

## **A comparison of hydrocarbon indicators derived from AVO analysis**

Hong Feng, Brian H. Russell and John C. Bancroft

### **ABSTRACT**

Numerous approaches have been published which derive fluid indicators (often called direct hydrocarbon indicators, or DHI) from AVO equations. The main idea behind these methods is to use the linearized Zoeppritz equations to extract petrophysical parameters such as P-impedance, S-impedance, bulk modulus, shear modulus, Lamé's parameters, Poisson's ratio, etc. and, from cross-plots of these parameters, infer the fluid content. Often, these indicators provide a good tool to quickly identify hydrocarbon zones. But the question of whether there is a best approach and, if so, which one it is, is still under debate. The purpose of this study is to examine which indicator can most easily discriminate a gas/oil sand from its background geology, and which indicator is most sensitive to pore-fluid content estimation.

### **INTRODUCTION**

The fluid factor ( $\Delta F$ ) was proposed by Smith and Gidlow (1987), and was derived by combining the linearized AVO equation with the mudrock line (Castagna et al., 1985). The authors also combined density and P-wave velocity changes by using Gardner's equation (Gardner et al., 1974). A version of the fluid factor which utilized density was introduced by Fatti et al. (1994). Goodway et al. (1997) suggested that Lamé's elastic parameters  $\lambda$  and  $\mu$  and their products with density could be useful tools in AVO analysis. Gray et al. (1999) showed how to estimate the parameters  $\rho\mu$  and  $\rho\lambda$  more directly by a new parameterization of the linearized AVO equation, as does Chen (1999). Russell et al. (2003) introduced the attribute  $I_p^2 - cI_s^2$  with  $c$  being a function of local  $(V_p/V_s)^2$ , where  $V_p$  and  $V_s$  are the dry rock P-wave and S-wave velocities.

In this study we used Gassmann fluid substitution to model changes in these parameters at given reservoir conditions in order to analyze the sensitivity of each fluid hydrocarbon indicator. We also examined the numerical example of the class 1, 2, and 3 sand models given by Hilterman (2001) to further examine the performance of these indicators.

### **METHODOLOGY**

The standard approach to AVO analysis is well known and was derived by Shuey (1985) based on the Aki-Richards (Aki and Richards, 1980) linearized formulation of the Zoeppritz equations. Their equation can be written as

$$R_p(\theta) = A + B \sin^2(\theta) + C \sin^2(\theta) \tan^2(\theta), \quad (1)$$

where  $A$  is the normal incident P-wave reflection coefficient  $R_p(0)$ ,  $B$  is the AVO gradient, and  $C$  is a curvature term that is often considered negligible. The product of  $A$

and B is often used to verify classical bright spots. This is based on the observation that low impedance gas sands encased in shale will have larger negative AVO intercepts (A) and a larger negative AVO gradient (B) than those not associated with gas. Thus A\*B should be an excellent indicator of a class III type gas sand. Furthermore, as shown by Swan (1993), product indicators have excellent S/N characteristics and may exhibit some degree of immunity to mild phase and velocity errors. On the other hand, A\*B will very effectively screen out Rutherford Class I (positive A, negative B) and Class II (A near-zero, negative B) gas sands. It can also be shown that, if  $V_p/V_s = 2$ , the scaled sum  $A+B$  gives an estimate of  $\Delta\sigma$ , the Poisson reflectivity where  $\sigma$  is Poisson's ratio, and the scaled difference  $A-B$  gives an estimate of  $R_s(0)$ , the zero offset shear-wave reflectivity.

The fluid factor as modified by Fatti et al. (1994) is defined by the equation  $\Delta F = R_p - g(t)R_s$ , where the terms  $R_p$  and  $R_s$  represent the P and S-wave normal incident reflectivities. The term  $g(t)$  represents a time varying scale factor given by the average ratio  $V_p/V_s$  multiplied by the linear coefficient in the mudrock line (Castagna et al., 1985). For wet reservoirs,  $\Delta F = 0$ , and for gas filled reservoirs it will be nonzero. It has been noted that the Fluid Factor and Poisson reflectivity are equivalent when  $V_p/V_s = 2$ . Lithologies that do not follow the  $V_p$  and  $V_s$  relationship for brine saturated clastics may produce a fluid factor anomaly. For example, coal produces an anomaly similar to gas sand reservoir. Since carbonates do not follow the criteria for clastic rocks, a carbonate equivalent to the "mudrock line" should be used for carbonate reservoirs.

Elastic constants are fundamental in seismology because the P and S-wave velocities depend on the elastic constants and density of the rock. Thus, the extraction of these elastic constants can help us to understand the rock behavior when the pore fluid changes.

Goodway et al. (1997) proposed the LMR ( $\lambda$ - $\mu$ - $\rho$ ) method, where L, M, and R represent:  $\lambda$ , the first Lamé parameter;  $\mu$  the shear modulus or coefficient of rigidity; and  $\rho$ , the density. Note also that  $\lambda = K - (2/3)\mu$ , where K is the bulk modulus. It was observed by Goodway (2001) that various combinations of  $\lambda$ ,  $\mu$ , and  $\rho$ , such as,  $\lambda\rho$ ,  $\mu\rho$ , and  $\lambda/\mu$ , show a better separation of gas sand from brine sands and shales than  $I_p$  and  $I_s$  where  $I_p$  is P-impedance and  $I_s$  is S-impedance. An effective fluid indicator can be found on a cross-plot of  $\mu\rho = I_s^2$  versus  $\lambda\rho = I_p^2 - 2I_s^2$ . The intuitive interpretation of this is that  $\lambda\rho$  and  $\mu\rho$  are fundamentally more orthogonal than  $I_p$  and  $I_s$ , stemming from the ambiguity in the  $V_p$  and  $V_s$  velocity relationships that share the same value rigidity  $\mu$ . Goodway (2001) also argued that the value of  $\lambda/\mu$  is a more sensitive indicator than  $\lambda$ ,  $\lambda\rho$ ,  $V_p/V_s$ , or  $\sigma$ , since the formulation  $\lambda/\mu = 2\sigma/(1-2\sigma)$  shows that for a given change in Poisson ratio we see an enhanced change in the  $\lambda/\mu$  value.

Batzle et al. (2001) proposed the fluid indicator  $K-\mu$ . It provides a direct look at the effect of pore fluid on the bulk modulus, and as such is quite similar to  $\lambda\rho$  and to the Gassmann fluid indicator described by Hilterman. This indicator works best in a sandstone regime where the clay content may vary from negligible to significant. This method comes from the observation that the bulk modulus for gas saturated or dry clastics is approximately equal to the shear modulus, which is consistent with the laboratory data of Han (1986).

Russell et al. (2003) proposed a method based on Biot-Gassmann theory to extract the fluid term  $\rho f = I_p^2 - cI_s^2$  from P-wave and S-wave impedances, where  $I_p$  and  $I_s$  are the P-wave and S-wave impedance respectively,  $\rho$  is density, and  $f$  represents the fluid term. The term  $c$  is the local  $(V_p/V_s)^2$ ,  $V_p$  and  $V_s$  are the P- and S-velocities for dry rock condition. The  $\rho f$  term is also equivalent to the  $\rho \Delta K$  term in Batzle et al. (2001), where  $\Delta K$  corresponds to the change in saturated bulk modulus due to fluid effects. As discussed by Russell et al., the Goodway (1997) formulation is a particular case of the  $I_p^2 - cI_s^2$  attribute, where  $c=2$  implies that the dry rock Poisson's ratio is 0. The fluid indicator  $K-\mu$  proposed by Batzle et al. (2001) is also a particular case of  $I_p^2 - cI_s^2$  attributes, corresponding to  $K_{dry} = \mu$  and  $c=2.333$ .

## SENSITIVITY ANALYSIS FOR FLUID INDICATORS

### Analysis using Gassmann fluid substitution equation

To analyze the sensitivity of each fluid indicator, we used the Gassmann fluid substitution equation (Gassmann, 1951) to model the changes of these parameters at given reservoir conditions. First, the bulk modulus ( $K_{sat}$ ) and shear modulus ( $\mu$ ) at in-situ conditions can be estimated from the wire line log data by the equations,

$$K = \rho \left( V_p^2 - \frac{4}{3} V_s^2 \right), \text{ and} \quad (2)$$

$$\mu = \rho V_s^2, \quad (3)$$

where,  $V_p$  and  $V_s$  are the P-wave and S-wave velocity,  $K$  and  $\mu$  are the bulk and shear moduli, and  $\rho$  is the mass density. Using Gassmann fluid substitution, the bulk modulus for dry rock can be derived by the equation

$$K_d = \frac{K_{sat} \left( \frac{\phi K_m}{K_f} + 1 - \phi \right) - K_m}{\left[ \frac{\phi K_m}{K_f} + \frac{K_{sat}}{K_m} - 1 - \phi \right]}, \quad (4)$$

where,  $K_{sat}$ ,  $K_d$ ,  $K_m$ ,  $K_f$  are the bulk moduli of the saturated rock, dry rock, mineral matrix, and pore fluid, respectively, and  $\phi$  is porosity. Then, with known  $K_d$ , we can estimate  $V_p$  and  $V_s$  when the pore fluid changes. Finally, various hydrocarbon indicators can be estimated.

We examined the effect of fluid properties on the fluid indicators. The initial state is  $V_p=11000\text{ft/s}$ ,  $V_s=6500\text{ft/s}$  and  $\rho=2.2 \text{ g/cm}^3$ . Figures 1 and 2 show the results for each indicator calculated from the Gassmann fluid substitution equation, including  $\lambda$ ,  $\mu$ ,  $\sigma$ ,  $k-\mu$ ,  $I_p^2 - cI_s^2$ ,  $R_p - gR_s$  and  $\lambda/\mu$ . From Figure 1 and Figure 2, we notice that  $I_p^2 - cI_s^2$  is the term most sensitive to the fluid change, followed by  $K-\mu$  and  $\lambda$ .

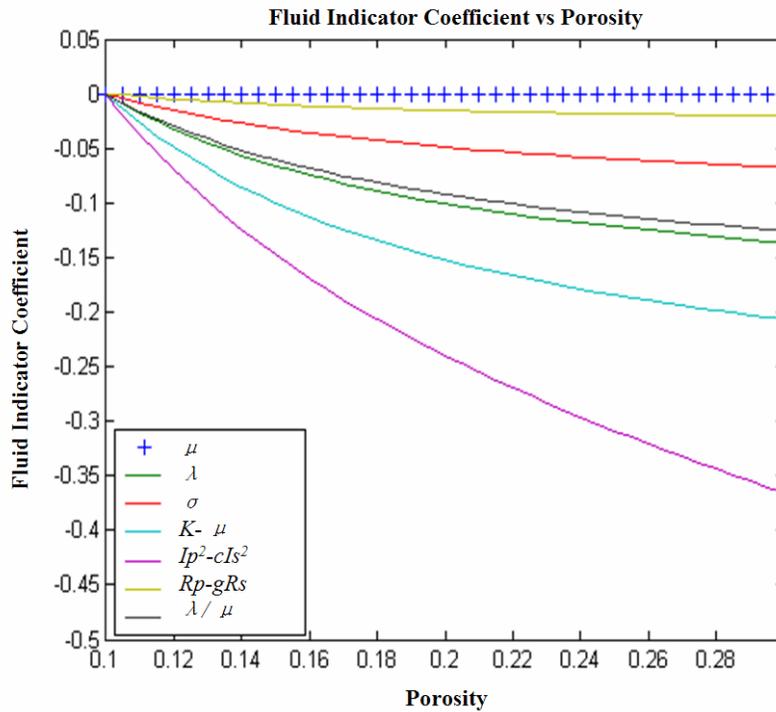


FIG. 1. Fluid indicator coefficients versus porosity. These indicators are calculated using the Gassmann fluid substitution equation with the initial state:  $V_p=11000\text{ft/s}$ ,  $V_s=6500\text{ft/s}$ ,  $S_w=0.2$ ,  $\rho=2.2\text{ g/cm}^3$ . From these curves, the indicator  $Ip^2-cl_s^2$  is most sensitive to the porosity change.

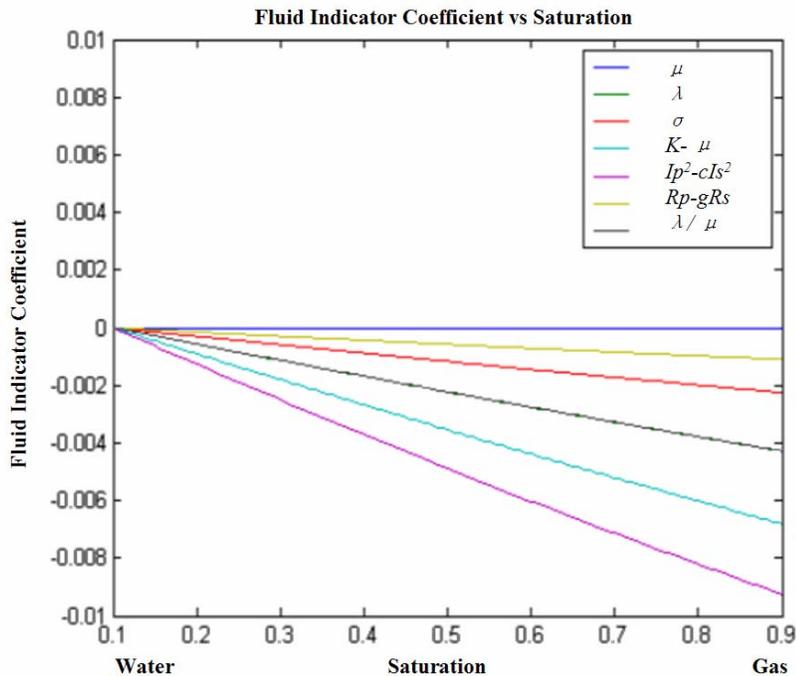


FIG. 2. Fluid indicator coefficients versus saturation. These indicators are calculated using the Gassmann fluid substitution equation with the initial state:  $V_p=11000\text{ft/s}$ ,  $V_s=6500\text{ft/s}$ , Porosity=0.2 and  $\rho=2.2\text{ g/cm}^3$ . From these curves, the indicator  $Ip^2-cl_s^2$  is most sensitive to the saturation change.

**Analysis using numerical example**

To diagnose the sensitivity of each indicator, let us examine the class I, II, and III sand models given by Hilterman (2001). These models were derived from Gulf of Mexico data. The following twelve indicators:  $\Delta\alpha/\alpha$ ,  $\Delta\beta/\beta$ ,  $\Delta\rho/\rho$ ,  $\Delta K$ ,  $\Delta\mu$ ,  $\Delta\lambda$ ,  $\Delta Y$ ,  $\Delta\sigma/\sigma$ ,  $\Delta(Vp/Vs)/(Vp/Vs)$ ,  $K-\mu$ ,  $Ip^2-cl_s^2$  and  $\Delta F$ , were chosen for discrimination. The  $c$  term in the  $Ip^2-cl_s^2$  fluid indicator was chosen to be 2.85. Each fluid indicator coefficient diagnoses the sensitivity to fluid discrimination and is defined as the difference between shale and wet sand or shale and gas sand divided by the value related to the shale reference. From Figures 3 through 5 we can observe that the indicators  $Ip^2-cl_s^2$ ,  $\Delta K$ ,  $K-\mu$  and  $\Delta\lambda$  are more effective than other indicators. Also, there is good separation between wet sand and gas sand for Class I. However, for Class 2 and Class 3 sands, the values decrease and the difference between wet and gas sand decrease, making it very difficult to separate gas sand from wet sand.

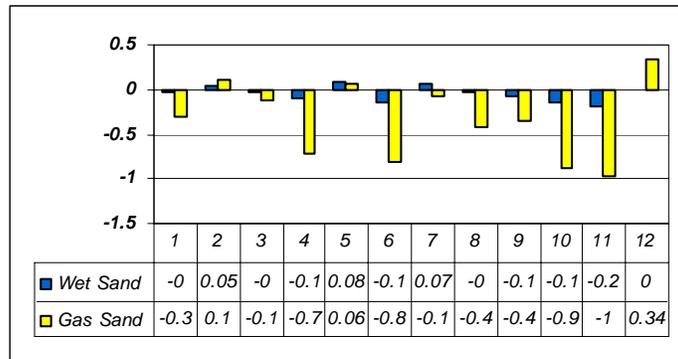


FIG. 3. Calculated indicators for Class 1 sand model (4000-ft depth). 1.  $\Delta\alpha/\alpha$  2.  $\Delta\beta/\beta$  3.  $\Delta\rho/\rho$  4.  $\Delta K$  5.  $\Delta\mu$  6.  $\Delta\lambda$  7.  $\Delta Y$  8.  $\Delta\sigma/\sigma$  9.  $\Delta(Vp/Vs)/(Vp/Vs)$  10.  $K-\mu$  11.  $Ip^2-cl_s^2$ , 12.  $\Delta F$ . Observe that the indicators  $Ip^2-cl_s^2$ ,  $\Delta K$ ,  $K-\mu$  and  $\Delta\lambda$  are more effective than other indicators. Also, there is good separation between wet sand and gas sand.

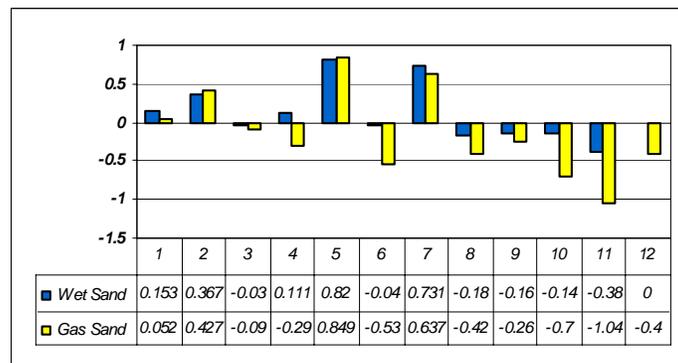


FIG. 4. Calculated indicators for Class 2 sand model (9000-ft depth). 1.  $\Delta\alpha/\alpha$  2.  $\Delta\beta/\beta$  3.  $\Delta\rho/\rho$  4.  $\Delta K$  5.  $\Delta\mu$  6.  $\Delta\lambda$  7.  $\Delta Y$  8.  $\Delta\sigma/\sigma$  9.  $\Delta(Vp/Vs)/(Vp/Vs)$  10.  $K-\mu$  11.  $Ip^2-cl_s^2$ , 12.  $\Delta F$ . Notice that absolute values of  $Ip^2-cl_s^2$ ,  $\Delta K$  and  $\Delta\lambda$  are still more effective than other indicators. However, their values decrease compared with the class 1 sand model, and the difference between wet sand and gas sand decreases.

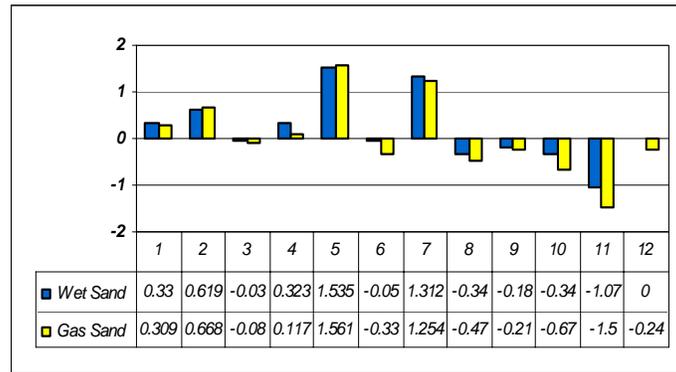


FIG. 5. Calculated indicators for Class 3 sand model (14000-ft depth). 1.  $\Delta\alpha/\alpha$  2.  $\Delta\beta/\beta$  3.  $\Delta\rho/\rho$  4.  $\Delta K$  5.  $\Delta\mu$  6.  $\Delta\lambda$  7.  $\Delta Y$  8.  $\Delta\sigma/\sigma$  9.  $\Delta(Vp/Vs)/(Vp/Vs)$  10.  $K-\mu$  11.  $I_p^2-cI_s^2$  12.  $\Delta F$ . Notice that the values of  $I_p^2-cI_s^2$ ,  $\Delta K$  and  $\Delta\lambda$  become less compared with class 1 and class 2 sand models. It is not easy to separate wet sand from gas sand just from these indicators, except for  $\Delta\lambda$  and  $\Delta F$ .

### Analysis using Castagna and Smith’s examples

Castagna and Smith (1994) presented a set of 25 worldwide measurements of P- and S-wave velocities and densities in associated shales, brine-saturated sands and gas-saturated sand. It is instructive to examine this data set in the cross-plot domain. The generated fluid indicators include  $I_p$ ,  $I_s$ ,  $\lambda$ ,  $\mu$ ,  $\sigma$ ,  $k-\mu$ , and  $I_p^2-cI_s^2$ . Crossplot analysis was conducted to see their discrimination ability between gas-saturated sand, brine-saturated sand and shale.

From the crossplot of  $\mu\rho$  versus  $I_p$  (Figure 6), there are bad separations in both directions between shale, gas sand and wet sand. However, there is good separation in the  $K-u$  direction and bad separation in the  $I_p$  direction between shale, gas sand and wet sand (Figure 7). For the crossplot of  $\lambda\rho$  versus  $\mu\rho$  (Figure 8), there is good separation in the  $\lambda\rho$  direction and bad separation in the  $\mu\rho$  direction between shale, gas sand and wet sand. Also, from Figure 9, there is overlap in the  $ps$  direction. However, there is a good separation in the  $\rho f$  direction between shale, gas sand and wet sand. From the crossplot analysis it can be observed that it is easy to discriminate gas sand from wet sand or shale, but it is hard to separate the wet sand from shale.

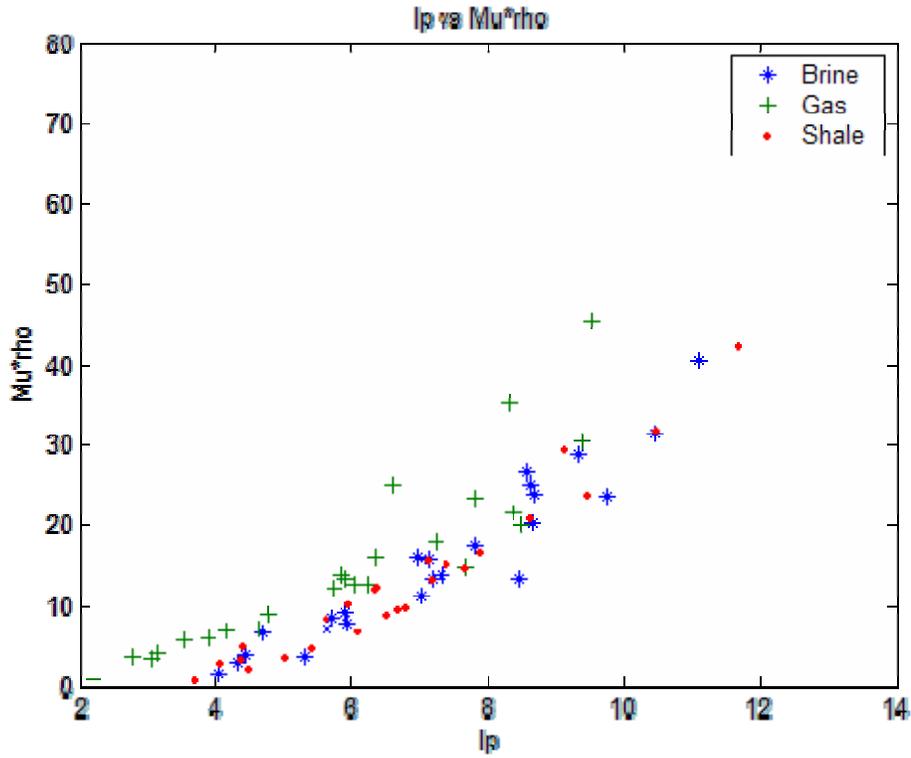


FIG. 6 Crossplot of  $I_p$  versus  $\mu\rho$

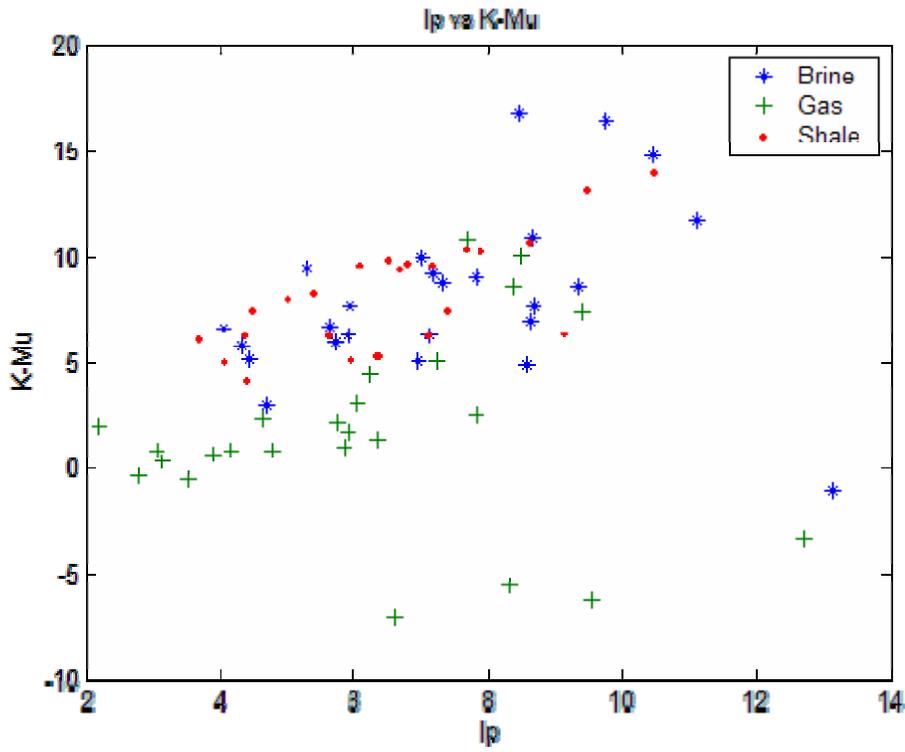


FIG. 7 Crossplot of  $K-u$  versus  $I_p$

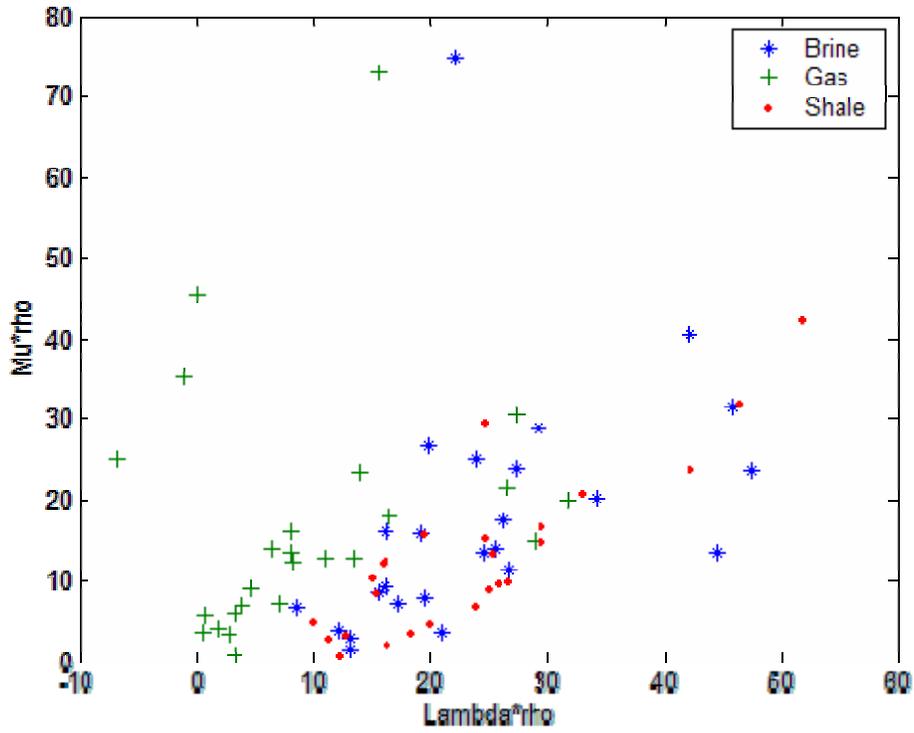


FIG. 8 Crossplot of  $\lambda\rho$  versus  $\mu\rho$

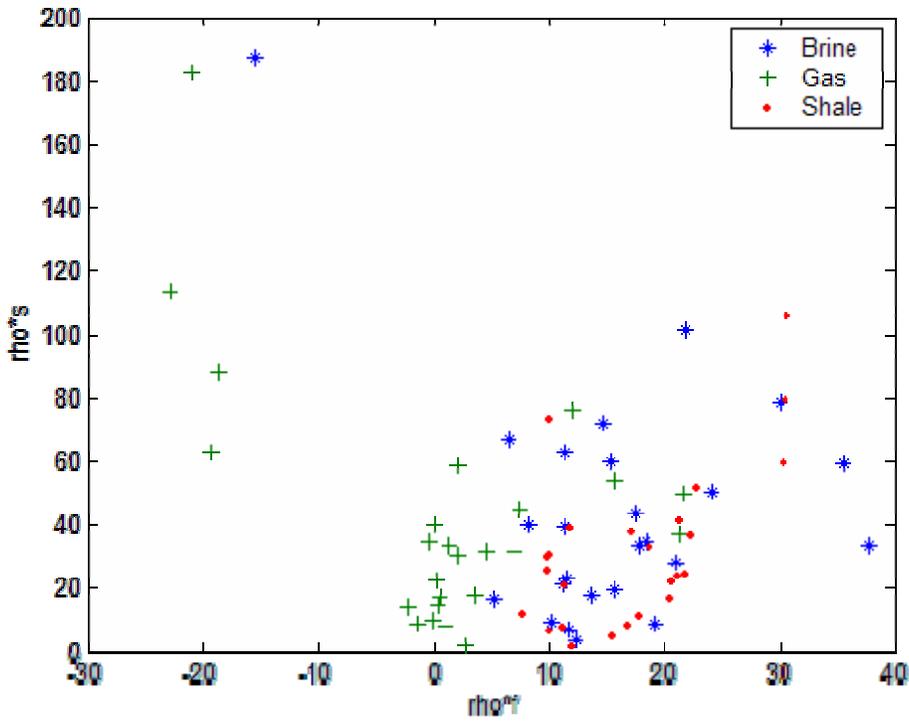


FIG. 9 Crossplot of  $\rho^f$  versus  $\rho^s$

## CONCLUSIONS

Various combinations of rock properties have been proposed as hydrocarbon indicators and it can be concluded that there is a great deal of equivalence between fluid indicators  $I_p^2 - cI_s^2$ ,  $K - \mu$ ,  $\lambda\rho$ ,  $\lambda/\mu$ . For sandstones, the difference  $I_p^2 - cI_s^2$  may be the most sensitive in absolute terms. However, most of these indicators give similar results in magnitude, and each can give insight into the meaning of the other. The best indicator needs to be calibrated and tested for the local situation.

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