

INFERRING INJECTION RETURNS FROM CHLORIDE MONITORING DATA

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ABSTRACT

Reservoir chloride concentration and injection rate data have been used to identify injection return flow paths in the reservoir. General trends of chloride and injection rate with time were isolated from their respective short-term variations using the wavelet transformation approach. Multiple regression techniques were then used to correlate the isolated short-term variations in chloride with corresponding short-term fluctuations in injection rates and subsequently to quantify the degree of connectivity between injectors and producers. Data from Palinpinon-I field in the Philippines were analyzed using the outlined method and results were verified successfully against tracer test data and qualitative field observations.

Results of analysis of chloride and injection rate data from Dixie Valley field in Nevada illustrated that multilinear modeling is not suitable for analyzing data sets that lack sufficient time variability or "texture".

1.0 INTRODUCTION

Traditionally, tracer tests are used to establish the degree of connectivity between wells. However, for wells that are only weakly connected these tests may need to be conducted over long periods of time using huge amounts of tracer of sufficient stability to obtain meaningful data. In such cases tracer tests can be too costly and impractical.

On the other hand, there are substances occurring naturally in the reservoir that can behave as tracers. One such substance is chloride. In Palinpinon geothermal field in the Philippines, some injectors and producers are strongly connected so that changes in injection rates result in corresponding increase or decrease in chloride concentrations measured in production wells. Data from one such injector-producer pair in Palinpinon is shown in Fig. 1. The magnitude of the changes in chloride concentration thus reflects the degree of communication between wells. Moreover, Chloride is stable, reasonably conservative and it is free. Therefore, we may be able to extract the same, if not more, information from chloride data as we could from traditional tracer tests and at lower cost.

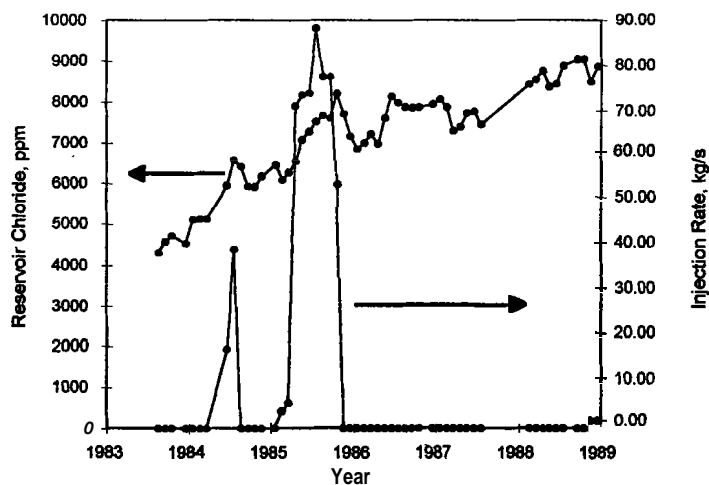


Figure 1. Example chloride and injection data from Palinpinon-I.

The following sections summarize how the method of wavelets and multiple regression techniques were used to analyze chloride and injection data and consequently, identify injection return flow paths; and, how the permeability of these paths were ranked by quantifying the degree of connectivity between injectors and producers.

20 PRELIMINARY LINEAR MODELS

As part of an optimization problem, an earlier work by Macario (1991) proposed several correlations for modeling the reservoir chloride and applied these models to data from Palinpinon-I geothermal field in the Philippines. Of the models tested by Macario (1991), the linear combination model came closest to reproducing field observations. In the first phase of this project therefore, we chose to expand on that model and test it further.

Following is the original linear Combination model proposed by Macario (1991):

$$Cl_p = a_0 + a_1Q_{I1} + a_2Q_{I2} + a_3Q_{I3} + \dots + a_nQ_{In} \tag{1}$$

Based on this model, the strength of the connection between the modeled producer P and an injector I_i is assessed by the magnitude of the coefficient, a_i, of that injector in the model; high values of a correspond to strong connections.

Aside from the extent of reinjection fluid returns, other factors could also affect the chloride level in the reservoir. Extensive boiling and steam separation within the reservoir and natural recharge of higher mineralized fluid are processes that could increase chloride concentrations (Harper and Jordan, 1985). The first process, boiling and steam separation, is a natural reservoir response to exploitation. The chloride concentration may therefore be expected to increase with time as the reservoir is produced. To model this variation with time, a linear time term was added to model (1), thus:

$$Cl_p = a_0 + a_1Q_{I1} + a_2Q_{I2} + a_3Q_{I3} + \dots + a_nQ_{In} + bt \tag{2}$$

More than anything, it was simplicity that guided our choice of the form (linear) of the time term. Solution saturation limits could be expected to put a cap on the maximum chloride concentration and cause it to level off late in the life of the reservoir. For practical purposes, however, we assumed that the chloride concentrations being modeled were far from the maximum limit and increased linearly with time. The question of how chloride concentration in the reservoir actually varies with respect to time will be addressed in more detail in a later section.

We also hypothesized that the reinjection returns' effect on reservoir chloride is governed not just by the rates of injection but also by the chloride concentration of the reinjected fluid. Hence, we have proposed the following modification to model (2):

$$Cl_p = a_0 + a_1Q_{I1}Cl_{I1} + a_2Q_{I2}Cl_{I2} + a_3Q_{I3}Cl_{I3} + \dots + a_nQ_{In}Cl_{In} + bt \tag{3}$$

The additional parameter Cl_{I_n} refers to the chloride concentration of the fluid being injected to injector I_n.

2.1 Results and Discussion

The original and extended models were applied to analysis of both the data set from Palinpinon-I previously used by Macario (1991) and another data set from the Dixie Valley field in Nevada. Qualities of the fit to the data were assessed by inspecting both the calculated values of the multiple regression coefficient, R², and plots of model predicted chlorides against actual data. The multiple regression coefficient, R², represents the proportion of variation in the modeled variable (in this case, chloride concentration) that is predictable from the model. It is therefore desirable to have high values of R². Only the model which best fitted the data or equivalently, had the highest value of R² was subjected to further tests.

For any model to be considered relevant it was deemed necessary that that model be able to account for variations in chloride at any time interval in the data set regardless of which portion of the data set was used to

calculate the linear coefficients. Thus we assessed model relevance by examining how well the model predicts later chloride measurements using coefficients that were calculated from earlier portions of the data set.

The following section discusses the results of application of models (1), (2), and (3) to the Dixie Valley and Palinpinon-I data sets. Model (3) was not used to analyze the Palinpinon-I data set due to the lack of injectate chloride data from that field

Dixie Valley Case

At Dixie Valley, injection rates were recorded daily while chloride concentrations were measured much less frequently, thus, it was the amount of chloride data that set the limit to the number of data points (cases in which chloride concentrations and injection rates were available simultaneously) used for regression.

Table 1 lists R^2 values for models (1), (2), and (3) obtained for each production well. Except for wells 27-33 and 28-33, model (2) gave the highest value of R^2 for all production wells. Addition of the time term to model (1) resulted in 2 to 35% increase in R^2 while inclusion of injectate chloride concentration in model (3) did not result in any significant change in R^2 values. Figure 2 shows the effect of a 35% difference in R^2 on the quality of data fit for well 84-7; also, it illustrates the very minor effect that the injectate chloride term had on the quality of the match. Based on these results we chose to subject model (2) to further testing.

Subsequently, the last six data points were excluded from the regression. Model (2) was then used to predict these values using the coefficients calculated based on the truncated data set. Figure 3 plots the results of the truncated series analysis for well 27-33 which had a 9% maximum deviation of predicted chloride from actual data - the highest deviation observed among all the production wells. Other wells had as little as 1% deviation (Fig. 4).

Inspection of the calculated coefficients revealed one possible reason for the relatively good predictive capacity displayed by model (2) (see Table 2). For this data set, the time term dominates the correlation; the coefficient of the time term is several orders of magnitude (3 to 5, even 8 times!) greater than the injection rate coefficients. This discrepancy was enough to render the injection rate terms trivial: excluding an injection rate term from model (2) resulted in only tiny changes in the quality of the data fit. Figure 5 shows the chloride match for well 74-7 when the chloride is predicted using model (2) but with injection rate term corresponding to injector INJ5218 excluded. That the

Table 1. R^2 values for Dixie Valley wells.

	R^2		
27-33	0.917	0.963	0.965
28-33	0.852	0.936	0.940
4533	0.935	0.970	0.966
63-7	0.826	0.828	0.815
73-7	0.774	0.952	0.952
74-7	0.755	0.968	0.967
76-7	0.930	0.947	0.943
82-7	0.764	0.969	0.967
84-7	0.716	0.978	0.978

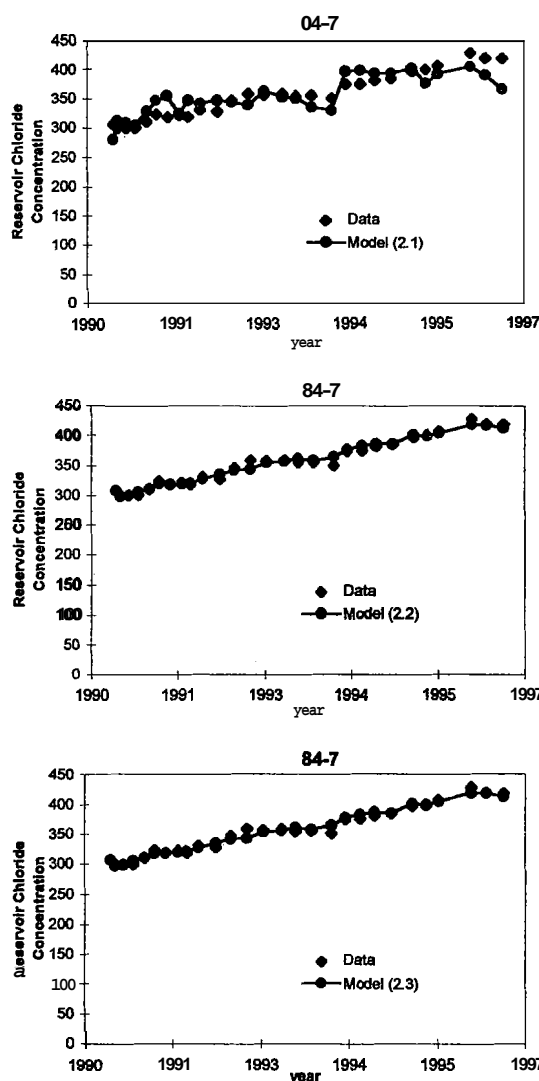


Figure 2. Predicted vs. measured chloride concentration for well 84-7, Dixie Valley.

Table 2. Model (2) coefficients for Dixie Valley production wells.

Model Parameter	Production Well									
	27-33	28-33	45-33	63-7	73-7	74-7	767	82-7	84-7	
a_0	314.58	354.39	271.46	283.77	272.43	318.32	384.12	254.39	271.60	
Injection wells:										
INJ255	8.61E-05	-4.84E-04	5.06E-04	1.80E-03	2.94E-04	1.51E-05	3.60E-04	-3.05E-06	3.99E-04	
INJ455	-2.20E-04	-5.56E-04	3.33E-04	1.33E-03	2.70E-04	-1.62E-04	-1.37E-04	7.57E-05	3.89E-05	
INJ3218	-1.21E-03	-4.30E-04	-1.54E-03	2.69E-03	2.12E-04	-5.89E-04	-3.59E-04	6.67E-04	-1.24E-03	
INJ4118	4.58E-04	6.02E-04	3.54E-04	-4.40E-04	4.35E-05	7.31E-05	2.87E-04	-2.31E-04	9.76E-07	
INJ5218	-1.20E-03	-7.39E-04	-1.18E-03	-2.95E-04	4.37E-04	-2.50E-04	-5.42E-04	7.43E-04	-1.42E-04	
INJ6518	3.56E-03	3.40E-03	2.65E-03	-2.77E-03	-6.31E-04	9.02E-04	-5.63E-04	-3.33E-04	2.07E-03	
INJ_SWL1	-3.99E-03	-3.74E-03	-3.33E-03	8.45E-04	5.39E-04	1.68E-04	4.18E-04	6.27E-04	-7.24E-04	
INJ_SWL3	1.49E-03	9.69E-04	1.33E-03	-4.17E-03	-1.35E-03	-2.18E-04	-3.49E-05	-1.43E-03	-3.27E-04	
t	8.48	10.37	7.61	2.30	17.44	12.95	4.82	26.25	19.04	

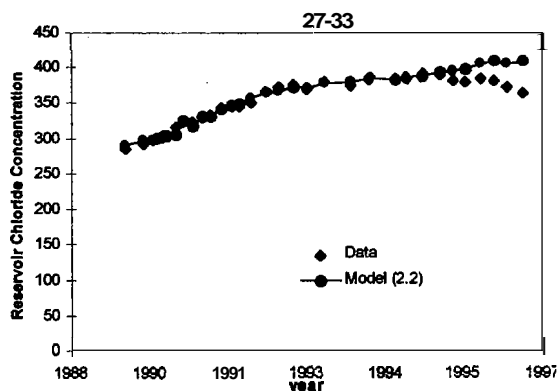


Figure 3. Predicted vs. measured chloride concentration for well 27-33, Dixie Valley; model (2) coefficients calculated with last six data points excluded.

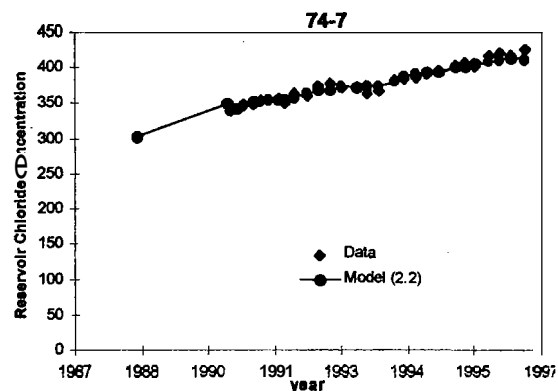


Figure 4. Predicted vs. measured chloride concentration for well 74-7, Dixie Valley; model (2) coefficients calculated with last six data points excluded.

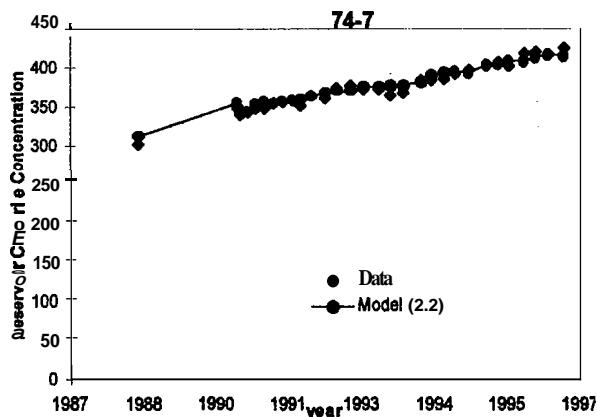


Figure 5. Predicted vs. measured chloride concentration for well 74-7, Dixie Valley; predicted values calculated using model (2) with one injection rate term excluded

injection rate terms are inconsequential to chloride prediction was also evident from inspection of the chloride data: for the most part, the chloride increased linearly with time and response to changes in injection rates was not readily evident. Hence, once the variation of chloride with time was captured in the analysis of the early portion of the data set and there was little deviation observed in the succeeding predictions. It was noted however, that although the deviations were small some of them showed a tendency to increase (Fig. 3). This was true for wells whose chloride ceased at some point to vary linearly with time.

At this point, it is worthwhile to recall that the goal of this project was not prediction but rather, correlation. Although for this specific data set model (2) matched and predicted chloride data relatively well, the dominance of the time term rendered the injection rate coefficients meaningless and ultimately made this model unsuitable for comparing the effects of injection wells on production wells in the field being considered here.

The preceding results lead us to conclude that for the purpose we have set for this project, multiple regression is not a suitable analysis tool for chloride data sets which lack texture.

Polinpinon-I Case

In this case, injection rate data were available as monthly average values; thus, chloride data were converted to monthly average values prior to analysis. As with the Dixie Valley data set, the amount of chloride data set the limit on the number of data points used for regression. Only the portion of the data set from 1983 to 1989 was initially available for use in the initial inspection of the linear models; thus, the following results pertain to the analysis of that early portion of the data set.

The effect on R^2 of adding the time term to model (1) was even more drastic for the Palinpinon-I data set: a maximum increase of 80% in R^2 was observed (Table 3). The effect of a 60% increase in R^2 on the quality of the match for well PN-16D is shown in Fig. 6.

As was done previously in the analysis of the Dixie Valley data set, in the subsequent regression using model (2) the last six points in the chloride series were not considered in the calculation of the linear coefficients. The excluded chloride values were then predicted using the coefficients calculated based on the truncated data set. Deviations of

Well Name	R^2	
	Model (2.1)	Model (2.2)
OK-7D	0.783	0.956
OK-9D	0.717	0.902
OK-10D	0.490	0.535
PN-15D	0.824	0.993
PN-16D	0.606	0.964
PN-17D	0.519	0.939
PN-18D	0.718	0.930
PN-19D	0.559	0.903
PN-23D	0.736	0.958
PN-24D	0.706	0.895
PN-26D	0.728	0.922
PN-27D	0.696	0.944
PN-28D	0.643	0.895
PN-29D	0.817	0.948
PN-30D	0.710	0.832
PN-31D	0.625	0.946

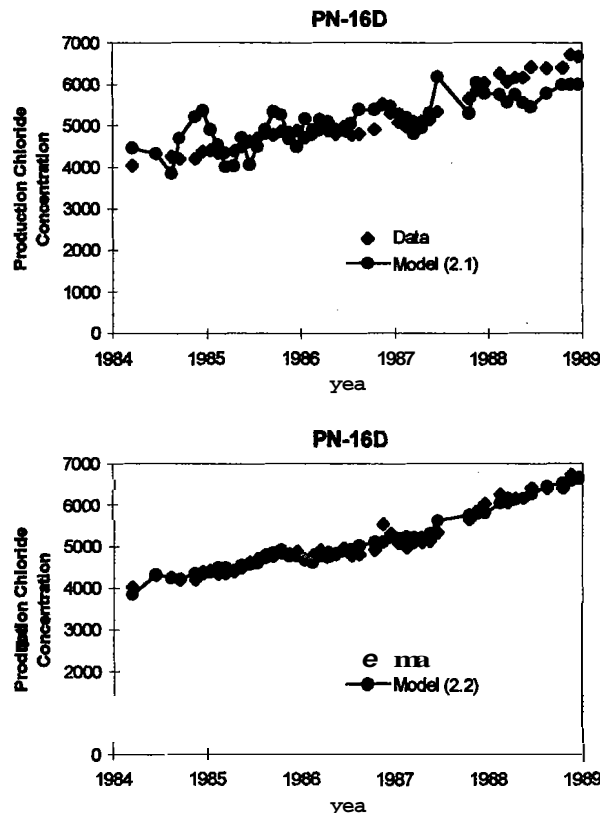


Figure 6. Predicted vs. measured chloride concentration for well PN-16D, Palinpinon-I.

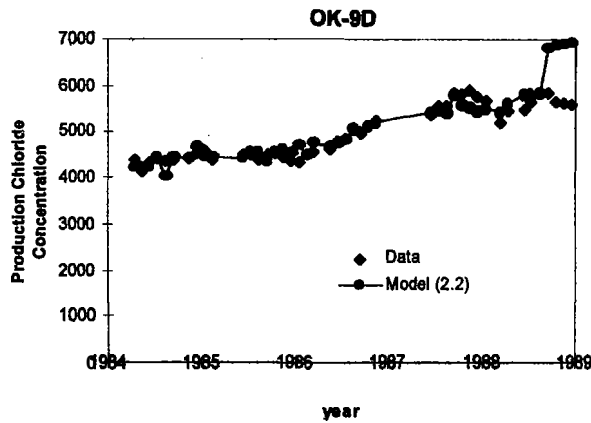


Figure 7. Predicted vs. measured chloride concentration for well OK-9D, Palinpinon-I; model (2) coefficients calculated with last six data points excluded.

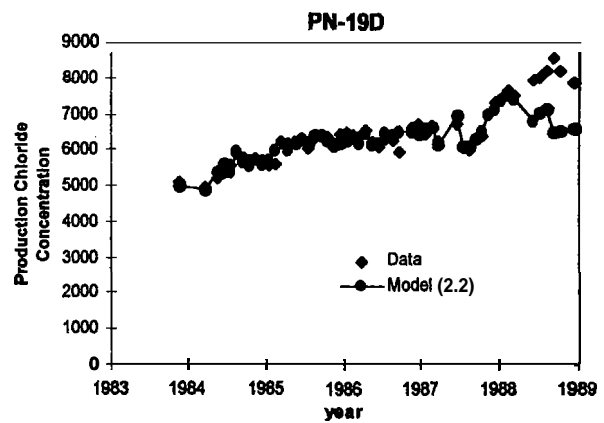


Figure 8. Predicted vs. measured chloride Concentration for well PN-19D, Palinpinon-I; model (2) coefficients calculated with last six data points excluded.

predicted chloride values from actual data for the Palinpinon-I data set were relatively high compared to those of Dixie Valley with a maximum of about 20%. Model (2) overpredicted the data for well OK-9D (Fig. 7) and underpredicted the chloride for well PN-19D (Fig. 8).

As with the Dixie Valley data set, the increasing deviations may be explained by the fact that the linear form of the time term does not account properly for the general trend in chloride with time. Moreover, the relatively high values of the deviations suggest that the injection rate terms contribute significantly to the model but that their contribution has not been assessed adequately.

Table 4 shows that the time term coefficients for this data set are only one to two orders of magnitude higher than the injection rate coefficients, as compared to five to eight orders of magnitude in the Dixie Valley data set. This is due to the more textured nature of the Palinpinon-I data; superimposed on the general increasing trend in chloride are marked dips and bumps. Since the dips and bumps which are accounted for by the injection rate terms are of substantial magnitude, the injection rate coefficients had high absolute values compared to those calculated for the relatively untextured Dixie Valley data set.

It is also important to note that contrary to expectation, some of the injection rate coefficients had negative values. This implies that the operation of injection wells corresponding to those negative coefficients would actually lessen the percentage of injectate being produced. One explanation is that the injectors with negative coefficients could be diverting the flow from the other injectors away from the production well. It is also possible that increased injection to the well with negative coefficient prevents inflow of natural recharge fluids with higher chloride concentration.

2.2 Improvements

The linear form of the time term in model (2) was a very convenient assumption we made despite the nonlinear trend in chloride that was readily apparent from the data. Use of the linear time term in the previous section gave us an idea of how much the time variable accounted for variations in chloride. For both Dixie Valley and Palinpinon-I the previous results showed that the time term contributed a very significant part to the chloride model as implied by the high values of time term coefficients. From these observations came the motivation to set aside convenience and identify the correct form of the time term or equivalently the general trend in chloride with time. We have chosen to use wavelet analysis for this task.

Based on the results of our analysis of the Dixie Valley data set, we concluded that the effects of individual injection rate terms on chloride were trivial because their corresponding coefficients were very small compared to the coefficient of the time term. But is the significance of a variable's contribution to the regression solution really

Table 4. Model (2) Coefficients for Palinpinon-I production wells.

Model Parameter	Production Wells							
	OK-7D	OK-9D	OK-10D	PN-15D	PN-16D	PN-17D	PN-18D	PN-19D
a_0	5292.15	4290.85	3875.44	4761.51	3824.35	4917.71	4758.59	5043.92
<i>Injection wells</i>								
PN-1RD	4.49E+00	1.15E+00	3.31E+00	2.56E+00	1.09E+00	1.04E+01	1.81E+00	-8.37E-01
PN-2RD	-2.49E+01	-7.06E+00	-2.27E+00	-7.73E+00	-5.82E-01	-1.13E+01	-2.27E+01	1.01E+00
PN-3RD	1.37E+00	5.97E+00	-2.64E+00	1.88E+00	-3.81E+00	-3.48E+01	-1.40E+00	-9.77E+00
PN-4RD	-1.75E+00	6.28E+00	9.97E-01	-1.63E+01	-5.25E-01	1.04E+00	-6.84E+00	-3.38E+00
PN-5RD	1.38E+01	-1.46E+01	-3.48E+00	1.01E+01	-4.68E-01	2.47E+01	7.15E+00	7.57E+00
PN-6RD	-4.51E+00	8.91E-01	3.62E+00	1.97E+00	-4.20E-01	2.72E+00	-1.75E+00	-4.68E+00
PN-7RD	5.69E+00	-5.69E+00	1.15E+01	1.67E+01	7.65E+00	9.72E+00	-4.55E+00	4.97E-01
PN-8RD	2.56E+00	-1.64E-01	6.77E-01	9.05E-01	2.36E+00	-1.49E+01	4.69E+00	4.92E+00
PN-9RD	1.07E+01	5.81E+00	-1.99E+00	-9.65E+00	9.92E-01	-1.04E+01	2.15E+00	-1.61E+00
t	683.31	322.12	123.82	707.56	590.02	881.05	670.13	621.24

Cont.

Model Parameter	Production Wells							
	PN-23D	PN-24D	PN-26D	PN-27D	PN-28D	PN-29D	PN-30D	PN-31D
a_0	4434.14	3770.09	5552.84	3949.51	5843.37	5233.71	4360.31	4365.67
<i>Injection wells</i>								
PN-1RD	9.54E-02	-2.48E-01	7.56E+00	9.03E-01	6.16E+00	2.01E+00	-4.14E-01	1.33E+00
PN-2RD	-6.39E+00	-4.68E+00	-1.22E+01	-6.89E+00	-1.90E+01	-1.99E+01	-2.95E+00	2.01E-01
PN-3RD	4.71E+00	-5.60E+00	1.04E+00	1.54E+00	-1.06E+00	3.81E+00	3.96E+00	-2.84E+00
PN-4RD	3.02E+00	-7.31E+00	-1.15E+00	1.06E+01	-2.98E+00	-1.92E+00	8.92E-01	-5.57E+00
PN-5RD	-1.46E+00	2.85E+00	1.36E+01	-1.42E+01	5.99E+00	3.28E+00	3.24E+00	5.36E+00
PN-6RD	-1.92E+00	4.09E+00	4.53E-01	8.73E-01	-4.56E+00	-4.24E+00	-2.65E-01	-1.63E+00
PN-7RD	-3.10E+00	2.32E+01	6.82E+00	-5.24E+00	6.31E+00	-3.36E+00	-3.00E+00	-2.40E+00
PN-8RD	-1.86E+00	7.24E+00	3.87E+00	-6.15E-01	2.28E+00	1.31E+00	-3.07E+00	4.79E+00
PN-9RD	1.71E+00	9.74E+00	3.93E+00	1.11E+01	6.52E-01	9.01E+00	-1.15E-01	6.85E+00
t	475.82	661.44	685.94	569.911	708.20	836.911	189.521	677.22

reflected by the magnitude of its coefficient in the regression equation? Would the comparison of those small coefficients from Dixie Valley result in as meaningful and valid conclusions as those derived from comparison of the bigger coefficients in Palinpinon-I?

Regression using model (2) gave us R^2 values that are very close to unity, signifying that the variation in chloride is almost entirely predictable from the model. Ironically, we also obtained regression coefficients that could not be generalized from the early to the later portion of the data set; that is, we observed poor predictive capacity of the model. How could we reconcile these results?

To answer these questions we turned to statistics and the next section outlines our findings.

3.0 MULTIPLE REGRESSION

The high values of R^2 coupled with poor predictive capacity observed in the previous chapter lead us to doubt the validity of the results of the previous regression work. A survey of materials/texts on multivariate statistics suggested that what we had was a case of overfitting: we had a solution that provided a wonderful fit to the sample (the early portion of the data set) but did not generalize to the population (the entire chloride series). According to Tabachnik (19%) overfitting is a result of having a sample size that is too small relative to the number of variables

in the linear model. To illustrate the point, let us take the case of bivariate regression where a straight line ($y = mx + b$) is fitted through the data points available. When calculating the parameters m and b , the square of the prediction error (graphically, the deviation of the data points from the 'best fit' line) is minimized. In the extreme case where only two data points are available, the minimization problem reduces to a deterministic problem; m and b are calculated exactly based on the two data points and the solution becomes perfect (and meaningless).

Tabachnick (1996) suggests the following rules of thumb for the required sample size for multiple regression:

$$N \geq 50 + 8m \quad \text{for testing } R^2$$

$$N \geq 104 + m \quad \text{for testing individual coefficients}$$

Tables 5 and 6 show that for both Dixie Valley and Palinpinon-I the number of data points used in the previous regression analysis was significantly below the required amount of data suggested by Tabachnick (1996). This asserts that the R^2 values and coefficients calculated previously were only artifacts of the data analyzed and do not generalize to extensions of the chloride series. The restriction on the amount of data required for analysis prevented us from using the Dixie Valley data set. Fortunately, however a larger data set from Palinpinon-I was made available to us by PNO-EDC and we used this extended data set for succeeding analyses. In instances where even the extended data set was short of the required amount of data, the only possible solution was to reduce the number of terms in the model to include only those which contribute significantly to the regression solution. The procedure for choosing the important terms is discussed later in this section.

It is worth noting here that the addition of the time term to model (1) pushed the regression problem towards the deterministic region as it lowered the data-to-parameter ratio; hence, the increase in R^2 values observed previously.

Another issue raised earlier was that of the suitability of inferring the contribution of injection rate terms to the regression solution from the sizes of the coefficients alone. According to Tabachnick (1996), interpretation of the multivariate solution based on the sizes of the coefficients alone is strictly possible only in the case where all the independent variables or IVs (injection rates and time in the case of model 2) are indeed independent from each other. Disregarding the interdependence between IVs, there are statistical tests that allow us to tell whether the unique contribution of an IV as represented by its coefficient is significantly different from zero or not; that is, it tells one whether to accept or reject the hypothesis that the coefficient of an IV is zero. One such test is the probability or P-test. According to this test, there is a $(100-x)\%$ probability that an IV is important to the regression solution or equivalently, its coefficient is not equal to zero if its P-value is less than or equal to $x\%$. It is common practice to set x to 5%; hence, there is a 95% certainty that the coefficient of an IV is not equal to zero if its P-value is less than or equal to 0.05. Calculation of P-values is discussed by Bowerman and O'Connell (1990) and is done automatically by the Microsoft Excel regression macro that we used. Note again that the P-test does not take into account the interdependence between IVs.

Table 5. Number of data points in Dixie Valley data set.

Well Name	$m = 9, \text{ model (2.2)}$		
	actual N	$(50 + 8m)$	$(104 + m)$
27-33	32	122	113
28-33	31	122	113
45-33	36	122	113
63-7	44	122	113
73-7	39	122	113
74-7	31	122	113
76-7	56	122	113
82-7	42	122	113
84-7	28	122	113

Table 6. Number of data points in Palinpinon-I data set

Well Name	$m = 10, \text{ model (2.2)}$		
	actual N	$(50 + 8m)$	$(104 + m)$
OK-7D	53	130	114
OK-9D	44	130	114
OK-10D	55	130	114
PN-15D	25	130	114
PN-16D	47	130	114
PN-17D	24	130	114
PN-18D	46	130	114
PN-19D	52	130	114
PN-23D	54	130	114
PN-24D	30	130	114
PN-26D	37	130	114
PN-27D	37	130	114
PN-28D	36	130	114
PN-29D	54	130	114
PN-30D	52	130	114
PN-31D	50	130	114

Considering the need to eliminate unimportant terms in the linear model to meet the data requirement as discussed previously and taking care not to exclude IVs whose importance are masked by their interdependence with other IVs, we have proposed the following procedure for succeeding application of multiple regression analysis:

1. To economize on IVs, temporarily set aside variables with P-values higher than 0.05;
2. Also, eliminate IVs with P-values lower than 0.05 and low values of simple correlation, r ;
3. Inspect IVs which were eliminated in step 1 and put those with high r back to the model;
4. Perform another regression using the reduced model and interpret the results of this regression.

There are several possible variations to the preceding procedure and the one outlined above may not be the best but the important point to consider is the need to be aware of the possible complications that prevent straightforward interpretation of regression results based on the magnitude of coefficients alone.

4.0 WAVELET ANALYSIS

Recent work on the use of wavelet analysis (Jansen and Kelkar, 1997) in analyzing production data from oil fields prompted us to look into its applicability to our problem. In the course of our investigation we found out that wavelet analysis had a capability that serves our purpose of isolating the general trend of a signal (in our case, chloride concentration and injection rate) from its short-term variations. Wavelet analysis allows us to examine features of a signal of any size by decomposing the signal to different detail levels and a coarse approximation. The approximation retains the general trend with time while the details bear information on the signal's fluctuations at different time scales. Fig. 9 illustrates the concept using the chloride concentration signal from well OK-7D from Palinpinon-I. It is worth emphasizing that the approximation to OK-7D chloride shown in Fig. 9 demonstrates that the general trend in chloride is nonlinear, contrary to the assumption in model (2).

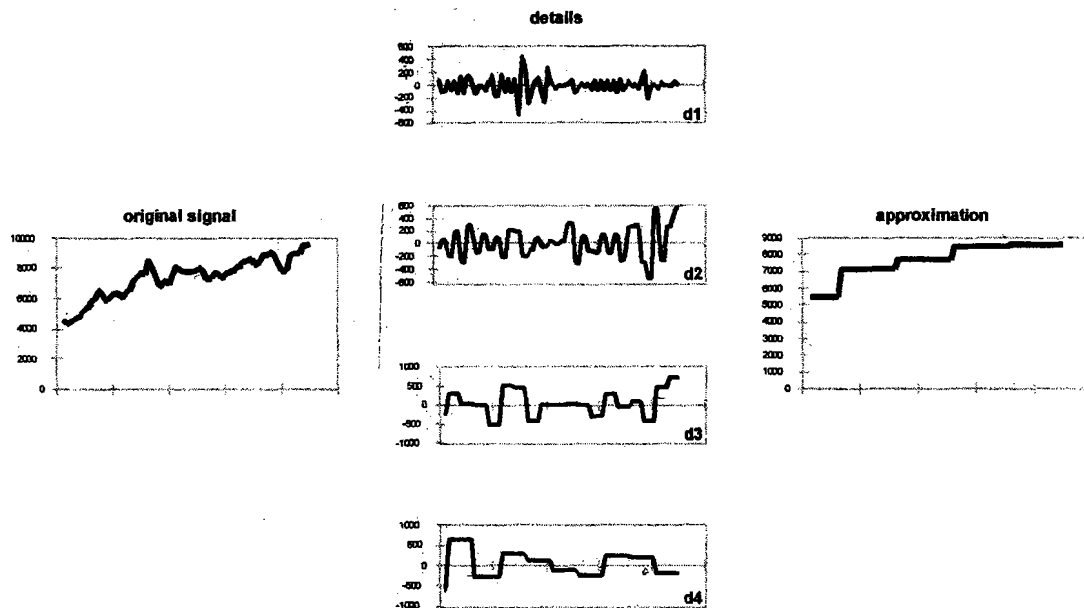


Figure 9. Wavelet decomposition: breaking a function down into a very coarse approximation, with an ordered sequence of detail functions making up the difference. (Chloride concentration, well OK-7D).

Since the effect of changes in injection rates is expected to manifest as short-term variations in reservoir chloride concentrations, it is more appropriate to analyze the detail functions described in the previous chapter instead of the approximation functions. Also, because the approximation functions isolate and retain the general trend in chloride with time, multiple regression of the details does not require a time term in the linear model. Thus, we used the following model:

$$Cl_p = a_1 Q_{11} + a_2 Q_{12} + a_3 Q_{13} + \dots \dots + a_n Q_{1n} \quad (5)$$

where Cl_p = chloride concentration detail in production well, P, Q_{In} = injection rate detail in injection well I_n , a_n = linear coefficient of well I_n

Comparison of coefficients obtained by using model (5) allows us to differentiate the degree of connectivity of different injectors to a given producer. Since details are deviations from local averages multiple regression using details ignores the differences in base chloride levels between producers. Regression results for different producers may therefore be intercorrelated; more specifically, the coefficients obtained may be used to compare the contributions of an injector to different production wells and consequently, to verify any conclusions drawn from the analysis against tracer test results.

The choice of modeling details over approximations was an obvious and straightforward decision. The appropriate detail level to model was less obvious, however. It seemed reasonable at first to assume that the best choice is the one that will give the highest R^2 value. Investigation of the R^2 values obtained from modeling the chloride details of OK-7D invalidated that assumption. Table 7 shows that at level 4 the regression coefficient becomes unity signifying a perfect correlation; and, correlation at succeeding levels remain perfect. As the decomposition level goes up, the detail will have longer time intervals with constant values. This effectively reduces the amount of data to be modeled and results to a perfect, meaningless correlation. The choice is thus narrowed down to levels 1, 2, and 3.

Table 7. R^2 values for multiple regression of OK-7D chloride detail.

Detail Level	R^2
1	0.202
2	0.530
3	0.853
4	1.000
5	1.000

Inspection of injection and chloride details showed that the correspondence between changes in chloride concentration and changes in injection rate is more readily visible at level 3. In Fig. 10 the level 3 details of injection wells PN-6RD and PN-9RD closely follow the detail of OK-7D chloride during intervals when injection to these wells are high. Some degree of correspondence at levels 1 and 2 is also apparent from Fig. 11 albeit not quite as obviously as in level 3. Thus, we have decided to analyze all three levels of detail.

As was done in previous analyses, the chloride data that were recorded at irregular time intervals were converted to monthly average values to put them in the same time basis as the injection rates. Since wavelet analysis requires that data be available in the entire time interval being analyzed, missing chloride data were linearly interpolated. Interpolation was done over maximum intervals of six months and only when no drastic fluctuations were apparent within six months of the interval where interpolation was to be done. Where interpolation was not possible, we analyzed only the longest continuous portion of the data series.

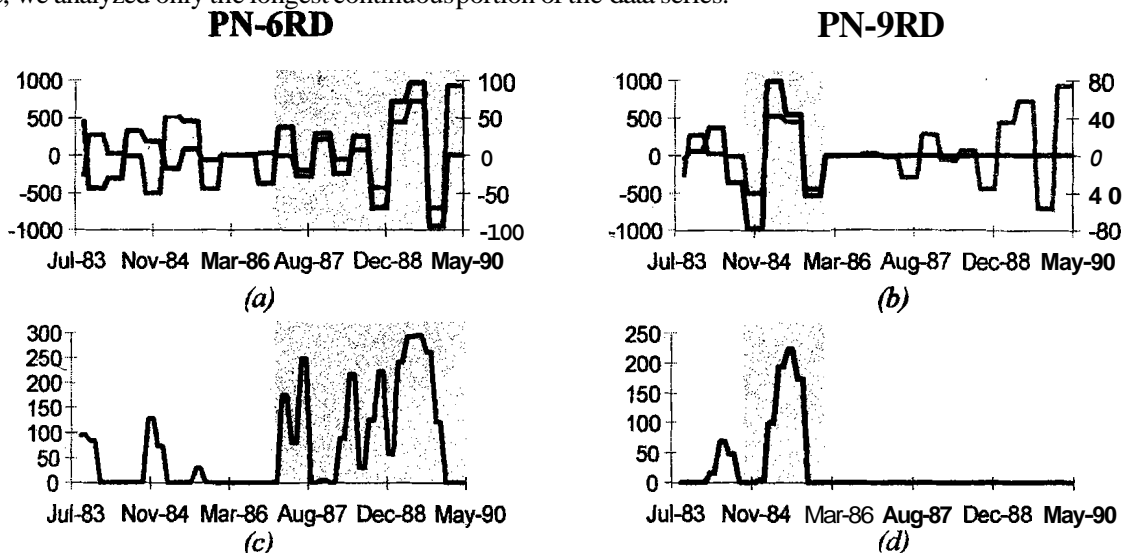


Figure 10. (a) Level 3 detail of OK-7D chloride - light line; level 3 detail of PN-6RD injection rate - dark line. (b) Level 3 detail of OK-7D chloride - light line; level 3 detail of PN-9RD injection rate - dark line. (c) PN-6RD injection rate. (d) PN-9RD injection rate.

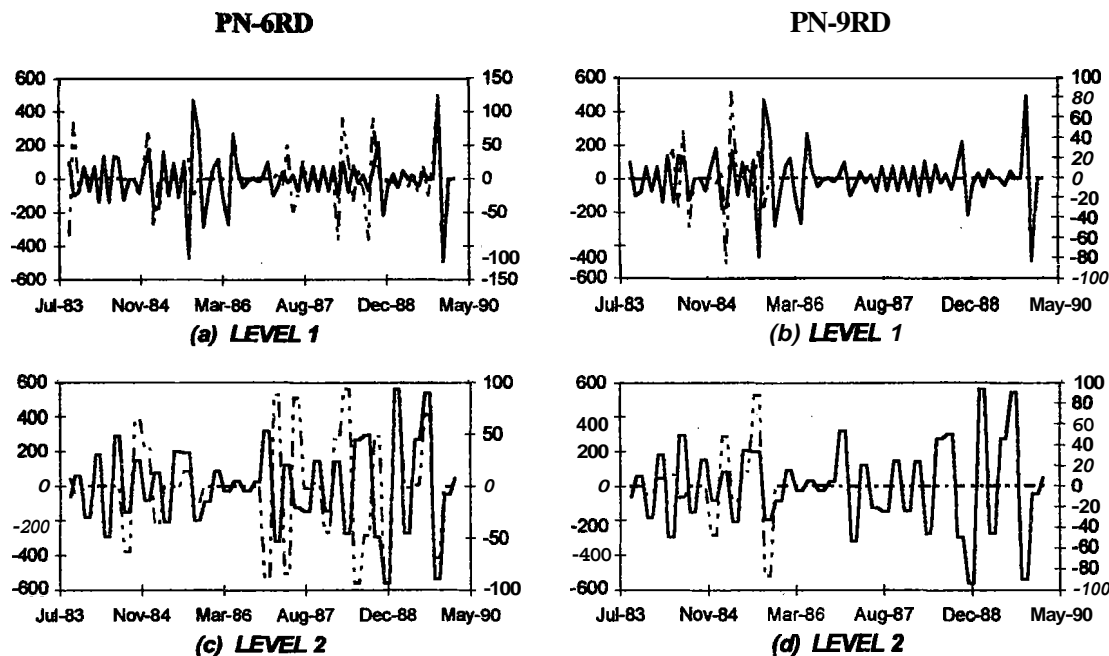


Figure 11. (a) Level 1 detail of OK-7D chloride- solid line; level 1 detail of PN-6RD injection rate -dashed line.
 (b) Level 1 detail of OK-7D chloride- solid line; level 1 detail of PN-9RD injection rate - dashed line.
 (c) Level 2 detail of OK-7D chloride- solid line; level 2 detail of PN-6RD injection rate -dashed line.
 (d) Level 2 detail of OK-7D chloride- solid line; level 2 detail of PN-9RD injection rate - dashed line.

When taking wavelet transforms of discrete data, the algorithms used require that the data set size be a power of two. A common way to precondition the data when this is not true is to "pad with zeroes," that is, to increase the size of the data set to the next larger power of two. Although this is a reasonable approach, it is problematic in that it "dilutes" the signal near the end of the original data set since wavelet coefficients will have zeroes averaged into their computation (Ogden, 1997). We believe that Matlab, the program we used to take the wavelet transform of our data, practices this data preprocessing procedure as evidenced by the inaccurate reconstruction near the end of the data series. It was therefore necessary to truncate the detail component functions used in regression analysis to eliminate the end effects of padding with zeroes. Plots of wavelet component functions retain the end effects but in the analysis, 3 to 6 data points were eliminated from the details series.

In about 1990, the bulk of the injection in Palinpinon-I was moved farther away from the production sector resulting in some injection wells being shut off and new ones being operated. Thus, based on operating time, injection wells in Palinpinon-I can be grouped into three: those which operated between 1983 and 1990, those which started to inject in 1990 and are still injecting, and those which have been injecting since 1983 and are still injecting. It is logical to assume that regression analysis will be best able to assess the degree of contribution of injection wells if all the wells are operating during the time interval over which the regression is done. So, where possible, the chloride series was divided into two time intervals, 1983 to 1990 and 1990 to present; and, regression analysis included only wells which were operating during those periods.

Regression was done for levels 1, 2, and 3 of the detail functions using the procedure outlined earlier. In some cases, that procedure had to be applied repeatedly until the number of terms left in the linear model is such that the data size requirement is met (or almost met). When eliminating injection well terms that had small r values, care was taken not to remove wells which operated only for very short periods of time (the small r values in these cases are unnatural effects of the scarcity of correlatable data).

4.1 Checking Results Against Tracer Test Data

Two sets of radioactive tracer test results were available for comparison with results of our analysis: that of the test conducted on well PN-9RD and one on OK-12RD. Both sets were reported by Macario (1991) and are

reproduced in Table 8. Macario (1991) defined mean transit time as the time it takes for half of the tracer return to reach the production well. Assuming that the mean transit time measures the degree of connectivity between the injector tested and a producer (lower transit times corresponding to stronger connections) Table 8 lists the production wells in order of decreasing connectivity with the injector. Correspondingly, Tables 9 and 10 lists the wells affected by OK-12RD and PN-9RD, respectively, in the order of decreasing coefficients based on regression on all three wavelet detail levels.

Table 10 shows, with the exception of one well, that all wells affected by PN-9RD had positive coefficients. Comparison of Table 10 with Table 8 shows that tracer return was indeed monitored in all wells shown by

Table 8. Radioactive tracer test results for PN-9RD and OK-12RD.

Monitored Well*	Mean Transit Time, days
PN-9RD Tracer Test	
OK-7	5.4
PN-26	13
PN-28	14
PN-29D	15.4
PN-30D	15.7
PN-23	15.8
PN-16D	16
PN-19D	16
PN-31D	16
PN-18D	17.2
OK-9D	monitored. no return
OK-12RD Tracer Test	
PN-15D	7.3
OK-10D	13.8
OK-7D	14.6
PN-29D	monitored. no return

* Only wells which have chloride data are reported here.

Table 9. Regression results for OK-12RD

OK-12RD	
(Affected Well)	Coefficient
PN-23D	2.46
PN-29D	-4.05
PN-31D	-10.82
PN-15D	125.27
PN-16D	-7.40
PN-29D	-3.34
PN-30D	6.15

Table 10. Regression results for PN-9RD.

PN-9RD	
Affected Well	Coefficient
d1	
PN-30D	5.74
PN-29D	3.99
PN-16D	1.47
d2	
PN-19D	4.87
PN-18D	4.06
OK-7D	2.96
PN-16D	2.02
PN-29D	-11.65
OK-7D	9.40
PN-29D	1.83
PN-16D	0.92
PN-23D	0.43

regression analysis to be affected by well PN-9RD, including PN-29D which had a negative coefficient. More importantly it shows that the order of the strength of connection **between** PN-9RD and the wells monitored in the tracer test was most closely mimicked by the results of regression **on** detail level 3 **with** OK-7D showing the strongest connection to PN-9RD and PN-29D, **PN-16D**, and PN-23D displaying **connections** of about the same strength.

On the other hand, comparison of Table 9 with Table 8 **shows** that tracer **return** was observed in **two** of the seven wells shown by regression analysis to be affected by well OK-12RD. **Four** of the seven wells were **not** monitored during the tracer test. As with PN-9RD, the well which is most connected to OK-12RD based **on** the tracer test had the highest coefficient at level 3 regression.

Based **on these** observations we have concluded that regression analysis of details at level 3 best **assesses** the degree of connectivity between wells: high positive coefficients correspond to **strong** connections **and**, negative and low positive coefficients correspond to weak connections.

Harper and Jordan (1985) reported the **following** observation: from May 1984 to **October** 1984 a large increase in reservoir chloride occurred in production wells **PN-19D, 23D, 29D, 31D, OK-7D** and OK-9D when reinjection was shifted to the wells PN-7RD and PN-8RD. This observation matches the results of level 3 detail analysis for well PN-8RD **as** outlined in Table 11: **OK-7D, PN-19D, and PN-31D** were all found to be strongly connected **with** PN-8RD. PN-23D, -29D, and OK-9D may have been receiving reinjection returns **from** OK-7D but no injection rate **data from** OK-7D was available to allow verification with regression results.

On the other hand, Amistoso and Orizonte (1997) reported that OK-10D and PN-20D experienced **enhanced steam** flows which they attributed to reinjection fluids intruding into the production sector at deeper levels. **They** cited the wells **TC-2RD, TC-4RD, PN-3RD** and PN-5RD to be wells that are providing pressure support to the reservoir due to deep injection but attributed the **enhanced steam** flow in OK-10D and PN-20D to **TC-2RD and TC-4RD**, specifically. Regression analysis results for **these** wells (Table 12), however, show that OK-10D is not affected by **TC-2RD**; rather it is affected by **PN-1RD, PN-2RD, and PN-3RD** between 1986 and 1990 and by **PN-3RD, TC-3R, N3 and OK-3R** between 1990 and 1996. It is worth noting that the effect of PN-3RD **on** OK-10D was found to be consistent between the intervals 1986-1990 and 1990-1996 **as** reflected by close r values for the **two** periods (-0.79 and -0.77). The large positive coefficient of well N3 is **suspect** however **as** it conflicts with its negative r value. PN-20D was also analyzed to be affected by PN-3RD. The effect of TC-2RD and TC-4RD on PN-20D could **not be ascertained from** regression analysis due to insufficient chloride data from PN-20D after 1990.

Pamatian (1997) reported that reinjection fluid **from** TC-2RD **neutralized** the fluid **acidity in** wells OK-10D and PN-13D. **Again**, the effect of TC-2RD **on** OK-10D was **not** substantiated by regression results but **its** effect **on** PN-13D was (Table 13). **Again**, terms **with** conflicting r and coefficient **signs** posed interpretation problems.

5.0 CONCLUSIONS AND RECOMMENDATIONS

Based on the results of regression analysis of chloride and injection rate **data** from Dixie **Valley**, we have concluded **that** multilinear modeling is **not** suitable **for** analyzing data **sets** which lack sufficient time variability or "texture".

A closer look at multiple regression techniques showed that what seemed to be highly **encouraging** results (**high** R^2 values) **from** **prior multilinear** modeling were **but** effects of the **scarcity** of **data** used in the **correlation**; hence, no meaningful physical interpretation may be **drawn** from them. Moreover, it showed that care should be **taken** not to base the **interpretation** of multiple regression results **on** straight-forward **comparison** of coefficients alone.

Table 11. Level 3 regression results for PN-8RD

PN-8RD	
Affected Well	Coefficient
d3	
OK-7D	3.14
PN-16D	0.64
PN-18D	2.76
PN-19D	4.93
PN-30D	-1.36
PN-31D	10.49

Wavelet analysis provided more useful results. Qualitative field observations and tracer test data agreed **best** with the results of regression on level 3 detail of chloride concentration and injection rates in Palinpinon-I: wells identified by tracer tests to be strongly connected had **high** positive coefficients and weak connections were indicated by negative and low positive coefficients at level 3 regression. This suggests that producer-injector interactions are **best** detected by correlating changes in chloride concentration over **periods** of four months (corresponding to level 3 resolution) with corresponding four-month fluctuations in injection rates. While the **good** correlation at such a relatively low level of time resolution may be explained as the result of the natural dispersion of chloride and injection rate signals as they propagate through the reservoir, it is also possible that this is due to the loss of information brought about by the use of monthly averaged **data** values in the analysis. It is therefore recommended that **both** chloride and injection rate data be recorded more frequently and the analysis be done **on this** larger **data set**. It is **also** possible that the Haar wavelet that **was** used in signal decomposition **was** too **coarse** in that it contributed to the loss of texture in the **data**. Investigation of the effects of using smoother wavelets is **also** recommended.

Table 12. Level 3 regression statistics for OK-10D and PN-20D.

OK-10D (1986-1990)

d3					
Regression Statistics					
Multiple R	0.839369312				
R Square	0.704540842				
Standard Error	79.20406691				
Observations	48				

	Coefficients	r (simple)	Standard Error	t Stat	P-value
pn1rd	1.0860989	0.742898166	0.432894148	2.506924886	0.01577982
pn2rd	-3.71087522	-0.72596153	1.775342121	-2.09023105	0.04227759
pn3rd	-4.4114713	-0.78948113	1.783793966	-2.4730634	0.01723762

OK-10D (1990-1996)

d3					
Regression Statistics					
Multiple R	0.835160095				
R Square	0.697492385				
Standard Error	150.1670809				
Observations	80				

	Coefficients	r (simple)	Standard Error	t Stat	P-value
on3rd	11.4979076	-0.77103302	1.485833395	-7.7383569	3.4832E-11
tc3r	1.704839524	0.522611064	0.404330736	4.216448003	6.7759E-05
n3	30.87773863	-0.65393407	6.498963053	4.751179278	9.3559E-05
ok3r	-798184387	-0.36088849	2.134566707	-3.44887229	0.00082135

PN-20D(1983-1989)*

d3					
Regression Statistics					
Multiple R	0.635764285				
R Square	0.404196201				
Standard Error	495.2038572				
Observations	80				

	Coefficients	r (simple)	Standard Error	t Stat	P-value
pn1rd	10.21445053	0.346132276	1.741845857	5.864152953	1.0654E-07
pn3rd	12.28448348	0.357197703	2.153358367	5.704802167	2.062E-07
pn6rd	4.493788629	0.105136293	1.637086662	2.744879074	0.00752785

*No chloride data was available from 1990 to 1993 and remaining data points were not sufficient for analysis.

Table 13. Level 3 regression statistics for PN-13D

PN-13D (1990-1996)

d3					
Regression Statistics					
Multiple R	0.962568329				
R Square	0.926537788				
Standard Error	74.57273176				
Observations	70				

	Coefficients	r (simple)	Standard Error	t Stat	P-value
tc2rd	4.826880146	0.403207897	0.459363951	10.5077687	1.4452E-15
tc3r	1.989170863	0.37912479	0.16882133	11.78269865	1.1114E-17
tc4r	52.80156771	-0.47305009	4.657792303	11.33617909	5.9844E-17
ml1rd	-191.386684	-0.73129226	16.95658741	-11.2875288	7.1995E-17
n3	11.67741779	-0.76944945	1.849731659	6.313033424	2.9328E-08
ok3r	27.49754181	-0.47241942	2.227450499	12.34484978	1.3816E-18

Emphasis is also placed on the need for continuous data measurements when doing wavelet analysis. Highly intermittent measurements result in **data** loss: since it is considered safe to interpolate only over short **periods** of time, the lack of **data** over long time intervals forces one to disregard the **data** collected prior to such periods when doing the analysis.

Another possible improvement to consider in **future** regression analyses is to take into account possible nonlinearity in the variation of chloride with injection rates. While nonlinearity does **not** invalidate the analysis, it certainly weakens it **as** the relationship **between** chloride concentration and injection rates is not completely captured by the coefficients of the linear model. Although regression analysis **uses** a **linear** model, effects of nonlinearity in the variation of chloride with injection rates may be incorporated into the model by using nonlinear terms: the model is **kept** linear even though the individual terms are not. Results of this modified analysis will **be** more difficult to interpret however, because the **strength** of interaction **between** producers and injectors will **be measured** not only by the magnitude of the coefficients but also by **the** exponent of each term.

6.0 NOMENCLATURE

a_0	=	a constant associated with local initial chloride concentration
a_n	=	linear coefficient of well L
b	=	linear time term coefficient
Cl_{in}	=	chloride concentration in injector well, In
Cl_p	=	chloride concentration/chloride concentration <u>detail</u> in production well, P
m	=	number of predictors
N	=	number of data points
Q_{in}	=	mass flow rate/mass flow rate <u>detail</u> to injection well, In
r	=	simple regression coefficient
R^2	=	multiple regression coefficient
S	=	standard deviation
SS_{reg}	=	sum of squared deviations of predicted Y from the mean
SS_Y	=	sum of squared deviations of Y from the mean
Y	=	dependent variable being modeled
Y'	=	predicted values of Y
\bar{Y}	=	average value of Y

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