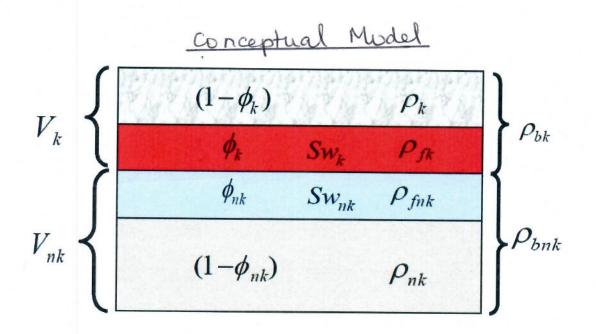
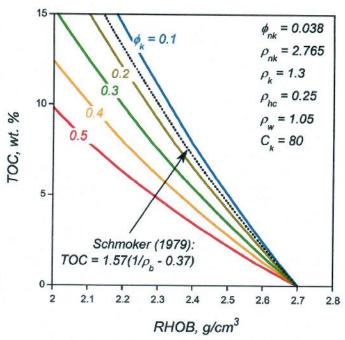
Petrophysical Model for Source Rocks.

In 2012, lev + I published an innovative cocrept "A New Petrophysical Model for organic Shales".

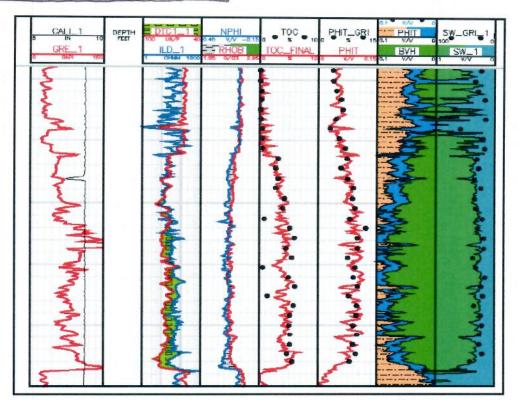
The main idea was hydrocarbons occupied organic porosity and water occupied inorganic porosity. This elemenated the need for Archie type resistivity based approach.



The model decoupled Kerogen porosity that is consistent with thermal maturity and Constraint by physical bounds from total porosity.

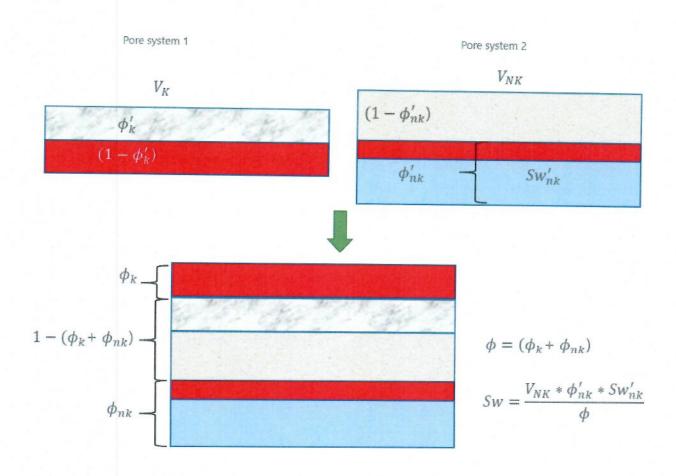


MODEL APPLICATION



while there was consensus about hydrocarbons occupying organic of, the fluids in the host rock was up for debate. Hence this update to the model allowing for bother HC & water to exist in the ineganic porosity.

The New Model (Update)



This model eliminates the issues faced by conventional Archie derivative models. The need for Rw, m, n, a, Rt, Vcl -> no longer is required. Rather two parameters are introduced to obtain a solution.

1. Swik 3 -> when Toc or Keregen = 0. 2: p'nk

There are ways to estimate this as described earlier or some may find creative ways to introduce these parameters. But the good there is this model can be used as a collibration tool to estimate them by calibrating to come data.

MODEL INPUTS

1. $C_k \Rightarrow \% = 0$ org. corbon in Konogen

2. $C_{nk} \Rightarrow density = 0$ inorg. rock

3. $C_k \Rightarrow density = 0$ kerogen

4. $C_{nk} \Rightarrow density = 0$ kerogen

5. $C_{nk} \Rightarrow density = 0$ kerogen

6. $C_{nk} \Rightarrow density = 0$ 6. $C_{nk} \Rightarrow density = 0$ 6. $C_{nk} \Rightarrow density = 0$ 7. $C_{nk} \Rightarrow density = 0$ 8. $C_{nk} \Rightarrow density = 0$ 9. $C_{nk} \Rightarrow density = 0$ 9

Computation.

$$\frac{\partial}{\partial x} = \frac{ay - bx}{cy - bx}$$

Where
$$a = Rm - R_b$$
 $b = Rm - R_{fnk}$
 $c = Rm - R_k$ $x = A \phi'nk$ $A = (1-k)$
 $y = 1 - \phi'nk + A \phi'nk$ $R_{fnk} = Sw'nk (Rw Rnc) + Rnc$

$$\Phi_{n\kappa} = \frac{A \phi_{n\kappa} (1 - \phi_{\kappa})}{(1 - \phi_{nic} + A \phi_{nic})}$$

(6)
$$\phi_{k}' = \frac{\phi_{k}}{\kappa(1-\phi)+\phi_{k}}$$

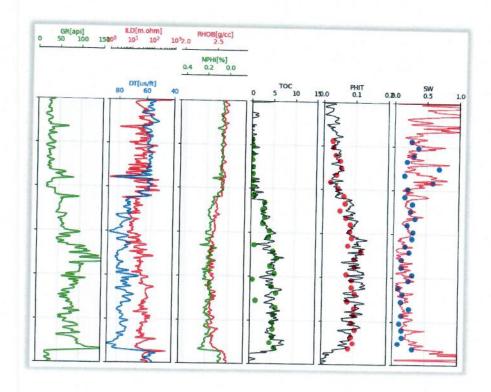
$$Sw = \frac{V_{nK} \phi_{nK} Sw_{nK}}{\phi}, \frac{V_{K} = K(1-\phi)}{(1-\phi)}$$

$$V_{nK} = \frac{V_{nK} - V_{nK}}{(1-\phi)}$$

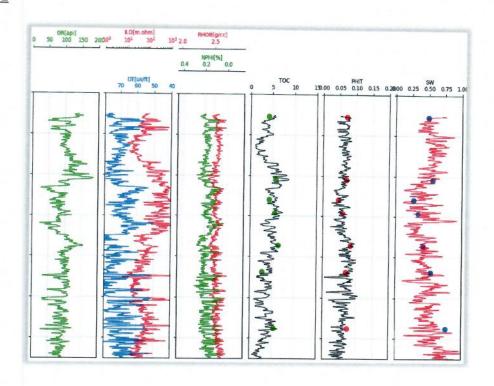
$$V_{nK} = \frac{V_{nK} - V_{nK}}{(1-\phi)}$$

Model Application:

Shale Play 1



Shale Play 2

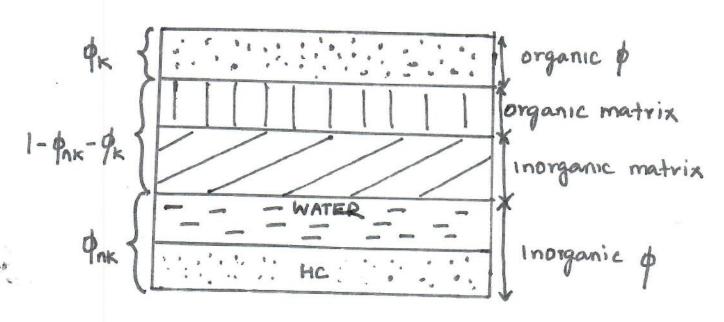


APPENDIX – Model Derivation

SOURCE ROCK PETROPHYSICAL MODEL

UPDATE: HYDROCARBONS IN ORGANIC PORDSITY
WATER + HYDROCARBONS IN INORGANIC POROSITY

Let us consider the total system System



Total porosity $\phi_E = \phi_K + \phi_{nK}$ Lets just take a subset of System (D) of only the matrix portion.

k → volume of kerogen (or any organic matter)

If $\ell_k = density of kerogen and$ $<math>\ell_{nk} = density of morganic rock (usually from XR)$ then from mass balance

expressing in wty.

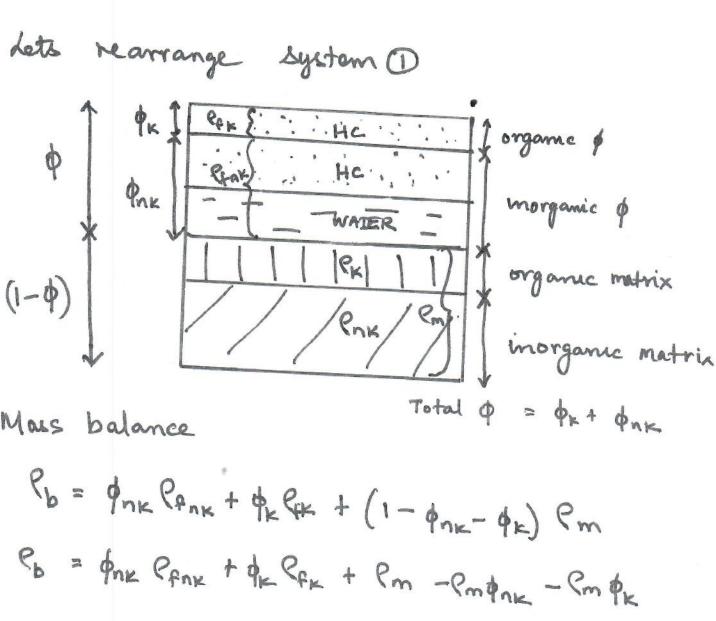
Too is the wty. of organic carbon. If Ck is the organic carbon 1. in kerogen,

can be combined to give.

So after computing Toc from logs,

1. Use egn (3) to compute karogen volume 2. Use Kor Toc to compute Pm from egns (1) or (3)

Pink -> obtained from XRD



The assumptions for fund.

O Kerogen/erg. of contains only HC

Cfk = Swk Pw + (1-Swie) Phc

Swk = 0

Pfk = Phc.

@ Ingry of contains both HC + water.

Penk = Swnk Pw + (1-Swnk) Phc

The "or prime represents values in the individual domain (explained later).

. Pfnx = Swnx (Pw-Pnc) + Phc.

If there is invasion in the system, it is assumed to invade only the th-organic of since org. of is HC wet.

of unvasion.

Pfnk = Sxo Pfile + (1-Sxo) [Swnk (Pw-Pnc) + Phc]

Lets define the domain system porosities Earlier we said

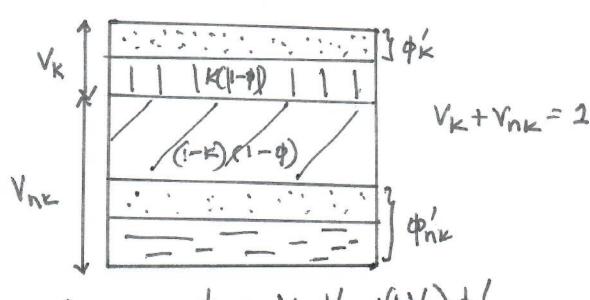
Pt = OK + PNK

of the day are the porosities in their def. of the total system, in essence

φ = organic φ + inorganic φ.

We now define something called domain specific system > denoted by "!"

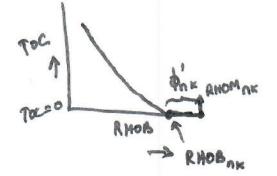
\$k' > kerogen domain \$p >> w.r.t. > V_K \$\delta_{nk}' > In. orgaine domain \$d. >> w.r.t > V_{nk}.

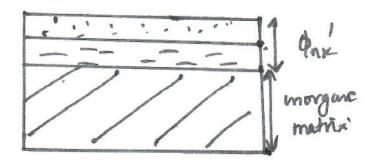


In this system &= VKOK+(1-VK) Ank

Pnk is what we get from cone data as we extrap

TOC = 0 at TOC = 0





(1- p/ + A p/x)

From (8)

$$\phi_{nk} = \frac{A \phi_{nk} (1 - \phi_{k})}{(1 - \phi_{nk} + A \phi_{nk})}$$
 where $A = (1 - k)$

Let
$$a = P_m - P_b$$

$$b = P_m - P_n \times P_n$$

·.
$$\phi_{k} = \frac{a - \phi_{k}b}{c} \Rightarrow \frac{a - bz(1 - \phi_{k})}{c}$$

$$\phi_{k} = \frac{ay - bx + bx \phi_{k}}{cy - bx}$$

$$\phi_{k} = \frac{ay - bx}{cy - bx} \rightarrow 9$$

So after computeries Volume of Kerogen K, matrix density Pm

estra polating TOC > 0 un TOC vs RHOB plot. (or even Toc vs op plot).

Again (extrapolate TOC -> 0 in core TOCKSW plot). Also one can use Vcl and establish a transform for Swink from Vcl.

Then

4 compute of from (9)

Lis once of is computed, compute onk from &

Then total porosity

\$ = PR+ PNK

once total porecity of is computed, we can then solve for Kerogen domain organic of, of.

$$\frac{\partial}{\partial k} = V_{k} \phi_{k}'$$

$$V_{k} = K(1-\phi) + V_{k} \phi_{k}'$$

$$\frac{\partial}{\partial k} = \frac{\partial}{\partial k}' K(1-\phi)$$

$$\frac{\partial}{\partial k} = \frac{\partial}{\partial k}' K(1-\phi)$$

$$\frac{\partial}{\partial k} = \frac{\partial}{\partial k}' \left[K(1-\phi) + \phi_{k} \right]$$

$$\frac{\partial}{\partial k}' = \frac{\partial}{\partial k}' \left[K(1-\phi) + \phi_{k} \right]$$

$$\frac{\partial}{\partial k}' = \frac{\partial}{\partial k}' \left[K(1-\phi) + \phi_{k} \right]$$

So from parevious steps we can solve for

the total swater saturation of the System Sw is given by.

Remember.

$$V_{K} = \frac{K(1-\phi)}{(1-\phi_{K}')} \delta V_{NK} = 1-V_{K}$$

Also

$$V_{k} = \frac{\phi - \phi_{nk}}{\phi_{k'} - \phi_{nk}}$$
(from $\phi = V_{k} \phi_{k'} + (1 - V_{k}) \phi_{nk}$)