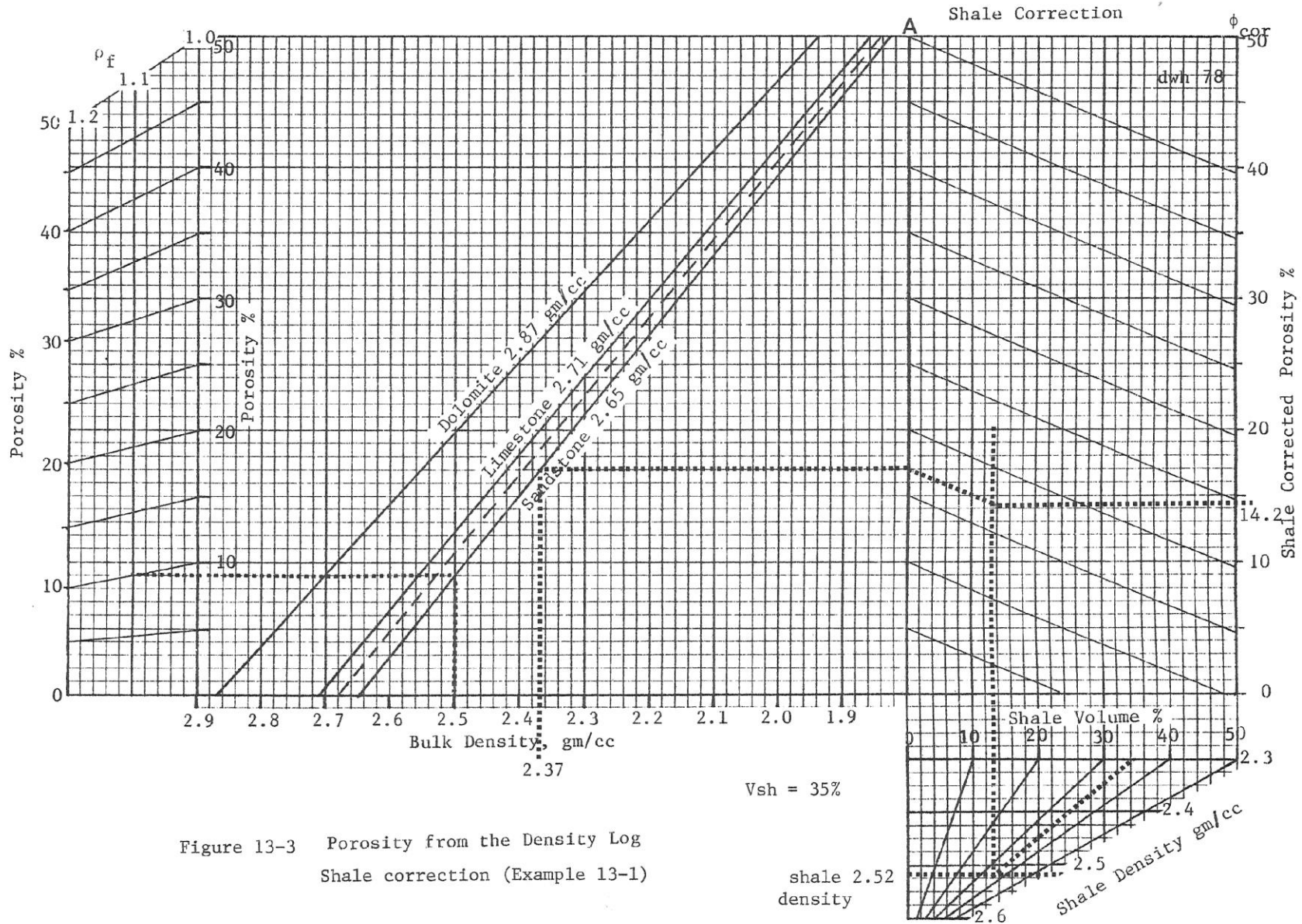


Figure 13-3 Porosity from the Density Log
Shale correction (Example 13-1)

Neutron log:

- The older neutron logs are either unscaled, recorded in "counts", or recorded in API units
- Most recent logs are recorded in *apparent* porosity units referred to either a sandstone, limestone, or dolomite scale. The default scale is limestone.

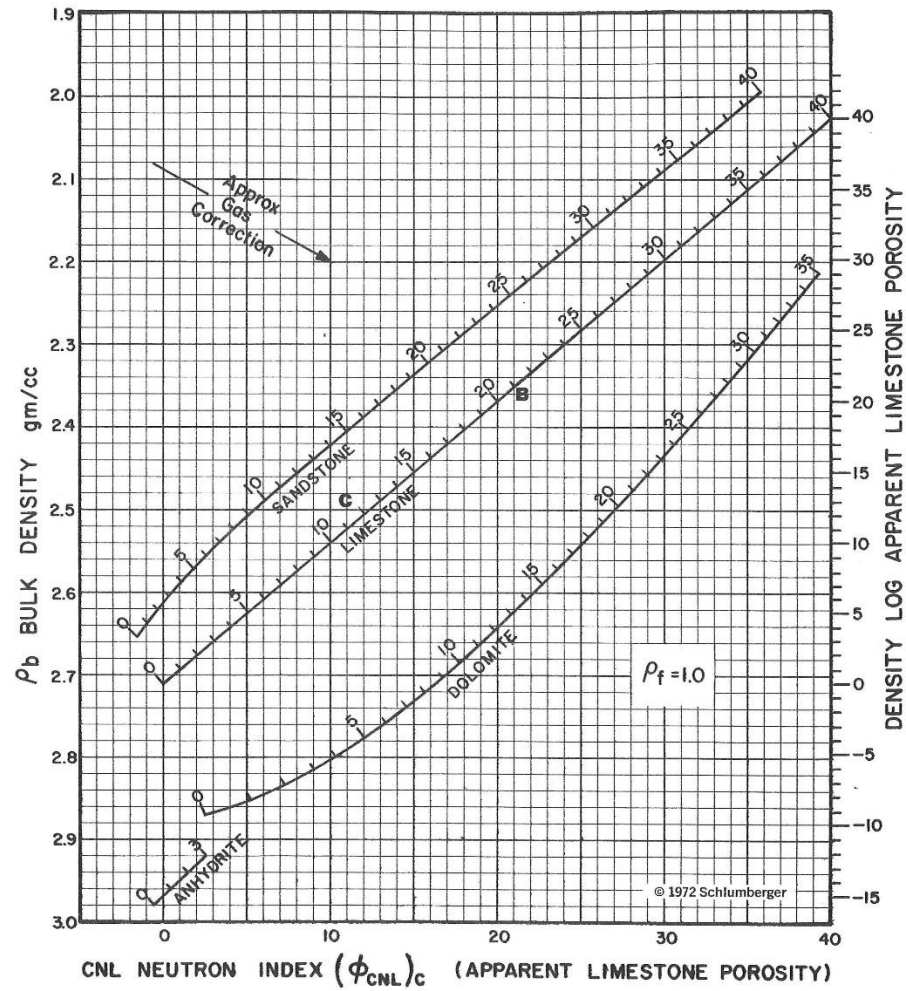
Neutron-density log combination:

When the neutron and density log are recorded as apparent porosity on a limestone scale, then a true volumetric porosity can be estimated by either:

$$\Phi = \frac{\Phi_n + \Phi_d}{2}$$

$$\Phi = \sqrt{\frac{(\Phi_n^2 + \Phi_d^2)}{2}}$$

Figure 10-2 Schlumberger density/compensated neutron crossplot
(courtesy Schlumberger)



Bob's Cheat Sheet

DENSITY LOG

<u>LITHOLOGY</u>	<u>Pma</u>	<u>3%</u>	<u>5%</u>	<u>6%</u>	<u>7%</u>	<u>8%</u>	<u>10%</u>	<u>12%</u>	<u>15%</u>
Dolomite	2.87	2.81	2.78	2.76	2.74	2.72	2.68	2.646	2.59
Limey Dolo.	2.80	2.75	2.71	2.69	2.67	2.66	2.62	2.58	2.53
Dolo. Limestone	2.75	2.70	2.66	2.65	2.63	2.61	2.575	2.54	2.49
Limestone	2.71	2.66	2.62	2.61	2.59	2.57	2.54	2.50	2.45
Limey Sandstone	2.68	2.63	2.60	2.58	2.56	2.546	2.51	2.48	2.43
Sandstone	2.65	2.60	2.57	2.55	2.53	2.52	2.485	2.45	2.40

<u>Mineral</u>	<u>Density - gm/cc</u>	$\phi = \frac{P_{ma} - P_b}{P_{ma} - P_f}$
Calcite	2.71	
Dolomite	2.87	
Quartz	2.65	
Anhydrite	2.95	
Gypsum	2.32	
Halite (NaCl)	2.16	
Sylvite (KCl)	1.98	
Coal	1-0-1.8	

SONIC LOG

<u>LITHOLOGY</u>	<u>tm</u>	<u>3%</u>	<u>5%</u>	<u>6%</u>	<u>7%</u>	<u>8%</u>	<u>12%</u>	<u>15%</u>
Dolomite	43.5	47.87	50.78	52.23	53.69	55.14	60.96	65.33
Limey Dolo.	44.5	48.84	51.73	53.17	54.62	56.06	61.84	66.18
Dolo. Limestone	45.5	48.8	52.68	54.1	55.55	56.98	62.7	67.03
Limestone	47	51.26	54.1	55.5	56.94	58.36	64.04	68.3
Limey Sandstone	52	56.1	58.85	60.22	61.59	62.96	68.44	72.55
Sandstone	55	59.02	61.7	63.04	64.4	65.7	71.1	75.1

<u>Sandstone</u>	<u>Carbonate</u>
$F = \frac{.62}{\phi^{2.15}}$ or $\frac{.81}{\phi^2}$	$F = 1/\phi^2$
$F = \frac{R_o}{R_w}$	$Sw = R_o/R_t = \sqrt{\frac{FR_w}{R_t}}$

Sonic log:

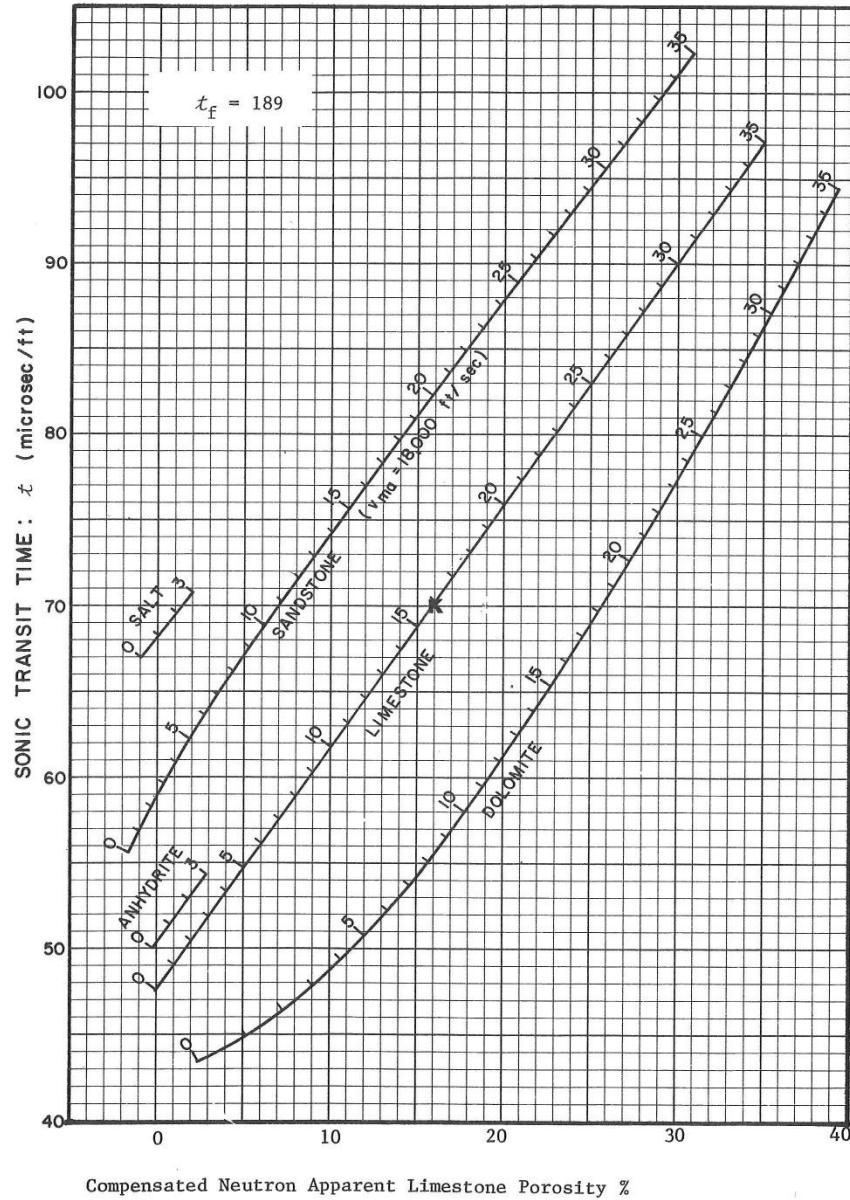
$$\Delta t = \Phi \cdot \Delta t_f + (1 - \Phi) \cdot \Delta t_{ma}$$

Wyllie time average equation:

$$\Phi = \frac{(\Delta t - \Delta t_{ma})}{(\Delta t_f - \Delta t_{ma})}$$

	Quartz	55.5 $\mu\text{sec/foot}$	
Δt_{ma} :	Calcite	47.5	Δt_f : 189
	Dolomite	43.5	

Figure 10-3 Schlumberger sonic/compensated neutron crossplot
(courtesy Schlumberger)



COUNTY	FORMATION	LOCATION		OPERATOR	LEASE	DEPTH	RW	TEMP	S	
		SEC	TWP							RANGE
CANADIAN	BIG LIME	11	13 N	6 W	MIDWEST OIL	HOLLOMAN 1	7360	0.038	100	P
CANADIAN	BIG LIME	11	13 N	6 W	MIDWEST OIL	HOLLOMAN 1	7359	0.038	100	P
CANADIAN	BIG LIME	12	13 N	6 W	MIDWEST OIL	HODGES 1	7355	0.040	100	P
CANADIAN	BOIS D'ARC	1	14 N	5 W			6997	0.041	100	U
CANADIAN	BOIS D'ARC	1	14 N	5 W			6997	0.041	100	T
CANADIAN	BOIS D'ARC	13	14 N	5 W			7065	0.042	100	U
CANADIAN	BOIS D'ARC	13	14 N	5 W			7065	0.038	100	T
CANADIAN	BOIS D'ARC	13	14 N	5 W			7065	0.042	100	T
CANADIAN	BOIS D'ARC	13	14 N	5 W			7065	0.038	100	U
CANADIAN	BOIS D'ARC	25	14 N	5 W			7110	0.044	100	U
CANADIAN	BOIS D'ARC	25	14 N	5 W			6984	0.047	100	T
CANADIAN	BOIS D'ARC	25	14 N	5 W			7110	0.042	100	T
CANADIAN	BOIS D'ARC	25	14 N	5 W			6984	0.047	100	U
CANADIAN	BOIS D'ARC	25	14 N	5 W			7110	0.042	100	U
CANADIAN	BOIS D'ARC	25	14 N	5 W			7110	0.044	100	T
CANADIAN	BROMIDE	15	12 N	6 W				0.030	100	U
CANADIAN	HUNTON	17	11 N	5 W	MAGNESS PETR	KINDL 17-1	8825	0.060	100	P
CANADIAN	HUNTON	7	11 N	6 W	RAMSEY ENGR.	KIRKEGARD #1	9657	0.057	100	U
CANADIAN	HUNTON	36	13 N	6 W	OKLA NAT GAS CO	OKLA #1		0.063	100	P
CANADIAN	HUNTON	1	14 N	5 W			7110	0.042	100	U
CANADIAN	HUNTON	1	14 N	5 W			7110	0.038	100	T
CANADIAN	HUNTON	1	14 N	5 W			7110	0.042	100	P
CANADIAN	HUNTON	1	14 N	5 W			7110	0.042	100	T
CANADIAN	HUNTON	12	14 N	5 W			7007	0.043	100	T
CANADIAN	HUNTON	12	14 N	5 W			7007	0.043	100	U
CANADIAN	HUNTON	13	14 N	5 W			7096	0.044	100	U
CANADIAN	HUNTON	13	14 N	5 W			7065	0.038	100	P
CANADIAN	HUNTON	13	14 N	5 W			7096	0.044	100	T
CANADIAN	MORROW	16	11 N	9 W	WOODS PETR	THOMAS 1		0.241	100	P
CANADIAN	MORROW	16	11 N	9 W	WOODS PETR	THOMAS 1		0.241	100	P
CANADIAN	MORROW	21	11 N	9 W	WOODS PETR	WILLARD 1		0.229	100	P
CANADIAN	MORROW	29	11 N	9 W	WOODS PETR	NUGENT 1	11971	0.254	100	T
CANADIAN	MORROW	1	11 N	10 W	JONES & PELLOW	RINEHART GROVES 1	11400	0.182	100	P
CANADIAN	MORROW	21	11 N	10 W			12500	0.323	100	P
CANADIAN	MORROW	21	11 N	10 W				0.318	100	U
CANADIAN	MORROW	18	12 N	9 W	AMOCO PROD	WITCHER UNIT	11079	0.237	100	P
CANADIAN	MORROW	18	12 N	9 W				0.190	100	U

19

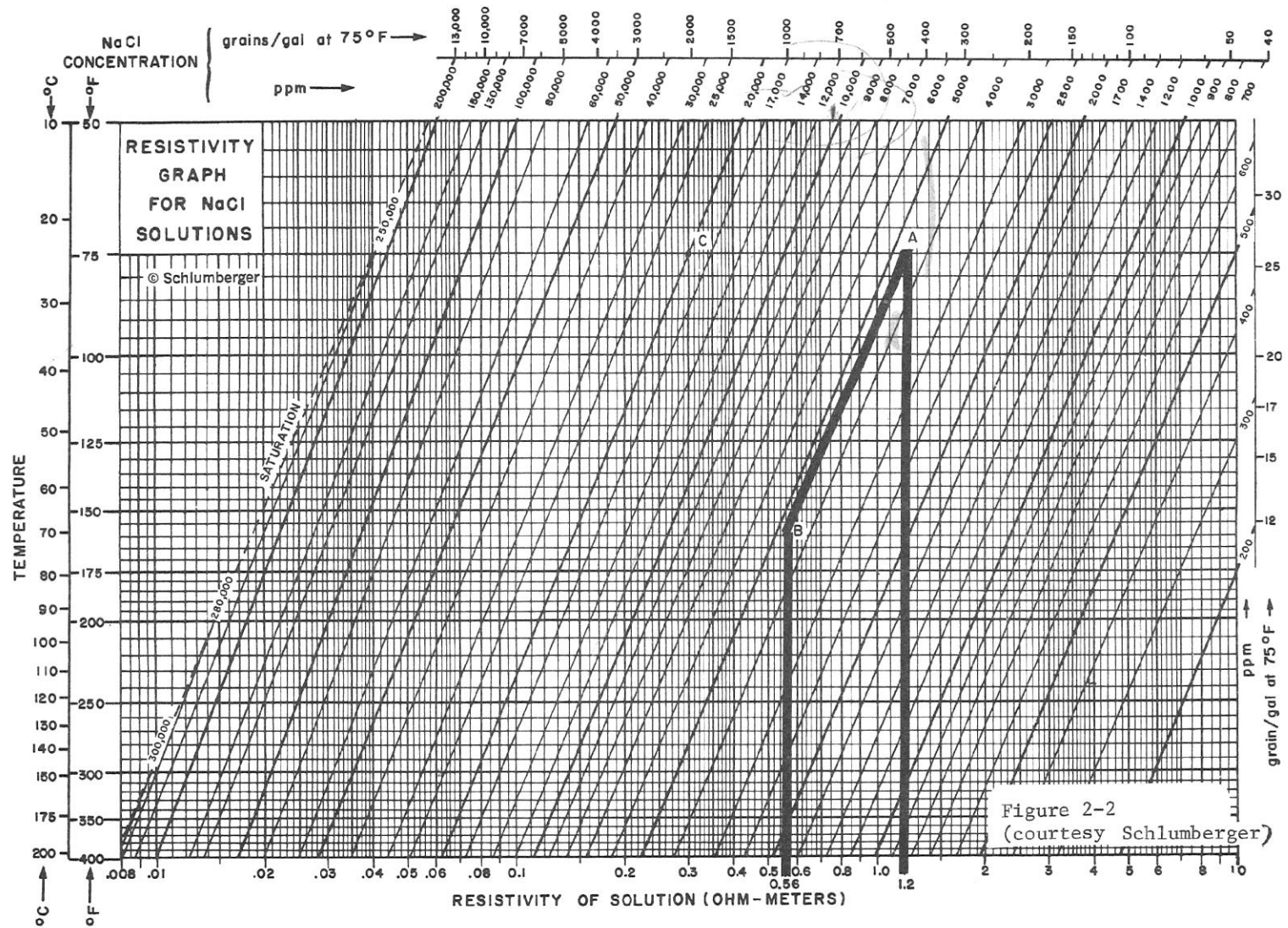


Figure 2-2
(courtesy Schlumberger)

Example: R_m is 1.2 at 75°F (point A on chart). Follow trend of slanting lines (constant salinities) to find R_m at other temperatures; for example, at Formation Temperature (FT) = 160°F (point B) read $R_m = 0.56$. The conversion shown in this chart is approximated by the Arps formula: $R_{FT} = R_{75} \times (75^\circ + 7) / (FT \text{ (in } ^\circ\text{F)} + 7)$.

Photoelectric factor log

