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Nuclear Magnetic Resonance (NMR) Outputs Generation for Clastic Rocks Using Multi Regression Analysis, Examples from Offshore Western Australia

Reza Rezaee 

Western Australian School of Mines, Minerals, Energy and Chemical Engineering, Curtin University, Perth, WA 6151, Australia; r.rezaee@curtin.edu.au

Abstract: A large database of nuclear magnetic resonance (NMR) logging data from clastic rocks of offshore oil and gas fields of Western Australia was used to assess the performance of multi regression analysis (MRA) to calculate NMR log outputs from conventional well logs. This short paper introduces a set of MRA equations for the calculation of the NMR log outputs using conventional well logs as inputs. This study shows that unlike machine learning methods the MRA approach fails to predict most of the NMR log outputs with acceptable accuracy but can provide Coates and SDR permeabilities with R2 of more than 0.75.

Keywords: nuclear magnetic resonance (NMR) logs; CMR; clastic rocks; multi regression analysis; permeability



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1. Introduction

The nuclear magnetic resonance (NMR) logging tool uses the response of hydrogen to a magnetic field through T_2 relaxation, which is the process by which the transverse components of hydrogen magnetisation decay [1]. In saturated porous media, the NMR T_2 relaxation rate is governed by the following equation which is a function of individual intrinsic (or bulk), surface, and diffusion relaxation processes [2]:

$$\frac{1}{T_2} = \frac{1}{T_{2B}} + \frac{1}{T_{2S}} + \frac{1}{T_{2D}} \quad (1)$$

where T_{2B} , T_{2S} , and T_{2D} , are the bulk, surface, and diffusion transversal relaxation time in ms, respectively.

Since in geological porous media transversal relaxation is chiefly controlled by surface relaxation, the T_2 relaxation time is related to surface to volume ratio (S/V) [3]:

$$\frac{1}{T_2} \approx \rho \frac{S}{V} \quad (2)$$

where ρ is the transversal surface relaxivity ($\mu\text{m/s}$).

As can be seen in Equation (2), the transverse relaxation rate, $1/T_2$, is proportional to the surface-to-volume ratio (S/V). As a result, the relaxation of hydrogen close to the solid surface is faster than that of those in the free-flowing fluid. Thus, the smaller pores where hydrogens are closer to solid surfaces have a faster relaxation rate than the larger pore spaces. This approach enables the NMR tool to provide some of the rock characteristics such as porosity, pore size distribution, and permeability that are explained in detail in many publications [1,4–12].

The T_2 distribution is generally represented by a set of eight bins porosities (BP1 to BP8) determined by a set of T_2 cutoffs. For clastic rocks, the cutoffs shown in Figure 1 are

generally used to differentiate the bin porosities and therefore to present several types of pore systems as below:

$$CBW = BP1 + BP2 \tag{3}$$

$$BVI = BP3 + BP4 \tag{4}$$

$$FFI = BP5 + BP6 + BP7 + BP8 \tag{5}$$

$$\varnothing_e = \sum_{i=3}^8 BP_i = BVI + FFI \tag{6}$$

$$\varnothing_t = \sum_{i=1}^8 BP_i = CBW + BVI + FFI \tag{7}$$

$$S_{wirr} = \frac{BVI}{BVI + FFI} = \frac{BVI}{\varnothing_e} \tag{8}$$

where BP = bin porosity; CBW = clay bound water; BVI = bulk volume irreducible; FFI = free fluid index; \varnothing_e = effective porosity; \varnothing_t = total porosity; and S_{wirr} = irreducible water saturation.

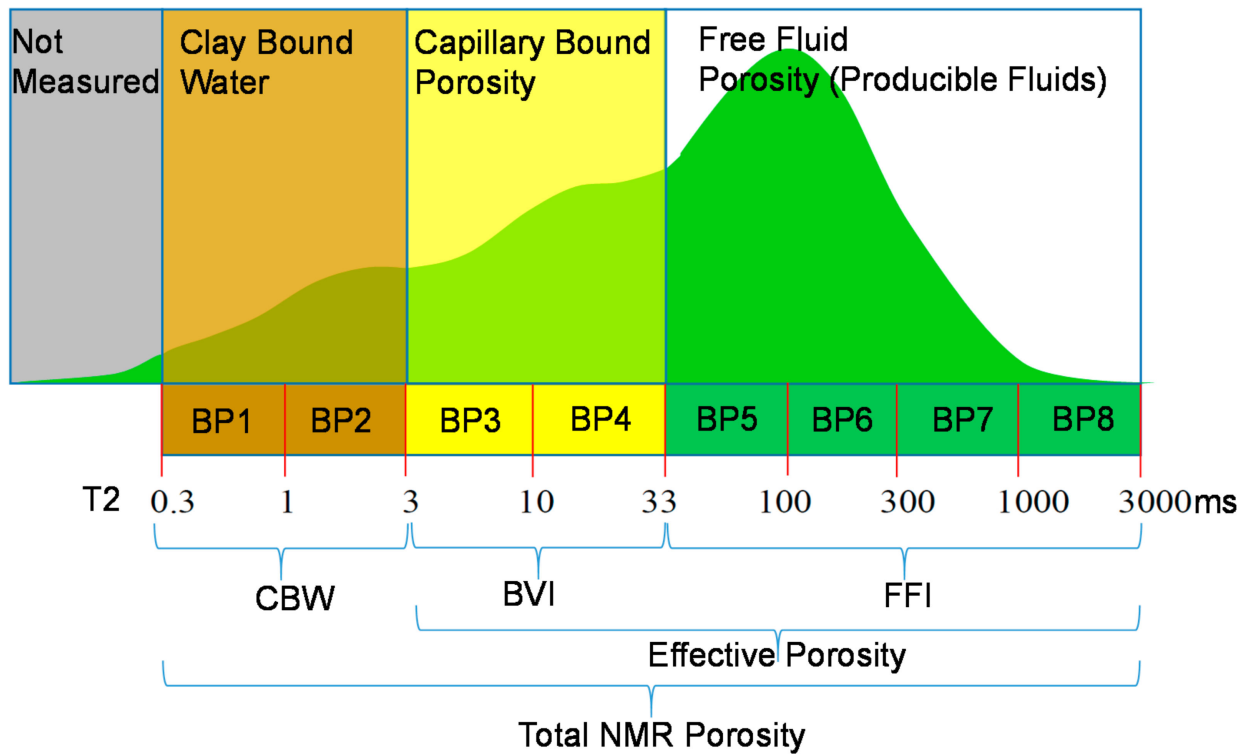


Figure 1. The T_2 distribution shows clay-bound water (CBW), capillary bound fluid (BVI), and free fluid (FFI). These properties are shown in relation to the T_2 signal. From [13].

The pore size distribution of the NMR logging tool has been used to calculate the rock’s matrix absolute permeability. The two most applied equations are Coates (Equation (9)) and the SDR (Schlumberger-Doll Research, Equation (10)) models [10,14–16].

$$k = \left[\left(\frac{\varnothing_e}{C} \right)^2 \left(\frac{FFI}{BVI} \right) \right]^2 \tag{9}$$

$$k = aT_{2LM}^2 \varnothing_e^4 \tag{10}$$

where k = matrix permeability (mD); \varnothing_e = NMR effective porosity (%); C = a constant specific to the formation that reflects the correlation between the rock’s pore throat and pore

size, and in fact, it is a function of pore geometry and is generally considered 10 as a default value for clastic rocks; T_{2LM} = the logarithmic mean of the T_2 distribution, milliseconds; and “ a ” = a coefficient that depends on formation type and has to be determined through calibration with core porosity. “ a ” is generally close to four for sandstone.

This study is an attempt to assess the ability of the multi regression analysis (MRA) method to generate NMR log outputs using conventional well logs as inputs for clastic rocks.

2. Materials and Methods

More than 16,000 data points from 16 formations of WA offshore basins were used to evaluate the performance of linear multi regression analysis (MRA) to calculate NMR logging tool outputs. The data were collected from formations with the age range of Permian to Tertiary from 14 wells with a complete set of CMR and conventional well logs. Wells are located within 3 main offshore basins of WA with a depth range of from 1188 to 4138 m (Table 1).

Table 1. The list of the basins, number of data points for each formation, age, depth interval, and type of lithology used in this study.

Basin	Formation	No of Data Points	Age	Depth (m)	Main Lithology
Browse	Bassett	380	Tertiary	1390–1408	Sandstone
Browse	Grebe	1571	Tertiary	1312–1385	Sandstone
Browse	Nome	167	Triassic	3823–3848	Sandstone
Browse	Plover	295	Jurassic	3776–3823	Sandstone
Browse	Vulcan	1130	Juras. to Cre.	3942–4138	Sandstone
N. Carnarvon	Angel	941	Jurassic	3400–3700	Sandstone
N. Carnarvon	Barrow Group	287	Cretaceous	1890–1936	Sandstone
N. Carnarvon	Brigadier	762	Triassic	3037–3162	Sandstone
N. Carnarvon	Forestier	406	Cretaceous	2983–3147	Claystone
N. Carnarvon	Muderong	145	Cretaceous	2960–2982	Shale
N. Carnarvon	Mungaroo	6868	Triassic	3045–3710	Sandstone
Perth	Cattamarra	697	Jurassic	2940–3052	Sandstone
Perth	Dongara	37	Triassic	1276–1281	Sandstone
Perth	High Cliff	1008	Permian	1310–1476	Sandstone
Perth	IRCM	498	Permian	1278–1475	Sandstone
Perth	Kockatea	833	Triassic	1188–1427	Shale

The well logs used in this study as MRA inputs are density (RHOB, g/cc), Neutron (NPHI, v/v), Photoelectric (PEF, b/e), resistivity (deep, shallow, and very shallow, ohm-m), and sonic (DT, us/ft). Instead of gamma-ray (GR), the volume of shale (V_{sh} , v/v) was used since GR may vary from well to well for the same formation. Effective porosity calculated from the density tool (PHIDE, v/v) was also included to increase MRA accuracy. Table 2 provides a brief explanation for each well log used for this study. Figure 2 shows the variation of well logs inputs.

Table 2. A list of typical well logs and their general applications.

Well Log Type	Applications
Gamma-ray (GR)	a measure of the natural radioactivity of the whole formation near the wellbore mostly to be used to calculate the volume of shale (V_{sh}).
Bulk density (RHOB)	a measure of the bulk density of the formation near the wellbore that can be used to calculate total porosity.
Neutron porosity (NPHI)	a measure of the bulk volume of hydrogen in the formation near the wellbore that provides total porosity.
Sonic (DT)	a measure of the travel time of sound waves in the formation near the wellbore that can be used to calculate porosity.
Photoelectric Factor (PEF)	a measure of photoelectric absorption that depends on the atomic number that can indicate lithology.
Resistivity logs (MSFL, LLS, and LLD)	a measure of the very shallow (MSFL), shallow (LLS), and deep (LLD) resistivities surrounding the tool-including contributions from the mud and the formation around the wellbore.

The volume of shale was calculated for all formations with the following general equation to normalise the GR values:

$$V_{sh} = \frac{GR_{log} - GR_{min}}{GR_{logmax} - GR_{min}} \quad (11)$$

The GR_{min} and GR_{max} values are the values of the sand and shale intervals, respectively, taken from the GR reading for each well.

The effective porosity was calculated by including a correction for the contribution of shale to the log measurements. The density log was used to calculate effective porosity for shaly formations using the following equation:

$$D_{shce} = \frac{ma - b}{ma - f} - V_{sh} \frac{ma - sh}{ma - f} \quad (12)$$

where $\varphi_{D_{shc}}$ = shale corrected density porosity; φ_e = effective porosity; ρ_b = bulk density (g/cc); ρ_f = fluid density (g/cc); ρ_{ma} = matrix density (g/cc); V_{sh} = volumetric fraction of shale; ρ_{sh} = shale bulk density (g/cc) that was identified for each well separately.

MRA outputs are BP1 to BP8, CBW, BVI, FFI, Swirr, T_{2LM} , Coates permeability (kCoates), SDR permeability (kSDR). To increase the performance of MRA, instead of using permeability values the logarithm of permeabilities (LogkCoates and LogkSDR) were used as outputs. Figure 3 shows the variation of major MRA outputs.

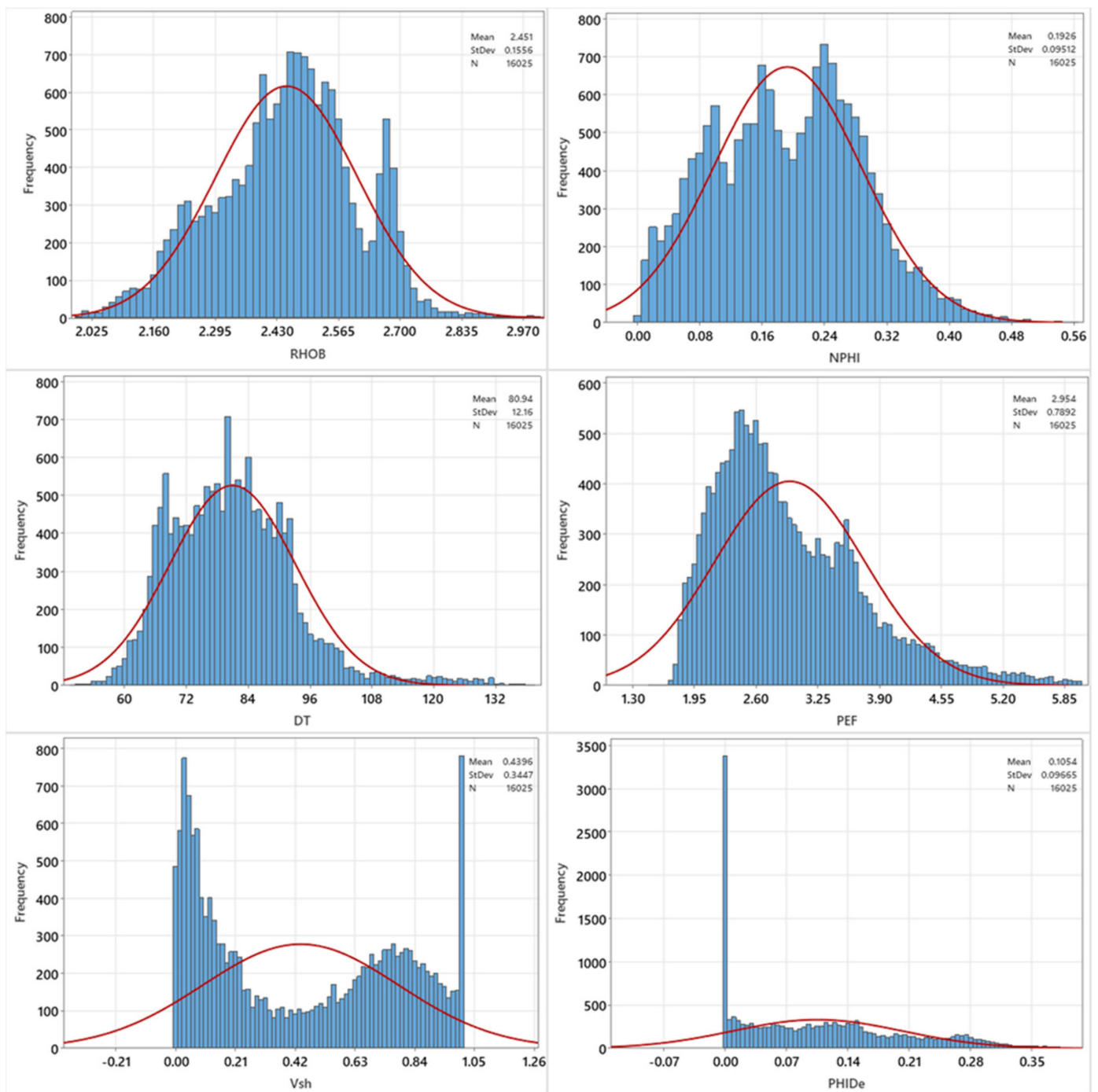


Figure 2. Variations of the major MRA inputs.

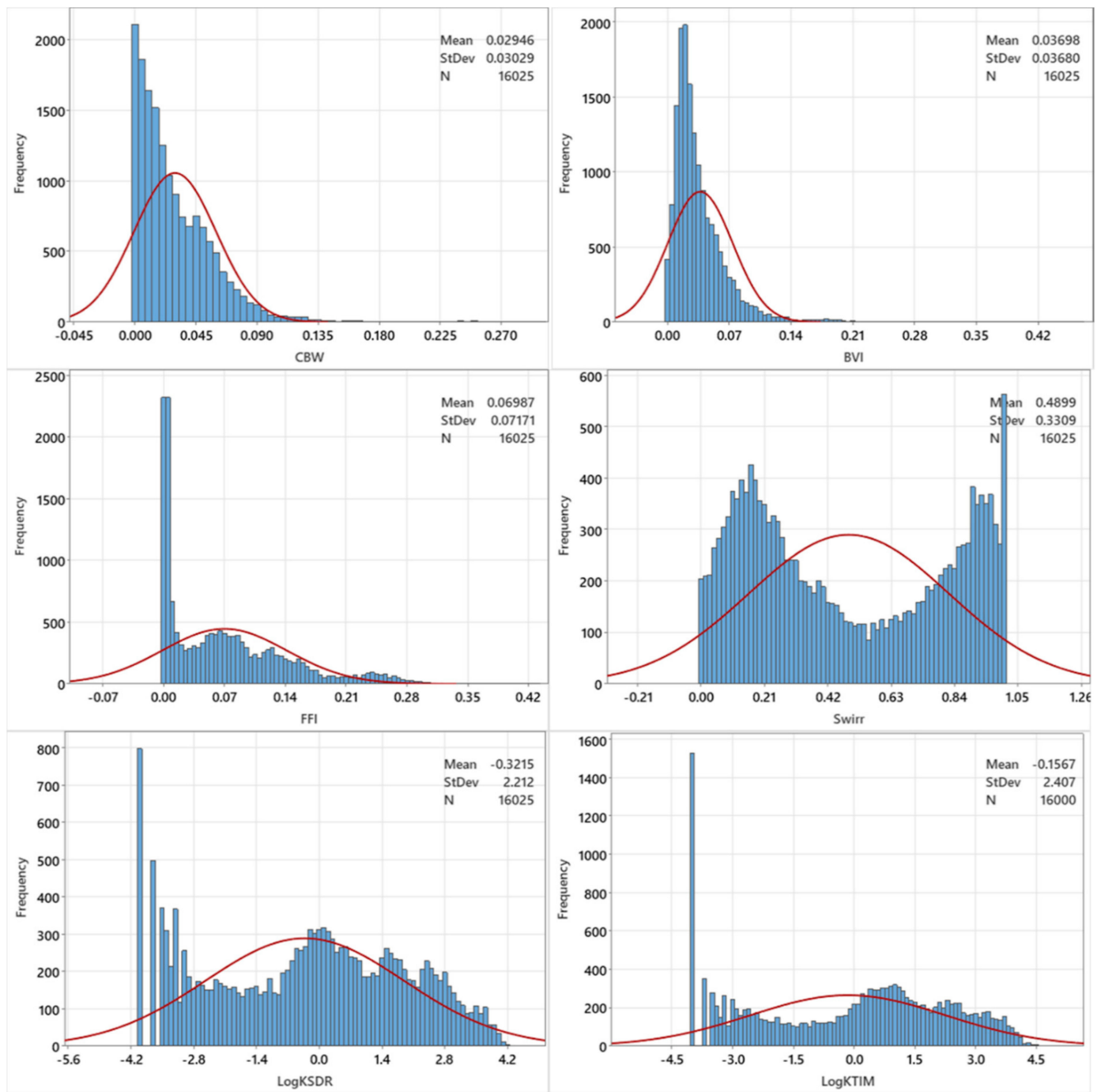


Figure 3. Variations of the major MRA outputs.

Before running the MRA, all well logs were quality controlled and checked for any possible errors. A careful depth match was conducted between NMR and other logs using GR that was available for both NMR and conventional well logs.

3. Multi Regression Analysis Results

At first, all available conventional well logs were used as MRA inputs to predict NMR log outputs separately. This needs to be noted that all bin porosities were also considered as NMR outputs although they were also used in the form of CBW, BVI, and FFI too (see Equations (3)–(5)).

Table 3 shows the generated equations and their corresponding R2 for each output. To optimise the equations by using fewer log inputs with nearly the same accuracy in terms of

R2, inputs with less influence were removed and new equations were generated for each output. Table 4 shows the optimised equations for each output with their R2.

Table 3. MRA equations for the calculation of NMR log outputs. Variables with ignorable influence are removed from the equations.

Equations	R2
BP1 = $-0.02629 + 0.0064RHOB + 0.0353NPHI + 0.00021DT - 0.000135PEF + 0.000004MSFL0 + 0.003167V_{sh} - 0.02648PHIDe$	0.34
BP2 = $0.0809 - 0.04334RHOB + 0.0382NPHI + 0.000475DT + 0.003513PEF + 0.000018LLS + 0.002914V_{sh} - 0.13883PHIDe$	0.46
BP3 = $0.3296 - 0.13588RHOB - 0.03444NPHI + 0.0003DT + 0.010233PEF + 0.00003LLS + 0.001917V_{sh} - 0.22825PHIDe$	0.24
BP4 = $0.27838 - 0.11193RHOB - 0.06877NPHI + 0.00023DT + 0.008134PEF - 0.000038MSFL - 0.003877V_{sh} - 0.14099PHIDe$	0.19
BP5 = $0.12283 - 0.05146RHOB - 0.06293NPHI + 0.000508DT + 0.001021PEF - 0.000040MSFL - 0.014323V_{sh} - 0.05593PHIDe$	0.28
BP6 = $0.06577 - 0.02436RHOB - 0.01841NPHI + 0.000189DT + 0.001628PEF - 0.000082MSFL - 0.018688V_{sh} + 0.07862PHIDe$	0.46
BP7 = $-0.09329 + 0.0464RHOB + 0.14057NPHI - 0.00052DT - 0.000217PEF - 0.00008LLS - 0.02807V_{sh} + 0.2707PHIDe$	0.53
BP8 = $-0.10033 + 0.04855RHOB + 0.11431NPHI - 0.000431DT - 0.00247PEF + 0.000051MSFL - 0.000072LLS - 0.016974V_{sh} + 0.2044PHIDe$	0.41
CBW = $0.0546 - 0.03695RHOB + 0.0735NPHI + 0.000686DT + 0.003378PEF + 0.000018LLS + 0.00608V_{sh} - 0.16531PHIDe$	0.45
BVI = $0.608 - 0.24781RHOB - 0.10321NPHI + 0.00052DT + 0.018366PEF - 0.00006MSFL + 0.00004LLS - 0.00196V_{sh} - 0.36923PHIDe$	0.21
FFI = $-0.005 + 0.01915RHOB + 0.17356NPHI - 0.000255DT - 0.00004PEF - 0.000072MSFL - 0.000128LLS + 0.000026LLD - 0.07806V_{sh} + 0.4978PHIDe$	0.72
Swirr = $-0.446 + 0.1997RHOB + 0.054NPHI + 0.00363DT + 0.0157PEF - 0.000166MSFL + 0.000547LLS - 0.000167LLD + 0.4621V_{sh} - 1.064PHIDe$	0.68
T2LM = $-584.1 + 270.4RHOB + 396.6NPHI - 1.15DT - 14.43PEF + 0.1854MSFL - 0.3293LLS + 0.0729LLD - 111.75V_{sh} + 1020.4PHIDe$	0.48
LogkSDR = $17.087 - 6.496RHOB + 2.3NPHI - 0.01037DT + 0.1473PEF + 0.0003MSFL - 0.004737LLS + 0.001246LLD - 3.6307V_{sh} + 1.220PHIDe$	0.76
LogkCoates = $18.046 - 6.608RHOB + 1.734NPHI - 0.01196DT + 0.1093PEF - 0.001087MSFL - 0.003939LLS + 0.000988LLD - 3.9892V_{sh} + 1.193PHIDe$	0.75

Table 4. Optimised MRA equations for the calculation of NMR log outputs.

Equations	R2
BP1 = $-0.02573 + 0.00617RHOB + 0.03409NPHI + 0.00021DT + 0.003236V_{sh} - 0.0264PHIDe$	0.34
BP2 = $0.08302 - 0.04386RHOB + 0.04394NPHI + 0.000475DT + 0.003567PEF + 0.00002LLS - 0.14572PHIDe$	0.46
BP3 = $0.33195 - 0.13671RHOB - 0.03013NPHI + 0.00029DT + 0.010242PEF + 0.000014LLS - 0.23291PHIDe$	0.24
BP4 = $0.27469 - 0.11082RHOB - 0.07688NPHI + 0.000231DT + 0.008091PEF - 0.000034MSFL - 0.13152PHIDe$	0.19
BP5 = $0.11091 - 0.04627RHOB - 0.06055NPHI + 0.000520DT - 0.000037MSFL - 0.013959V_{sh} - 0.05024PHIDe$	0.28
BP6 = $0.06874 - 0.02397RHOB + 0.000122DT + 0.001399PEF - 0.000056MSFL - 0.021843V_{sh} + 0.07788PHIDe$	0.46
BP7 = $-0.08797 + 0.04412RHOB + 0.14170NPHI - 0.000523DT - 0.000035LLS - 0.02870V_{sh} + 0.26720PHIDe$	0.53
BP8 = $-0.10246 + 0.04955RHOB + 0.11239NPHI - 0.000433DT - 0.00243PEF - 0.000041LLS - 0.0165V_{sh} + 0.20577PHIDe$	0.41
CBW = $0.0539 - 0.03664RHOB + 0.07289NPHI + 0.000685DT + 0.003393PEF + 0.000023LLS + 0.00624V_{sh} - 0.16483PHIDe$	0.45
BWI = $0.6095 - 0.24855RHOB - 0.10173NPHI + 0.000523DT + 0.018308PEF - 0.00244V_{sh} - 0.37009PHIDe$	0.21
FFI = $-0.004 + 0.01877RHOB + 0.17355NPHI - 0.000255DT - 0.000103MSFL - 0.000073LLS - 0.0782V_{sh} + 0.4968PHIDe$	0.72
Swirr = $0.4566 + 0.1991RHOB + 0.003828DT + 0.01645PEF + 0.000195LLS + 0.47167V_{sh} - 1.0569PHIDe$	0.68
T _{2LM} = $-590.3 + 273.3RHOB + 392.1NPHI - 1.154DT - 14.41PEF - 0.1411LLS - 110.79V_{sh} + 1023.5PHIDe$	0.48
LogkSDR = $17.075 - 6.490RHOB + 2.294NPHI - 0.01038DT + 0.1475PEF - 0.004542LLS + 0.001190LLD - 3.6281V_{sh} + 1.227PHIDe$	0.76
LogkCoates = $18.09 - 6.629RHOB + 1.774NPHI - 0.01192DT + 0.1085PEF - 0.004654LLS + 0.001192LLD - 3.9988V_{sh} + 1.166PHIDe$	0.75

To assess the performance of the MRA method for clay-rich and clay-poor lithologies, well logs were separated based on clay content. The dataset was classified based on V_{sh} ($V_{sh} < 50\%$ and $V_{sh} > 50\%$), to separate shale from shaly sand. A comparison of the results showed that generally there were no improvements in the MRA performance, and even, the performance was reduced for all of the outputs. Perhaps the performance deterioration could be due to the reduction of the number of data points.

4. Discussion and Conclusions

In this study, the MRA method was used to generate the NMR tool's outputs using a large database of conventional well logs. Figure 4 shows the success of linear MRA in providing NMR log outputs based on data in Table 4. As it can be seen from the figure, based on the coefficient of determination (R^2), the SDR and Coates permeabilities are successfully calculated followed by FFI and Swirr. The prediction of the remaining parameters with R^2 less than 0.6 is statistically unreliable.

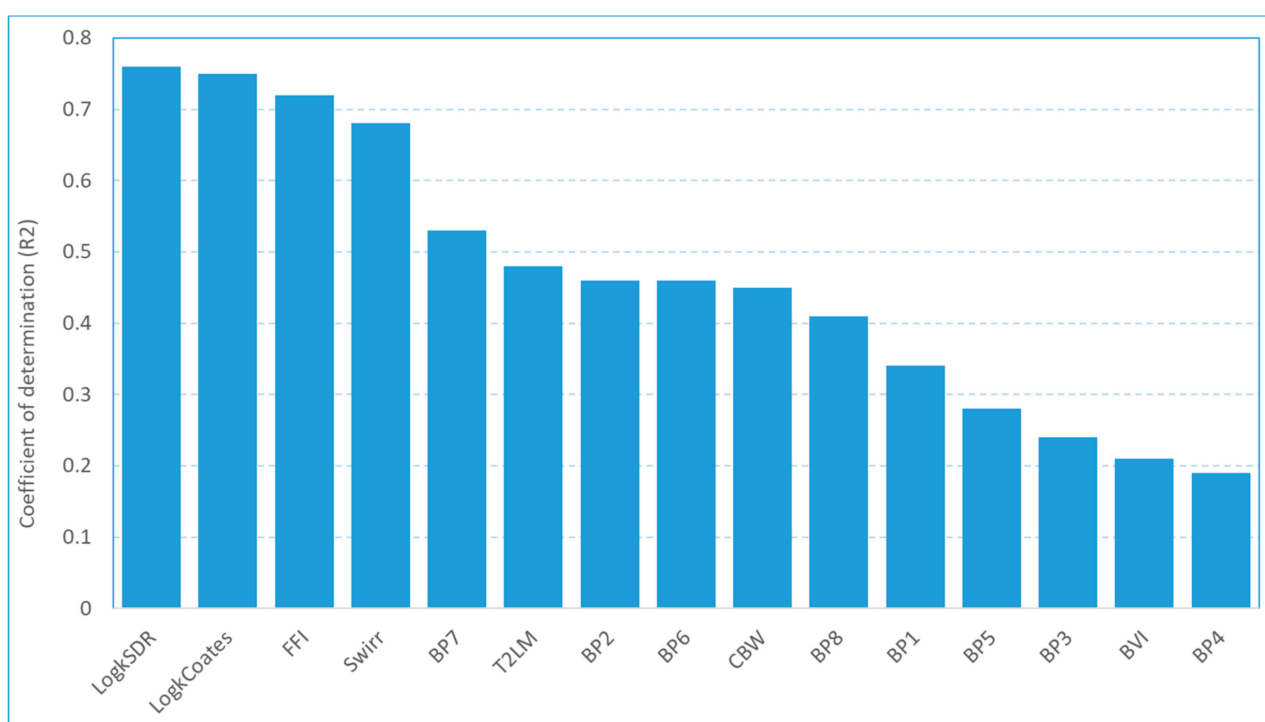


Figure 4. The coefficient of determination (R^2) or the success of MRA in predicting NMR log outputs.

The nature of the conventional logging tools is a way that each tool targets a specific property of rock and its pore fluid. GR detects natural emission of gamma radiation that in sedimentary rocks is mostly sourced from clays and therefore the tool is generally used for V_{sh} calculation. The density tool, RHOB, provides the bulk density of the rock based on the GR Compton scattering process and it is used to calculate total porosity. Neutron tool, NPHI, provides total porosity based on the hydrogen content that is believed to mostly reside in pore filling fluids. Sonic tool, DT, measures sound waves velocity that can be used to calculate porosity. Photoelectric tool, PEF, uses cross-sectional capture of low energy GR and thus can provide lithology. Resistivity tools measure the electrical resistivity of the rock which is the function of several parameters including pore water resistivity, formation factor, and fluid saturation. Therefore, the major application of resistivity data is to calculate fluid saturation. In shaly formations where clay is present the interpretation of well logs becomes complex. For example, in shaly formations all porosity tools (neutron, density, and sonic tools) overestimate porosity. In addition, clays create excessive conductivity due to their cation exchange capacity (CEC) and therefore result in reducing resistivity logs readings. This makes the relation between conventional logs and NMR more intricate.

On the other hand, the NMR logging tool uses hydrogen T_2 relaxation time and provides pore size distribution based on some empirical T_2 cutoffs. BP1 and BP2 make up a part of pore space that is occupied by CBW. BP3 and BP4 provide capillary pore spaces where immobile fluids exist and are named BVI. BP5 to BP8 provide large pore spaces where mobile pore fluid resides and is called FFI. This ability of the NMR tool to differentiate different pore spaces is out of reach of conventional logging tools where porosity tools such as RHOB, NPHI, and Sonic can just provide mostly total porosity regardless of their sizes. Such a difference between NMR and conventional logging tools could be the reason that they fail to predict bin porosities (BPs) accurately. On the other hand, it can be seen in Tables 3 and 4 that permeability, FFI, and Swirr are the only NMR outputs that are calculated with relatively high accuracy. The key point here is that conventional logging tools provide total porosity which is naturally close to FFI which is the dominant porosity in sandstone reservoirs (see Figure 3), the major rock type in this study. Consequently, conventional logging tools fail to differentiate and predict CBW and BVI. Since for NMR's permeability and Swirr, FFI plays an important controlling factor consequently conventional logging data can predict them with higher accuracy.

The same data set used in this study was used by the author to assess machine learning methods for generating NMR log outputs [13]. The study showed that Adaptive Boosting can predict CBW, FFI, permeability, T_{2LM} , and S_{Wirr} with an R2 of more than 0.9. This indicates that, unlike machine learning models that can solve complex relations between inputs and output, linear MRA fails to perform well where the relation between inputs and outputs are complex.

In conclusion, this study indicates that the MRA can be utilised to generate NMR permeability with relatively high accuracy from conventional logging tools for clastic rocks.

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