



# **SEMI-EMPIRICAL MODELING OF GAMMA- RAY DENSITY LOG WITH THE POSSIBILITY OF OBTAINING MORE INFORMATION**

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# TOPICS

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- INTRODUCTION**
- REVIEW OF EXISTING DENSITY LOG TREATMENT**
- THE FORWARD SCATTER MODEL APPROACH**
- MODEL VS. MONTE CARLO RESULTS**
- DISCUSSION, FUTURE WORK, AND CONCLUSIONS**

# REVIEW OF EXISTING DENSITY LOG TREATMENT



The text by Ellis describes the existing treatment of the density log. It consists of:

- ❑ a “spine and ribs” treatment in which the “spine” represents the average response and the “ribs” represent variations due to variations in mudcake density and composition.
- ❑ the bulk density is taken as equal to the porosity times the fluid density plus one minus the porosity times the rock matrix density:  $\rho_b = \Phi \rho_f + (1 - \Phi) \rho_m$   
 where  $\rho_b$ ,  $\rho_f$ , and  $\rho_m$  are the bulk, fluid, and rock matrix densities, respectively, and  $\Phi$  is the porosity of the rock.
- ❑ the main response is considered to be a simple exponential with no variation due to the photoelectric effect (rock composition) and no pre-exponential scattering term.  
 the equation used is:  $\Phi = \Phi_0 \exp[-\rho_b (Z/A) N_0 \sigma x]$

# REVIEW OF EXISTING DENSITY LOG TREATMENT, 2



- Ellis defines the electron density index to be:

$$\rho_e = 2(Z/A) \rho_b$$

- So that the tool response (or measured flux  $\Phi$ ) can be specified as proportional to:

$$\Phi \propto \exp(-\rho_e x)$$

- Ellis notes that the bulk density and electron density index are almost identical for the three most common rock minerals; calcite or limestone,  $\text{CaCO}_3$ , Quartz,  $\text{SiO}_2$ , and dolomite,  $\text{CaCO}_3\text{MgCO}_3$ , but both filling fluids of oil or water contain H with a  $Z/A \approx 1$  so  $\rho_e$  is not exactly proportional to  $\rho_b$ .

- So the transform between the measured  $\rho_e$  and the log measured density,  $\rho_{\log}$  is adjusted so that water and limestone are a linear relationship:

$$\rho_{\log} = 1.0704 \rho_e - 0.188$$

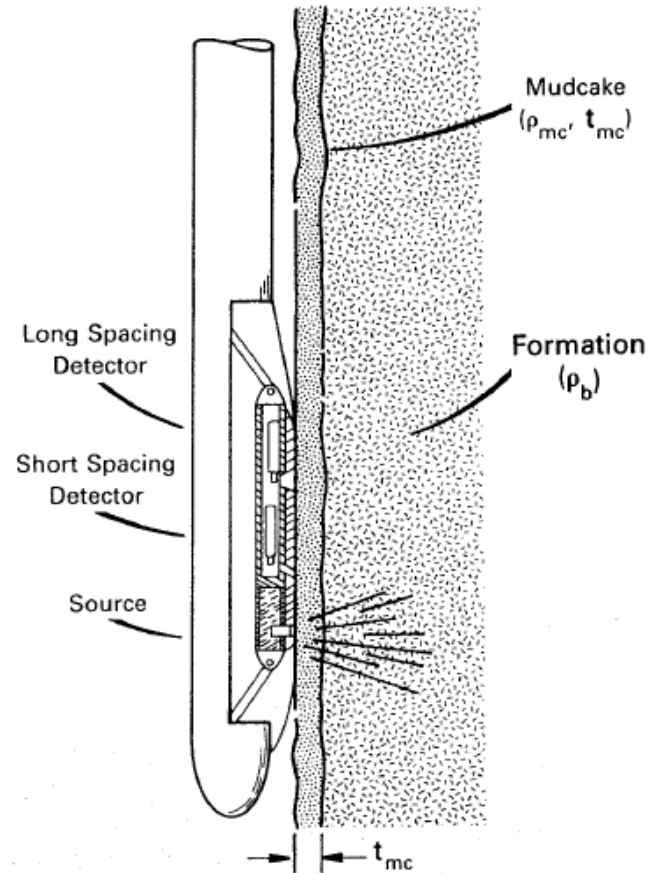
# REVIEW OF EXISTING DENSITY LOG TREATMENT, 3

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- ❑ Unfortunately, the Ellis treatment does not directly take into account the photoelectric effect (absorption) interaction, which is also very important and also does not take into account the scattering “in” that is also important.
- ❑ Ellis uses the simple exponential model:
$$N=N_0\exp(-\mu\rho x)$$
- ❑ At the time of the Ellis book “wireline” logs were being used exclusively and the main density measurement interference was mudcake from the drilling mud. The resulting treatment of this is shown in the following slides.

# REVIEW OF EXISTING DENSITY LOG TREATMENT, 4



**Figure 10-3.** A formation density device in the borehole situation applied to the borehole wall and separated from it by the thickness of the mudcake,  $t_{mc}$ . From Ellis et al.<sup>4</sup>

# REVIEW OF EXISTING DENSITY LOG TREATMENT, 5

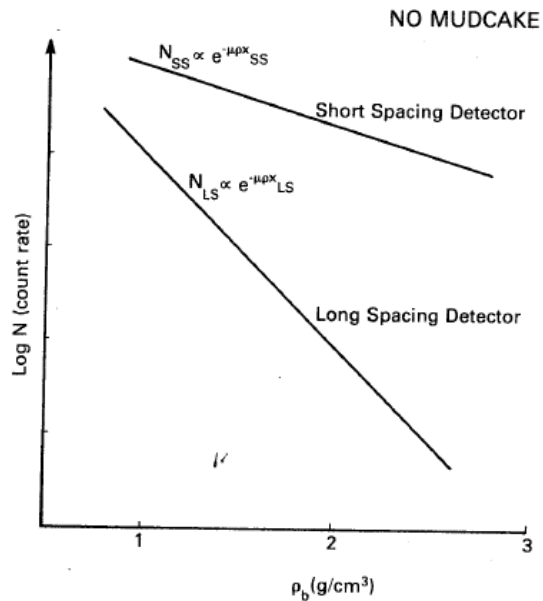


Figure 10-4. Idealized counting rate response of two detectors for variation of formation density in the presence of no mudcake.

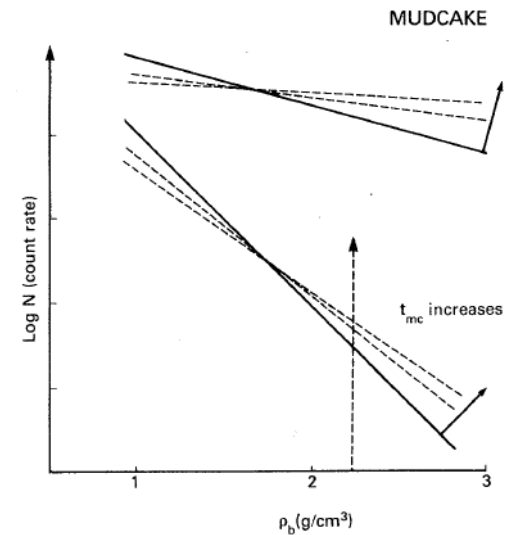


Figure 10-5. Idealized response of the two detectors for two different thicknesses of mudcake. Its density is approximately  $1.8 \text{ g/cm}^3$ . For a given thickness of mudcake and formation density/mudcake density contrast, both detectors experience approximately the same percentage of change in counting rate.

# REVIEW OF EXISTING DENSITY LOG TREATMENT, 6

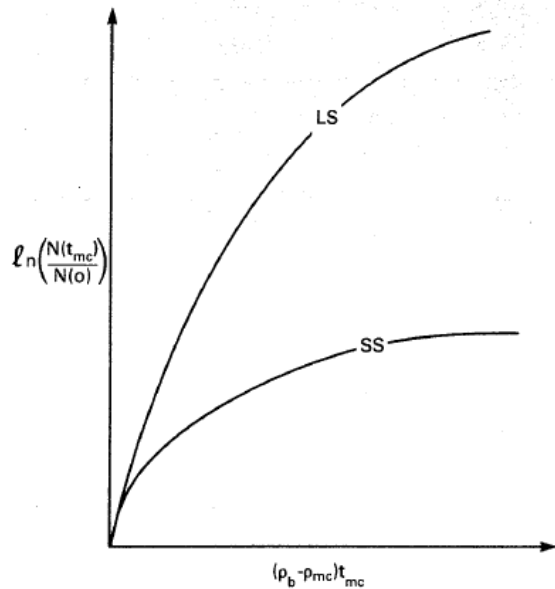


Figure 10-6. Effect of mudcake on the normalized counting rate of near and far detectors. The controlling parameter is the contrast in mudcake and formation density multiplied by the mudcake thickness.

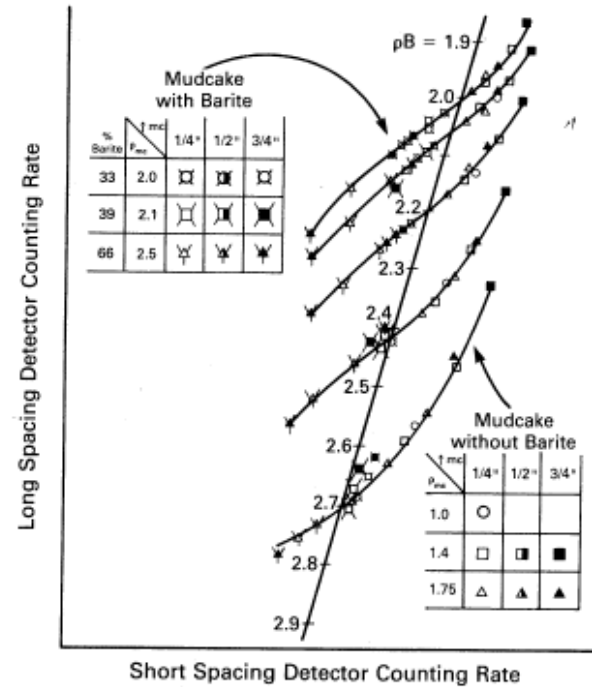


Figure 10-7. A "spine and ribs" representation of the response of a two-detector density device to formation density and mudcakes. Because of its characteristic outline, it is known as the "spine and ribs" chart. From Tittman et al.<sup>3</sup>



# REVIEW OF EXISTING DENSITY LOG TREATMENT, 7



- Ellis next treats “lithology” logging by first defining a new parameter called the “photoelectric index”,  $P_e$ , which is given by:

$$P_e = (Z/10)^{3.6}$$

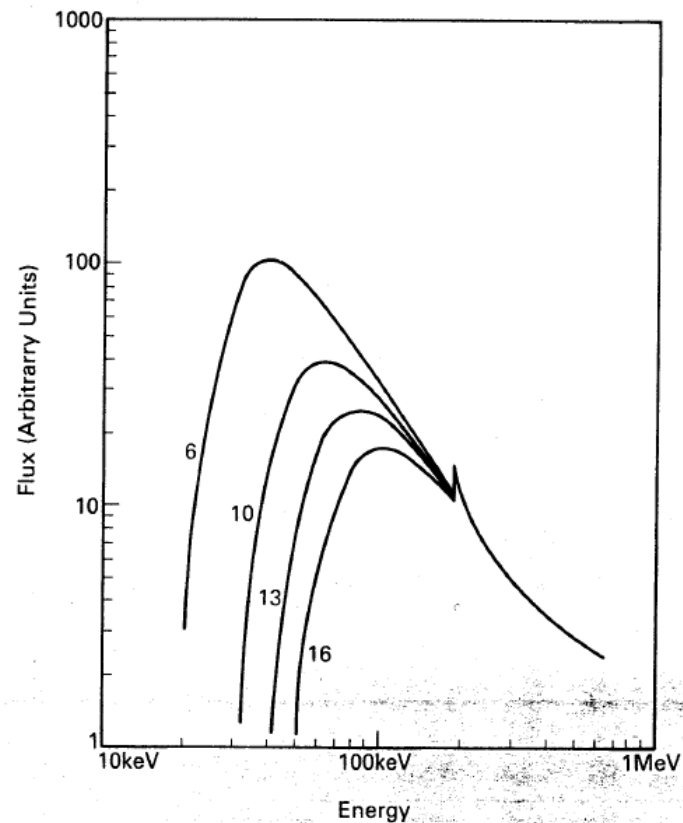
- Then the attenuation relationship can be given as:  $\Phi \propto \exp(-n_e P_e x)$

where  $n_e$  is the number density of electrons as in the case of Compton scattering.

- The gamma-ray spectra taken with recent logs yields the “lithology” as shown in the following three figures.



# REVIEW OF EXISTING DENSITY LOG TREATMENT, 9



**Figure 10–14.** A Monte Carlo calculation of the spectrum due to uniformly distributed sources, including the effect of photoelectric absorption. The evolution of the spectra is seen for materials of four different atomic numbers. As the atomic number increases, the low energy portion of the spectrum is attenuated. From Bertozzi et al.<sup>6</sup>

# THE FORWARD SCATTER MODEL APPROACH



- The basic response is taken as:

$$R=A_3X_S\exp(-A_1X_S-A_2X_A)+A_4$$

- Where the  $X_S$  and  $X_A$  are the scattering and absorption cross sections given by:

$$X_S= \rho \sum_i w_i Z_i/A_i$$

$$\text{and } X_A= \rho \sum_i w_i Z_i^{4.4}/A_i$$

- Note that the  $X_S$  term is essentially  $\rho/2$  for all elements other than hydrogen. For the common oil well logging rocks the variation in  $Z/A$  is less than 1% when no hydrogen is present.

# THE FORWARD SCATTER MODEL APPROACH, 2



- So if we express  $X_S$  in terms of either hydrogen and all other elements (or water or oil and all other elements), we then have:

$$X_S = 0.99209 \rho w_H + 0.50000 \rho (1 - w_H)$$

$$\text{or } X_S = 0.50000 \rho + 0.49209 w_H \rho$$

- Likewise we can express  $X_A$  in the same manner as:

$$X_A = (1^{4.4}/1.00729) \rho w_H + (1 - w_H) \rho \sum_i w_i Z_i^{4.4}/A_i \quad (i > 1)$$

$$\text{or } X_A = 0.99209 \rho w_H + (1 - w_H) \rho \sum_i w_i Z_i^{4.4}/A_i \quad (i > 1)$$

# THE FORWARD SCATTER MODEL APPROACH, 3



- Now we can express the near and far detector responses as:

$$R_N = A_3 X_S \exp(-A_1 X_S - A_2 X_A) + A_4$$

$$R_F = B_3 X_S \exp(-B_1 X_S - B_2 X_A) + B_4$$

- These two equations can be solved simultaneously for  $X_A$  to give:

$$\begin{aligned} X_A &= A_1 X_S / A_2 + (1/A_2) \ln[(R_N - A_4) / (A_3 X_S)] \\ &= B_1 X_S / B_2 + (1/B_2) \ln[(R_F - B_4) / (B_3 X_S)] \end{aligned}$$

- From which we can define  $f(X_S)$  as:

$$\begin{aligned} f(X_S) = 0 &= (B_2 A_1 - A_2 B_1) X_S + B_2 \ln[(R_N - A_4) / (A_3 X_S)] \\ &\quad - A_2 \ln[(R_F - B_4) / (B_3 X_S)] \end{aligned}$$

# THE FORWARD SCATTER MODEL APPROACH, 4

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- The differential of  $f(X_S)$  wrt  $X_S$  is given by:  
$$df(X_S)/dX_S = B_2A_1 - A_2B_1 - B_2/X_S + A_2/X_S$$
- Now the root of  $X_S$  can be obtained by iterating on:  
$$X_{S_{n+1}} = X_{S_n} - f(X_{S_n}) / (df(X_{S_n})/dX_{S_n})$$
  
until the change in  $X_S$  is less than a specified value.
- This is called the Newton Iterative approach.

# THE FORWARD SCATTER MODEL APPROACH, 5



- Now the value of  $\rho$  can be obtained from:  
$$X_S = 0.50000 \rho + 0.49209 w_H \rho$$
  
or 
$$\rho = X_S / (0.50000 + 0.49209 w_H)$$
- So if  $w_H$  is known – say from the porosity tool, then  $\rho$  is easily and accurately determined from the density tool measurement – using the near and far responses with this model.
- Now we can check to see if this works by using Monte Carlo simulation to generate  $R_N$  and  $R_F$  responses.



# THE FORWARD SCATTER MODEL APPROACH, 6

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- Thirteen responses each were generated by Monte Carlo simulation with MCNP5 for a range of densities (and corresponding water contents) of limestone, sandstone, and shale for both the near and far detectors of a density log.
- Then the model parameters were determined for both near and far detectors by using the data from all three types of rock for both the near and far detectors. The parameters obtained are given in the following table. Monte Carlo and model responses are shown plotted in the four following figures following the table.

# MODEL VS. MONTE CARLO RESULTS – MODEL PARAMETERS



**FAR ( $\chi_{\eta}^2=1.46$ )**

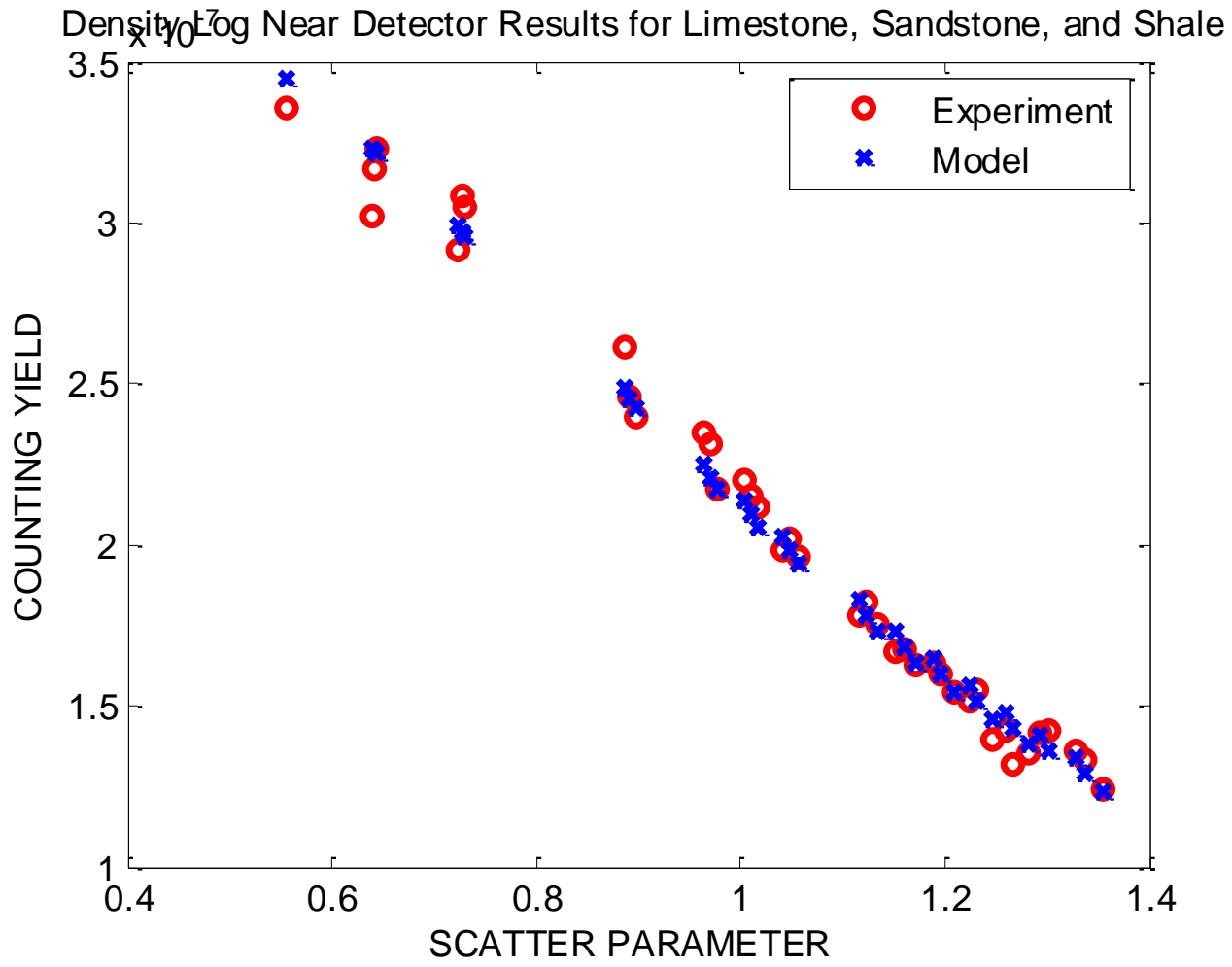
<b><math>A_1=5.85081</math></b>	<b><math>\sigma(A_1)=0.01015</math></b>
<b><math>A_2=5.38332E-5</math></b>	<b><math>\sigma(A_2)=2.542E-6</math></b>
<b><math>A_3=1.06696E-4</math></b>	<b><math>\sigma(A_3)=5.458E-7</math></b>
<b><math>A_4=1.14036E-8</math></b>	<b><math>\sigma(A_4)=1.012E-9</math></b>

**NEAR ( $\chi_{\eta}^2=0.870$ )**

<b><math>B_1=2.44637</math></b>	<b><math>\sigma(B_1)=0.01386</math></b>
<b><math>B_2=5.44215E-6</math></b>	<b><math>\sigma(B_2)=2.408E-6</math></b>
<b><math>B_3=2.27403E-6</math></b>	<b><math>\sigma(B_3)=4.596E-8</math></b>
<b><math>B_4=2.02098E-8</math></b>	<b><math>\sigma(B_4)=3.506E-9</math></b>

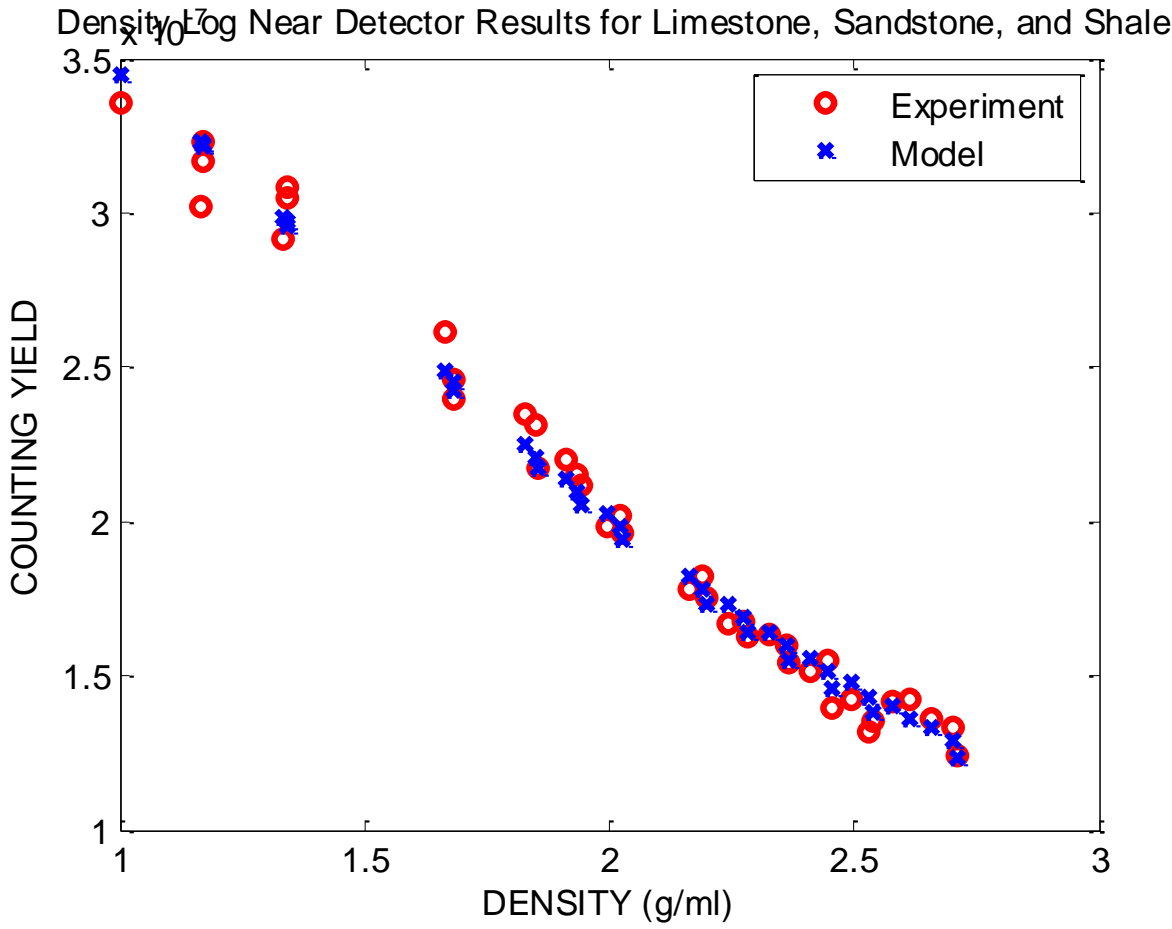


# NEAR DETECTOR RESULTS

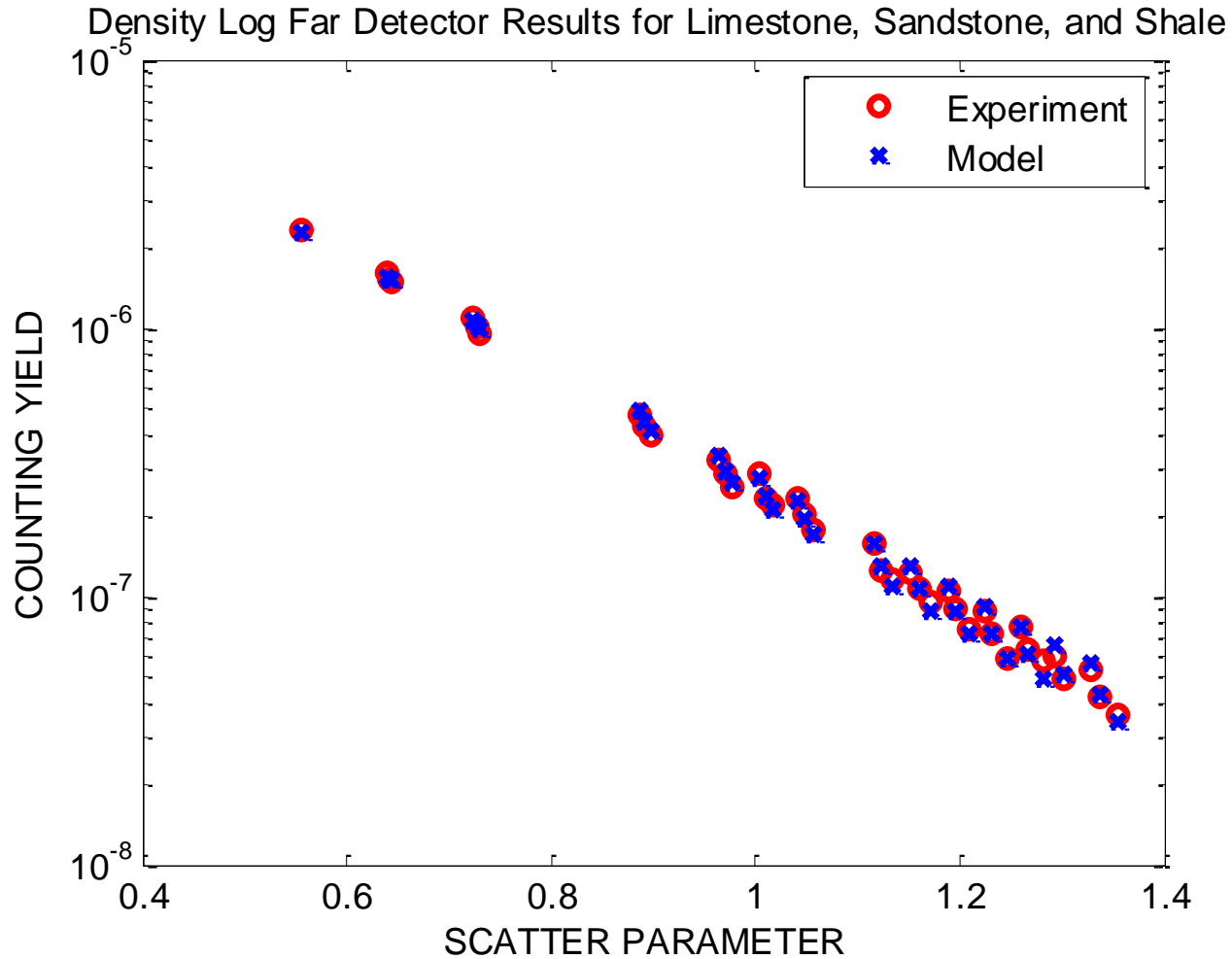




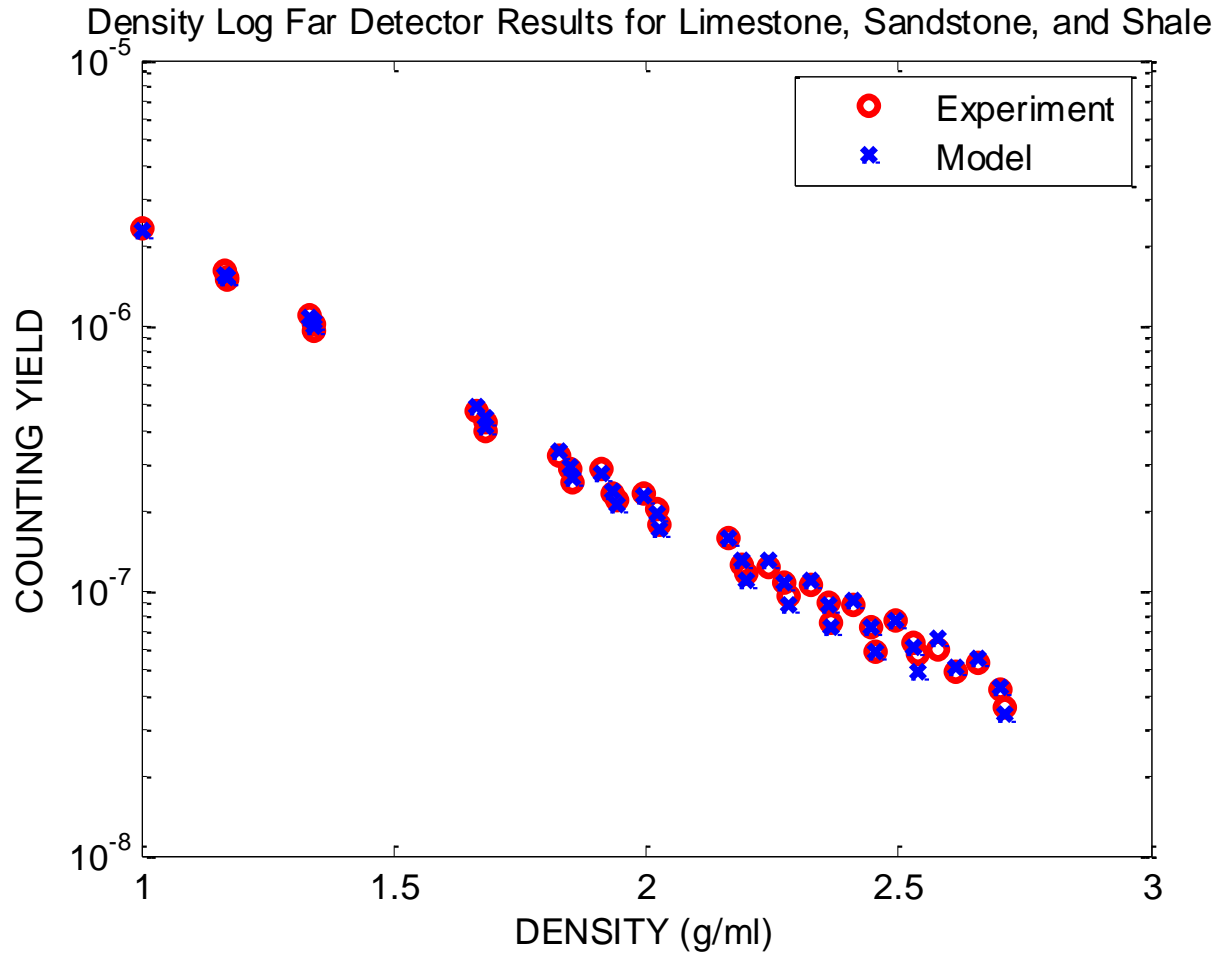
# NEAR DETECTOR RESULTS



# FAR DETECTOR RESULTS



# FAR DETECTOR RESULTS



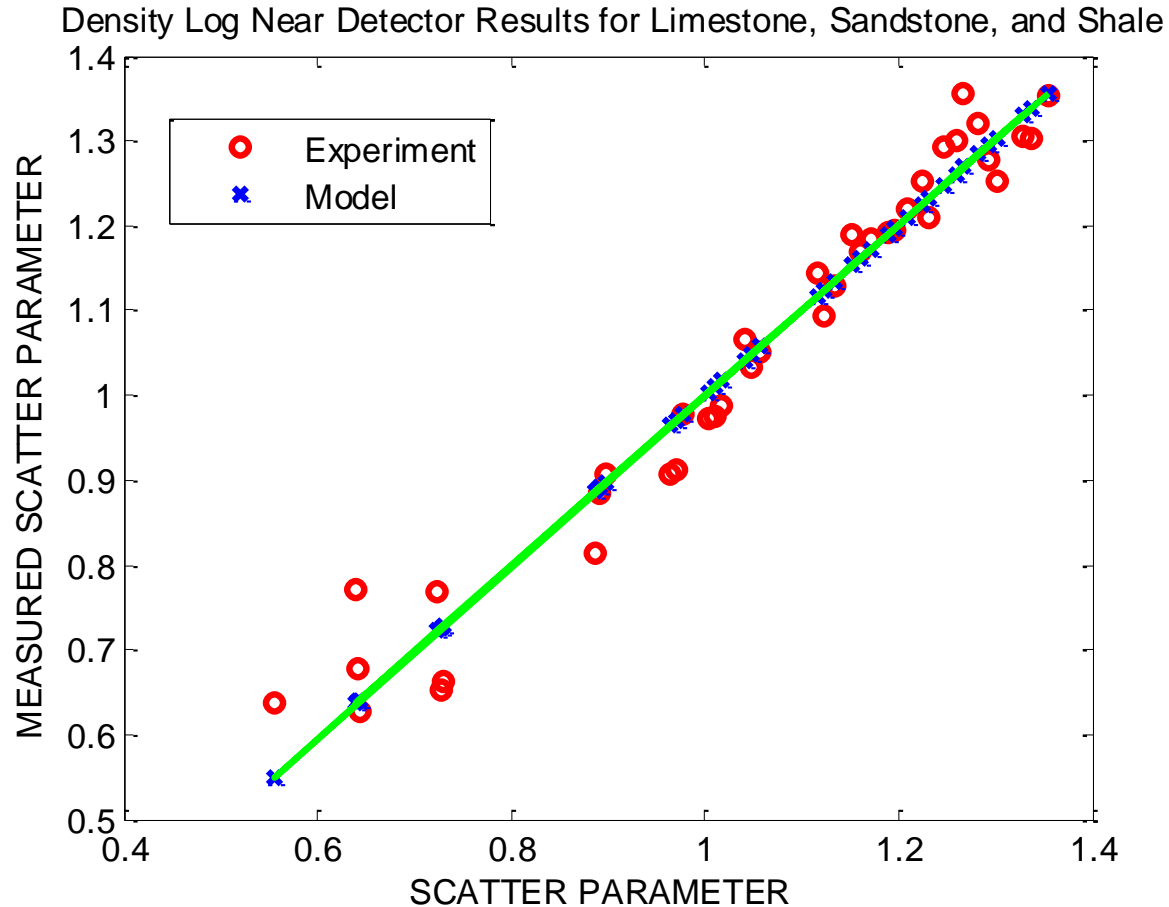
# THE FORWARD SCATTER MODEL APPROACH, 7

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- ❑ The scatter parameter  $X_S$  is found by the method outlined on slide 15.
- ❑ The scatter parameters for both the experimental and model based values are plotted in the following figure. Note that the scatter parameters for the model based case give perfect results as expected, while those for the experimental case give quite a bit of variation. We believe this variation is caused primarily by the near detector calculations, which exhibit fairly large variations. Calculations for a larger number of histories should be made for the near detector case.

# Determination of Scatter Parameter





# DISCUSSION, FUTURE WORK, AND CONCLUSIONS

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- The new approach has the major advantage that it can provide intuitive insight since it is a semi-empirical model rather than an empirical one and includes the effect of both scatter and absorption in an appropriate way.
- It is believed that the parameters that best describe the effect of scattering and absorption are  $X_S/\rho$  and  $X_A/\rho$  rather than  $\rho_e$  and  $P_E$ . The latter parameters include both density and the sum of  $Z/A$  and  $Z^n/A$ , while the new parameters separate the effect of composition and density.



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