

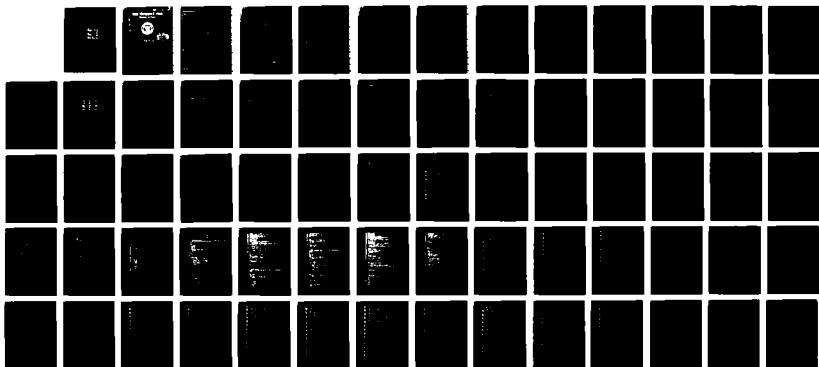
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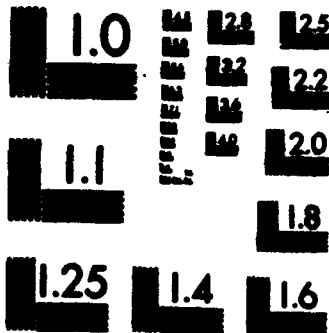
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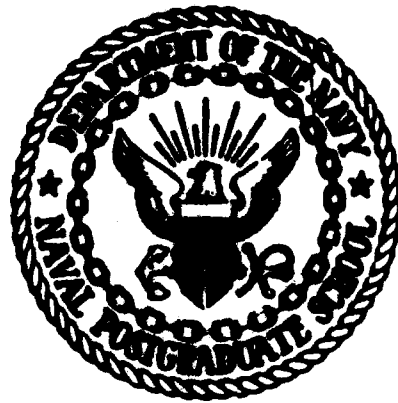
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THESIS

THE CURVE FITTING
OF
PORTABLE WEAR METAL ANALYSIS

by

Byung H. Min

March 1987

Thesis Advisor

Harold J. Larson

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The Curve Fitting
of
Portable Wear Metal Analysis

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Submitted in partial fulfillment of the
requirements for the degree of

MASTER OF SCIENCE IN OPERATIONS RESEARCH

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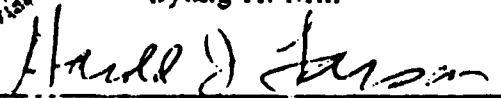
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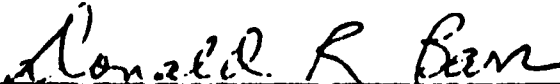


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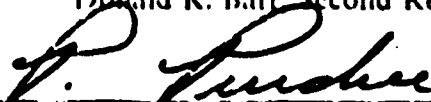
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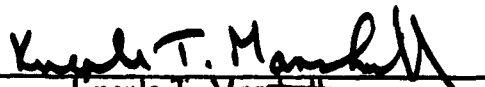
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ABSTRACT

This thesis describes the curve fitting algorithm that the Portable Wear Metal Analyzer used for calculating concentrations in ppm and compares this with some alternative algorithms. Each algorithm considered fits a curve to the three standards used, which were 20%, 50%, and 100% of the full scale for all nine wear metals. APL was used for all programs.

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TABLE OF CONTENTS

I.	INTRODUCTION	8
II.	PWMA AND DATA DESCRIPTION	9
	A. GENERAL	9
	B. THE PORTABLE WEAR METAL ANALYZER	10
	C. TEST PROGRAM AND DATA COLLECTED	11
III.	ALGORITHMS	16
	A. GENERAL	16
	B. ALGORITHM 1	17
	C. ALGORITHM 2	17
	D. ALGORITHM 3	20
	E. ALGORITHM 4	22
	F. ALGORITHM 5	23
IV.	ALGORITHM COMPARISON	27
	A. GENERAL	27
	B. SUM OF SQUARES METHOD.	27
V.	CONCLUSIONS AND RECOMMENDATIONS	39
	APPENDIX A: PWMA FIELD TEST BURN SEQUENCE	40
	APPENDIX B: DATA PPM401	42
	APPENDIX C: PROGRAMMING LIST	43
	APPENDIX D: SUMMARY OF RANK	48
	LIST OF REFERENCES	65
	INITIAL DISTRIBUTION LIST	66

LIST OF TABLES

1. ABSORBANCE DATA OF TUBE 401	12
2. DATA OF EACH LABORATORY	13
3. TABLE OF OIL SAMPLE MIXTURES	15
4. ESTIMATED CONCENTRATION BY 6 ALGORITHMS	32
5. ALGORITHM COMPARISON FOR FE	32
6. SUMMARY OF THE RANKS FOR TUBE 401	33
7. RANKS OF 24 TUBES FOR FE	34
8. RANKS OF 24 TUBES FOR AG	34
9. RANKS OF 24 TUBES FOR AL	35
10. RANKS OF 24 TUBES FOR CR	35
11. RANKS OF 24 TUBES FOR CU	36
12. RANKS OF 24 TUBES FOR MG	36
13. RANKS OF 24 TUBES FOR NI	37
14. RANKS OF 24 TUBES FOR SI	37
15. RANKS OF 24 TUBES FOR TI	38
16. THE BEST ALGORITHM FOR EACH ELEMENT	38

LIST OF FIGURES

2.1	Calibration Technique used by the PWMA	14
3.1	An Asymptote Line in Algorithm 2	20
3.2	Possible Observed Curve from Algorithm 3	21
3.3	The Curve Bending Downward at Lower Concentration by Algorithm 4	23
3.4	The Curve Bending Upward at High Concentration by Algorithm 4	24
3.5	The Curve Bending Downward at High Concentration by Algorithm 5	25
3.6	The Curve Bending Upward at Lower Concentration by Algorithm 5	26
4.1	The Horizontal Deviation of Residual by Algorithm 2	28
4.2	The Horizontal Deviation of Residual by Algorithm 4	30

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I. INTRODUCTION

Spectrometric oil analysis is used to determine the type and amount of wear metals in lubricating oil samples. The Portable Wear Metal Analyzer (PWMA) was developed for the USAF to perform engine oil analysis when aircraft are deployed away from their home bases. The analyzer can detect nine wear metals in parts per million concentration ranges. PWMA takes samples with known concentration to obtain absorbance areas, and uses those to determine a calibration curve. Absorbance areas obtained from unknown samples are compared to the calibration curve so that their concentration can be calculated.

This thesis describes the curve fitting algorithm that the PWMA used for calculating the concentration in ppm and compares this with some alternative algorithms for the curve fitting. We compare these algorithms to see which algorithm performs best. Each algorithm considered fits a curve to the three standards used 20%, 50%, and 100% of the full scale for all nine wear metals. APL was used for all programs.

Data gathered as the result of the field test of the PWMA are described in Chapter II. Three rational functions, and polynomials which are quadratic in absorbance and quadratic in concentration are suggested as alternatives in Chapter III. A squared error technique is employed to compare the algorithms' performances.

II. PWMA AND DATA DESCRIPTION

A. GENERAL

The PWMA test program was conducted at Naval Air Station, Pensacola FL, and three Air Force field test sites, Langley AFB, VA, Myrtle Beach AFB, SC, and Elmendorf AFB, AK in 1985. The purpose of this test plan was to determine the effectiveness and evaluate the functionality of the newly developed PWMA prototypes for measuring wear metal content in turbine engine lubricating oils under field conditions.

When aircraft are deployed from their home bases to austere sites where normal oil analysis capabilities are nonexistent, there still exists a requirement to determine wear metal concentration in turbine engine lubricants. Due to the critical nature of turbine engine lubrication system component wear, particularly in high performance systems such as the F-15, F-16, and A-10 aircraft, the requirement for oil analysis may be on an "after each flight" basis. Presently, the Air Force deploys Flame Atomic Absorption (FAA) spectrophotometers [Ref. 1: p. 50], which were not designed for mobile use. These instruments are sensitive and highly susceptible to damage when deployed for mobility exercises; such damage may lead to partial or complete loss of oil analysis capability in early stages of such exercises. In addition, the instruments require hazardous support accessories, such as compressed gases (nitrous oxide and acetylene) and solvents for dilution (and, in some past cases, acid solutions for digestion of wear metal particulates).

In 1980 HQ TAC issued TAF SON 305-80, establishing the requirement for a critically needed light-weight portable aircraft engine oil analysis unit. In response, AFWAL¹ undertook a contractual effort to determine the feasibility of developing such a device (Contract Number F33615-80-C-2037). Since no such spectrometer with required light-weight portable features was commercially available, a major development program was initiated to ruggedize and miniaturize the conceptual hardware for mobility use. The actualization of the portable wear metal analyzer has been achieved under Contract Number F33615-81-C-2080 by Perkin-Elmer Corporation, Applied Science Division.

¹Air Force Wright Aeronautical Laboratory

B. THE PORTABLE WEAR METAL ANALYZER

The PWMA features simultaneous measurement of nine wear metals in lubricating oils with an analysis cycle of four minutes (via air cooling): Silver (Ag), Aluminum (Al), Chromium (Cr), Copper (Cu), Iron (Fe), Magnesium (Mg), Nickel (Ni), Silicon (Si), and Titanium (Ti). Moreover, with the potential modification to water cooling on production units, this time can be reduced to two minutes. In addition, the PWMA may be operated on either 110/220 VAC, 50 or 60 Hertz, to accommodate use in both the U.S and Europe. The PWMA is packed in two aluminum "suitcases": the electronics/argon case, weighing forty pounds; and the furnace/optics case, weighing sixty pounds. The cases are readily set-up and interconnected by two multiconductor cables and one pneumatic hose. Developed for simple operation and reduced dependence on skill and labor of the operator, the PWMA is characterized as follows: ease of set-up and self-contained configuration; minimal consumables (sample tips, graphite tubes and argon gas supply); no hazardous gases or chemicals (inert compressed gas only); microprocessor-controlled software for calibration, diagnostic etc.; and elimination of transcription errors by automated read-out/print-out of data and results.

The PWMA operates on the principle of absorption of characteristic radiation by atoms of the wear metals of interest. Two multi-element hollow cathode lamps (HCLs), alternately pulsed for source combination and modulation, provide emission source lines for discrete line absorption by specific atoms, which are formed through heating in a preprogrammed ramped cycle (up to 3000° C) in the graphite furnace. All nine absorption lines can be separated simultaneously by the polychrometer and transmitted to respective discrete photomultiplier tubes (where the signal is transduced from energy to current). Next, each individual electrometer converts the current to a voltage signal, which is amplified and processed by the analog (A/D) converter. The output is printed in concentration (parts per million, ppm) for all measurements.

The operation of the PWMA is as follows. The operator first introduces an oil sample into the graphite tube of the Graphite Furnace Assembly. The PWMA start button, located on the control panel, is then depressed, and the entire analysis program is initiated and controlled by an "on board" microcomputer. As the graphite furnace executes the atomization cycles, the emission line spectra originating from the HCLs are focused through the graphite tube to the polychrometer where they are separated into individual spectral lines and are further conveyed to the exit slits. The spectral

lines pass through their respective exit slits where their spectral intensities are detected and measured by individual Photomultiplier Tubes (PMTs). Wear metals present in the oil sample will absorb the light associated with each metal's spectral line, thus reducing the intensities reaching the PMTs. The amount of light absorbed is related to the concentration of metal in the oil sample.

The PWMA has several function modes. In the calibration mode, a prepared sample with known concentrations is used to get the absorbance data. Analysis mode is selected to analyze the unknown samples. The prototype PWMA prints out two types of data in the analysis mode, concentration data (computed by the on-board microprocessor) and absorbance data (in parentheses). These data can be used to compare analysis algorithms.

C. TEST PROGRAM AND DATA COLLECTED

The four prototypes were tested at their respective field evaluation sites. The test followed the PWMA field test burn sequence, which is in Appendix A. The graphite tubes, used in the graphite furnace, deteriorate gradually with use, and should be replaced after every 160 sample burns. If a new graphite tube is being used, preconditioning of the tube is required. Laboratories were to follow the test sequence which included calibration samples, verification samples, correlation samples, random samples, reslope, and so on. Three calibration samples were prepared in MIN-L 7808 oil containing 20%, 50% and 100% of full scale for all nine metals. Perkin-Elmer provided these samples. Three verification samples were prepared in MIN-L 7808 oil containing 10%, 40% and 70% of full scale for all nine metals. Perkin-Elmer provided these samples, too. Correlation Samples used turbine engine oil samples. Each sample contained a minimum of 7 wear metals. Random samples were routine daily samples obtained at each evaluation site. This analysis employs only the calibration and verification samples.

The concentration data printed out by the PWMA are calculated by the microprocessor with the use of a calibration technique. Samples with known concentrations are analyzed first and the absorbance data obtained are used to construct calibration curves. Absorbance data obtained from unknown samples are then compared to the calibration curves so that their concentrations can be calculated. Later verification samples obtained from known samples give an indication of how valid the calibration curve is.

Table 1 shows data collected from one of the field test sites. P523 means "Pensacola May 23rd, 1985", the place and the date the data were collected, "401" is the tube sequential number. CAL IA and CAL IB stands for the calibration sample 20%. Cal II and Cal III are 50% and 100% of full scales respectively. In order to provide a more accurate result, a two-point average calibration was used for the field test, i.e., two CAL I samples were run at CAL I position; two CAL II samples at CAL II position; and two CAL III samples at CAL III position. For each calibration run, the numerical values are the absorbance data printed out (rounded to integers). The PWMA averaged the two samples at the same level, and stored them as the calibration points, but the average value was not printed out.

TABLE 1
ABSORBANCE DATA OF TUBE 401

P523 401		Element								
		Fe	Ag	Al	Cr	Cu	Mg	Ni	Si	Ti
CAL IA		58	53	10	38	23	44	40	10	10
CAL IB		56	52	10	34	22	46	38	8	9
CAL IIA		115	98	19	83	47	60	92	28	30
CAL IIB		111	100	21	84	43	61	96	18	29
CAL IIIA		143	129	26	104	70	72	119	33	42
CAL IIIB		145	131	24	104	71	74	120	35	44
VERIF 10	CONC	9	1	2	1	4	3	3	2	5
VERIF 10	ABS	28	28	5	17	12	32	19	4	11
VERIF 40	CONC	44	4	10	4	18	9	12	9	7
VERIF 40	ABS	105	90	17	70	42	56	79	20	20
VERIF 70	CONC	70	7	9	6	28	17	20	12	10
VERIF 70	ABS	129	114	16	88	56	66	105	26	29

Part of the data for tube 401 is shown in Table 1. The complete data for tube 401 is attached in Appendix B. VERIF 10, VERIF 40 and VERIF 70 are the verification samples. The data in the first row of VERIF 10 is the computed concentration and the second row is absorbance data.

Table 2 shows all the data available for this study, Langley by location tested 6 tubes, Myrtle Beach tested 5 tubes, Pensacola tested 11 tubes, and Elmendorf tested 6 tubes. All 28 tubes were used during the field test program.

TABLE 2
DATA OF EACH LABORATORY

Lab.	Tubes			Total
Langley	PPM201 PPM204	PPM202 PPM206	PPM203 PPM207	6
Myrtle Beach	PPM302 PPM305	PPM303 PPM306	PPM304	5
Pensacola	PPM401 PPM405 PPM409 PPM415	PPM403 PPM406 PPM412 PPM416	PPM404 PPM408 PPM413	11
Elmendorf	PPM501 PPM504	PPM502 PPM505	PPM503 PPM507	6

In the PWMA, three calibration standards of 20%, 50% and 100% are used to construct the calibration curve. The absorbance data in CAL IA through CAL IIIB are used to compute the calibration curve coefficients. A pair of two point rational equations were used to fit the calibration curves, one for each of the nine elements. The function used is given in equation 2.1.

$$C = K_1 A / (K_2 A - 1) \quad (2.1)$$

where K_1 , K_2 are unknown coefficients, C is concentration in ppm, and A is absorbance data.

The first part of the curve uses absorbance values obtained from 20% and 50% samples and the second part of the curve uses values obtained from 50% and 100%. We will explain how the concentration value of 9, the first line of the VERIF 10 in Table 1, was computed in the PWMA. Figure 2.1 shows the calibration technique used by the PWMA. In Figure 2.1, the first part of the curve for the FE is given by the solid line, the second part of the curve is given by the dashed line, and symbol + identifies the three calibration points. The verification absorption values of 129, 105

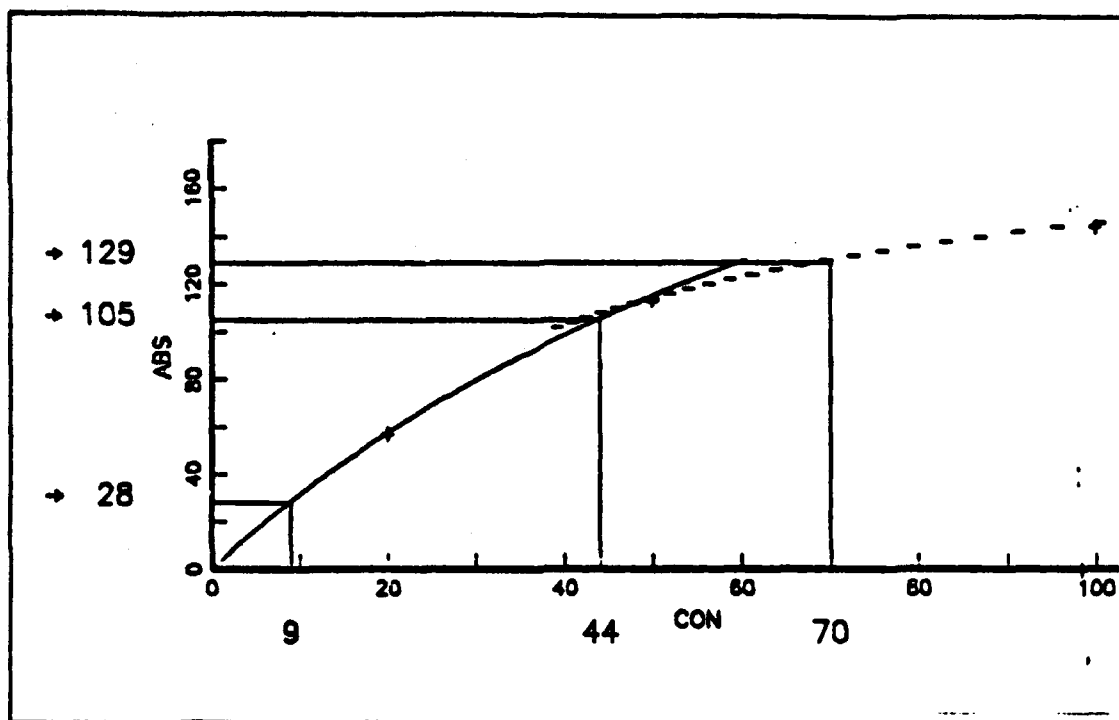


Figure 2.1 Calibration Technique used by the PWMA.

and 28 are applied to the curve, and the calibrated data computed by the PWMA are 70, 44 and 9 respectively. However, the correct values are 70, 40 and 10. Verification sample VERIF 10 which has known concentration 10 produces absorbance data 28, as shown in Table 1. If the absorbance data 28 was fitted to the curve from the absorbance axis, then its concentration should be close to 10. However, it shows some deviation because the curve point projected to the concentration axis makes concentration 9, as shown in Figure 2.1. When the observed absorbance is within the range of 0 to the absorbance value at 50% concentration, the first curve is applied to compute the concentration. For the absorbance data greater than this, the second curve is applied to compute the concentration. The computer selects the proper curve to calculate data, dependent upon the calibrating concentrations of all nine elements. [Ref. 2: p. 67]

The PWMA has a multi-element capability, the 9 elements already listed in Table 1, Iron, Silver, Aluminum, Chromium, Copper, Magnesium, Nickel, Silicon and Titanium.

The full scale capabilities for these elements are shown in Table 3. For example, the full scale value for Fe is 100, and is 10 for Ag. The columns labeled 20%, 50% and 100% are CAL I, CAL II and CAL III samples respectively. The 10%, 40% and 70% columns are VERIF 10, VERIF 40 and VERIF 70 respectively. The concentration varies with the sample and the element.

TABLE 3
TABLE OF OIL SAMPLE MIXTURES

METAL	concentration (ppm)					
	10%	20%	40%	50%	70%	100%
FE	10	20	40	50	70	100
AG	1	2	4	5	7	10
AL	2.5	5	10	12.5	17.5	25
CR	1	2	4	5	7	10
CU	4	8	16	20	28	40
MG	2.5	5	10	12.5	17.5	25
NI	3	6	12	15	21	30
SI	2	4	8	10	14	20
TI	2	4	8	10	14	20
		CAL I		CAL II		CAL III

III. ALGORITHMS

A. GENERAL

If one increment of concentration C causes an equal increment of absorbance A at all levels of C , then we say that there is no curvature in the analytical curve expressing the relation between C and A . In reality, in the presence of stray light the analytical curve bends towards the concentration axis. [Ref. 3: p. 681]

Calibration curvature can be controlled to an extent by the instrument design. The degree of curvature for iron at the 248.3-nm wavelength, for example, is partially controlled by the spectral bandwidth since most of the curvature can be attributed to overlap of nearby nonabsorbed lines in the iron spectrum. This curvature effect is analogous to stray light interferences observed in molecular spectroscopy. However, at practical instrument settings, it is not possible to eliminate this curvature completely.

B. E. Limbeck investigated transformations of the data from AA (Atomic Absorption) instruments to forms that would produce more easily handled curves. Graphic trials on several possible transformations revealed that a plot of A/C vs A or C/A vs C produced pseudocalibrations that were much less curved than the A vs C calibrations, even for the extreme cases. He concluded that over a limited concentration range, two or three standards were sufficient to produce excellent fits for AA calibrations with the rational algorithm. [Ref. 4: p. 96]

As already mentioned, the PWMA uses a rational function which uses two standards; to cover the full range this equation is used twice. This same function will be used twice in the following analysis. The PWMA on board computer uses single precision arithmetic and establishes its calibration curve internally. The data available for this study was printed out in the PWMA paper tapes, which rounded the observed absorbance numbers. Thus, if one uses the PWMA equation to establish a curve, and then reads the rounded absorbance figures back through this curve, one does not exactly arrive at the concentration printed out. Thus the first algorithm should give essentially the same results as those labeled PWMA, which were the rounded figures from the internal curve. The other two rational functions have three unknown coefficients so that they are used one time with three calibration standards. We will examine three rational functions and two quadratic functions, which are defined below.

$$\text{Algorithm 1 : } C = A / (K_1 + K_2 A)$$

$$\text{Algorithm 2: } C = (K_1 A + K_3 A^2) / (K_2 A - 1)$$

$$\text{Algorithm 3: } C = A / (K_1 + K_2 A + K_3 A^2)$$

$$\text{Algorithm 4: } C = K_1 + K_2 A + K_3 A^2$$

$$\text{Algorithm 5: } A = K_1 + K_2 C + K_3 C^2$$

B. ALGORITHM 1

This is useful when a simple linear approximation gives a good fit to the plot of A/C versus A. The following rational function was chosen as algorithm 1:

$$A / C = K_1 + K_2 A \quad (3.1)$$

It is clear by a simple rearrangement of the formula that the concentration is a function of absorbance:

$$C = A / (K_1 + K_2 A) \quad (3.2)$$

For brevity, the calibration algorithm using this linear rational function is referred to as the "rational method" and is the one actually employed in the PWMA. This algorithm uses a pair of two-point rational equations 3.2 to fit data for all nine elements. The first part of the curve uses absorbance data obtained from 20% and 50% samples, and the second part of the curve uses data obtained from 50% and 100% samples.

C. ALGORITHM 2

William B. Barnett found that the equation,

$$C = K_0 \frac{K_1 A + K_3 A^2}{K_2 A - 1} \quad (3.3)$$

was applicable in fitting the analytically useful concentration range of atomic absorption calibration curves. [Ref. 5: p. 829]

In this equation, C is the concentration, A is the observed absorbance, and the K_j 's are coefficients which must be determined during calibration. The coefficient K_0 is the reslope constant which is set equal to 1.0 during the initial standardization, which gives

$$C = \frac{K_1 A + K_3 A^2}{K_2 A - 1} \quad (3.4)$$

In equation 3.4, concentration has been expressed as a function of absorbance. This is the reverse of the usual representation where absorbance is considered to be a function of the standard concentrations. However, expressing concentration as the dependent variable facilitates its computation without the need for extracting roots of equations as others had done. Equation 3.4 can actually be used in three forms: as shown in the PWMA, with K_3 being set equal to zero,

$$C = K_1 A / (K_2 A - 1) \quad (3.5)$$

or with K_2 and K_3 both being set equal to zero. This latter case is, of course, the equation for a straight line which passes through the origin.

When the number of standards matches the number of coefficients to be determined, it is a simple matter to determine the coefficients by the solution of simultaneous equations. We will employ simultaneous equations, to show how to get the coefficients in APL.

Suppose

$$C = (c_1, c_2, c_3), \quad A = (a_1, a_2, a_3).$$

If we express both sides of equation 3.4 in terms of components and then equate corresponding components, we obtain the system

$$c_i = \frac{K_1 a_i + K_3 a_i^2}{K_2 a_i - 1} \quad i = 1, 2, 3 \quad (3.6)$$

which leads to

$$K_2 a_i c_i - c_i = K_1 a_i + K_3 a_i^2 \quad i = 1, 2, 3 \quad (3.7)$$

i.e.,

$$c_i = -K_1 a_i + K_2 a_i c_i - K_3 a_i^2 \quad i = 1, 2, 3 \quad (3.8)$$

Define

$$K = (K_1, K_2, K_3) \text{ and } P = \| p_{ij} \|$$

with

$$p_{11} = -a_i,$$

$$p_{12} = a_i c_i,$$

$$p_{13} = -a_i^2.$$

Then

$$C = P K \quad (3.9)$$

The solution for K is

$$K \leftarrow C \blacksquare P$$

where \blacksquare is the matrix inverse operator in APL [Ref. 6: pp. 82, 83].

While this curve may well faithfully represent the Absorbance - Concentration relationship over the useful range, there are certain possible observed values where it does not. For example algorithm 2 will not perform well when the observed pairs (c_i , a_i) lead to $K_2 = 1/a$, where $a_1 < a < a_3$, since the denominator of equation 3.6 becomes zero. For example, with $C = (20, 50, 100)$, $A = (43, 117, 152)$, the calibration coefficients K are $(-0.4693, 0.00694, 0.003322)$. The fitted curve defined by these coefficients is shown in Figure 3.1. The horizontal line is an asymptote line at A equal to 144.16, which is the reciprocal of K_2 . If the observed absorbance is slightly less than 144.16, the computed concentration will be a large negative number, and when the absorbance data is slightly greater than 144.16, then the computed concentration will be a large positive number, a behavior which would not be acceptable.

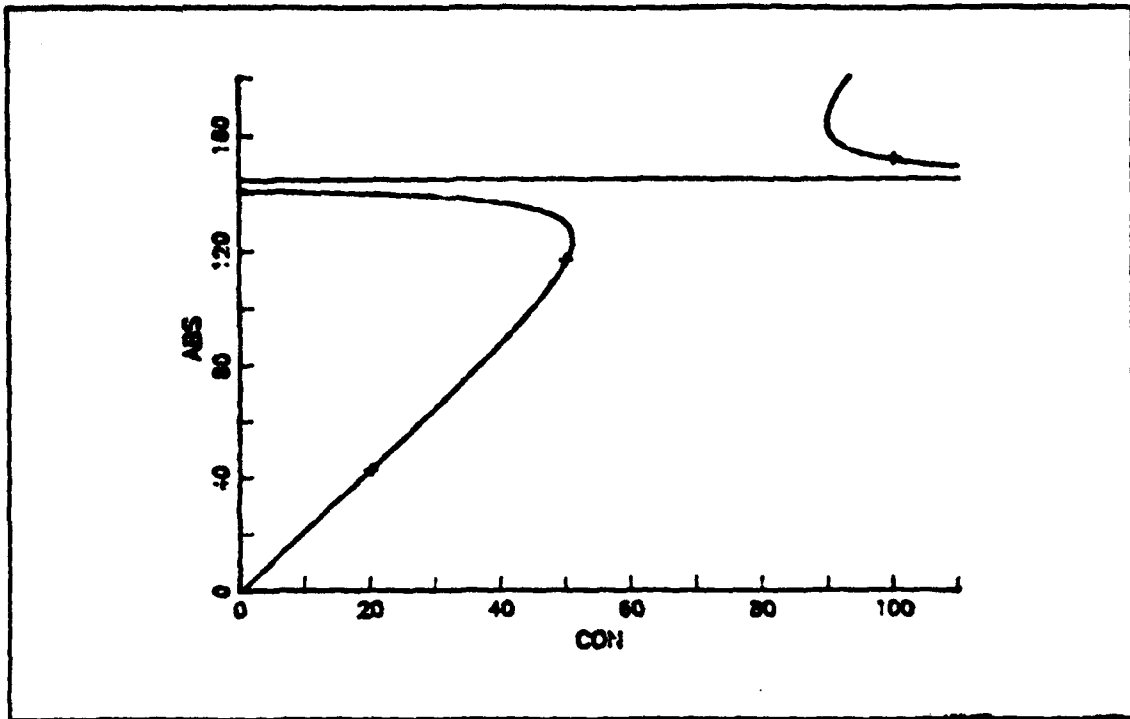


Figure 3.1 An Asymptote Line in Algorithm 2.

D. ALGORITHM 3

This algorithm was suggested by Limbeck [Ref. 4: p. 96].

$$C = \frac{A}{K_1 + K_2 A + K_3 A^2} \quad (3.10)$$

This is a different extension of algorithm 1, namely, $A/C = K_1 + K_2 A$. In equation 3.10, the quadratic term is included. To get the coefficients in APL, we again employ the notation of simultaneous equations.

$$K_1 + K_2 a_i + K_3 a_i^2 = a_i/c_i \quad i = 1, 2, 3 \quad (3.11)$$

Now, let

$$C = (a_1/c_1, a_2/c_2, a_3/c_3) \text{ with}$$

$$P_{i1} = 1,$$

$$P_{i2} = a_i,$$

$$P_{13} = a_i^2.$$

and again $C = P K$, so the solution is

$$K = C \cdot P^{-1}$$

This algorithm can also produce a curve with an interior asymptote, like algorithm 2, when the denominator is zero.

Another type of undesirable shape can occur with this case (as well as algorithm 2). This is illustrated by taking

$$C = (20, 50, 100),$$

$$A = (43, 75, 152).$$

Then, we find

$$K = (3.632, -0.0425, 0.000188).$$

The resulting curve is pictured in Figure 3.2. Increased absorbance will cause lower

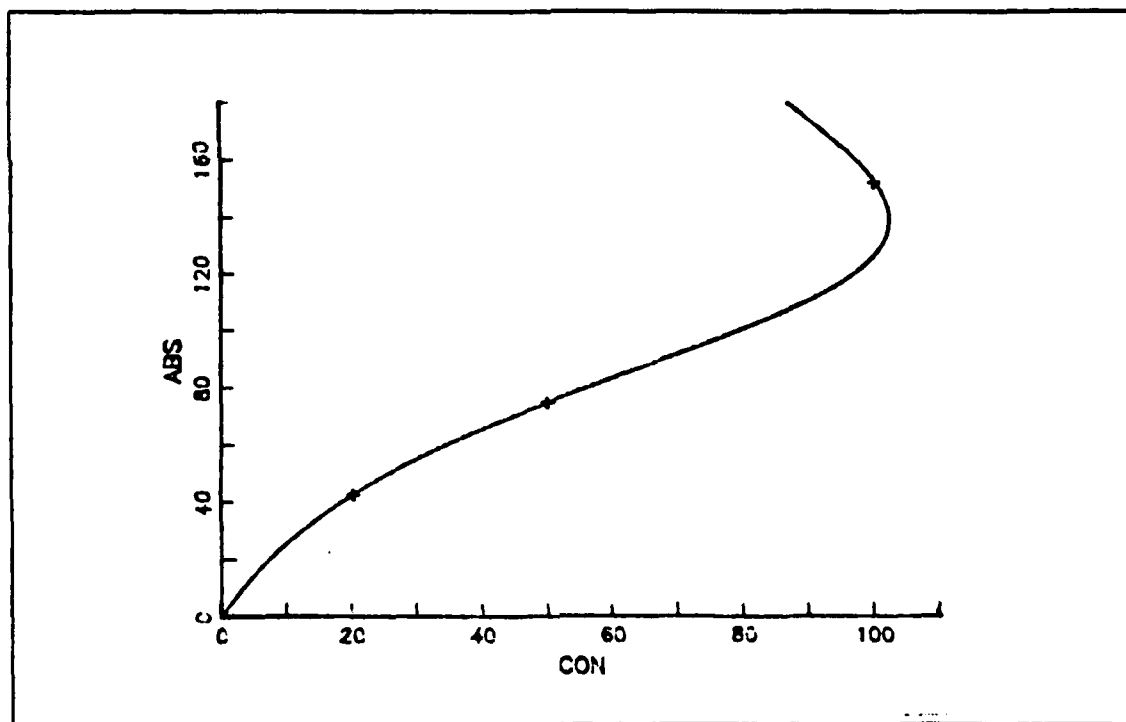


Figure 3.2 Possible Observed Curve from Algorithm 3.

estimated concentration for $A > 140$ with this example.

E. ALGORITHM 4

A simple quadratic equation

$$C = K_1 + K_2 A + K_3 A^2 \quad (3.12)$$

might be a reasonable alternative for a calibration curve.

Equation 3.12 has been expressed as a function of absorbance. This is the reverse of the usual representation where absorbance is considered to be a function of the standard concentrations. However, expressing concentration as the dependent variable facilitates its calculation without the need for extracting roots of equations as is needed for algorithm 5.

This algorithm may give a satisfactory shape, depending on the A values which occur during calibration burns.

It has an undesirable shape for

$$C = (20, 50, 100),$$

$$A = (40, 130, 170),$$

$$K = (43.3, -0.865, 0.007).$$

Pictured in Figure 3.3, here decreasing A, for $A < 60$, gives increasing C.

When

$$C = (20, 50, 100),$$

$$A = (40, 60, 170),$$

$$K = (-59.3, 2.3, -0.008),$$

algorithm 4 gives the curve shown in Figure 3.4, in which increasing A, for $A > 140$, leads to decreasing C.

Note that neither of these two curves pass through the origin, since the polynomial equation will not necessarily give zero concentration with zero absorbance.

In APL the coefficients K are given by

$$K \leftarrow C \left[A \circ . \star 0 \ 1 \ 2 \right]$$

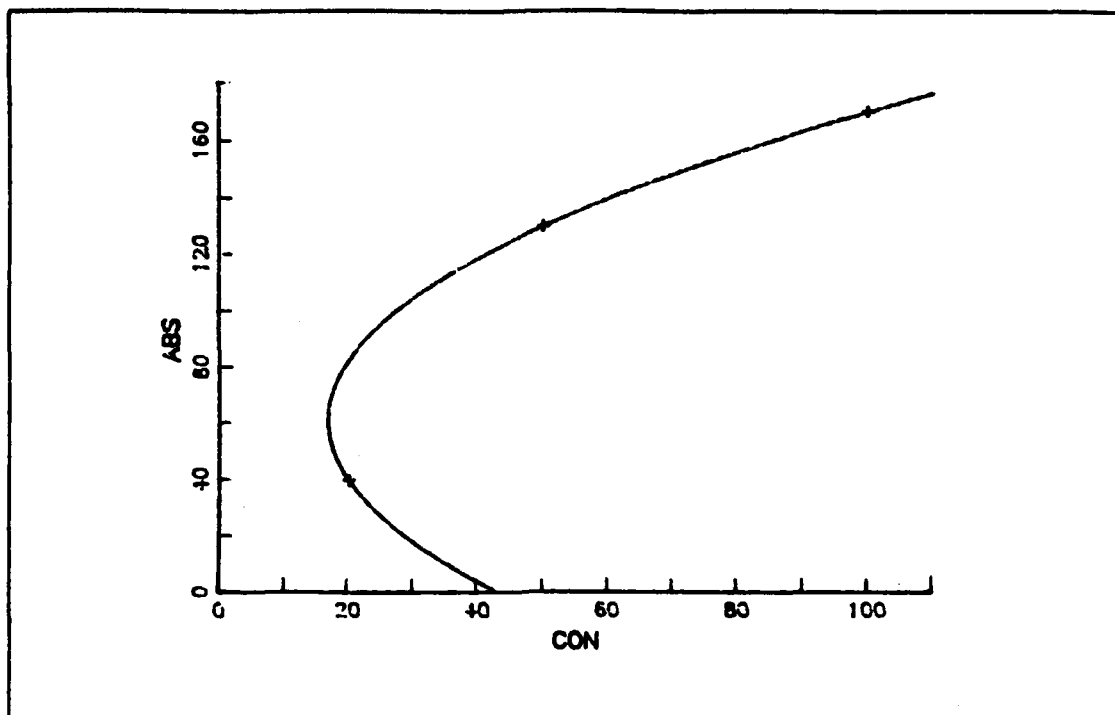


Figure 3.3 The Curve Bending Downward at Lower Concentration by Algorithm 4.

where $C = (c_1, c_2, c_3)$ are the concentrations used and $A = (a_1, a_2, a_3)$ are the observed absorbance numbers.

F. ALGORITHM 5

The final curve to be considered is

$$A = K_1 + K_2 C + K_3 C^2 \quad (3.13)$$

In equation 3.13, absorbance has been expressed as a function of concentration which is the natural representation. However, expressing absorbance as the dependent variable requires the solution of a quadratic equation to find concentration. Equation 3.13 can be expressed like this,

$$K_3 C^2 + K_2 C + K_1 - A = 0 \quad (3.14)$$

Solving the resulting equation for C , we obtain the formula,

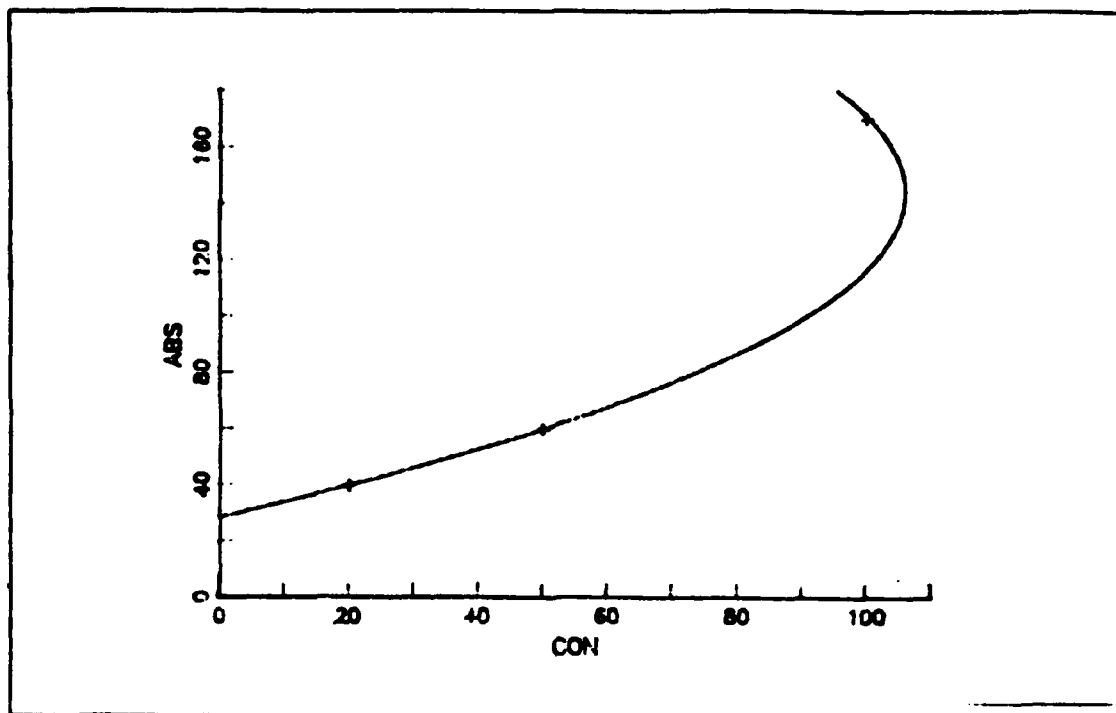


Figure 3.4 The Curve Bending Upward at High Concentration by Algorithm 4.

$$C = \frac{-K_2 \pm \sqrt{K_2^2 - 4K_3(K_1 - A)}}{2K_3} \quad (3.15)$$

Whenever $K_2^2 - 4K_3(K_1 - A)$ is negative, there is no concentration value C for the given absorbance A .

As with algorithm 4, two different undesirable shapes may occur; if the initial calibration data is

$$C = (20, 50, 100),$$

$$A = (40, 140, 170),$$

then

$$K = (-60.8, 5.7, -0.03),$$

giving the curve pictured in Figure 3.5.

If the initial data is

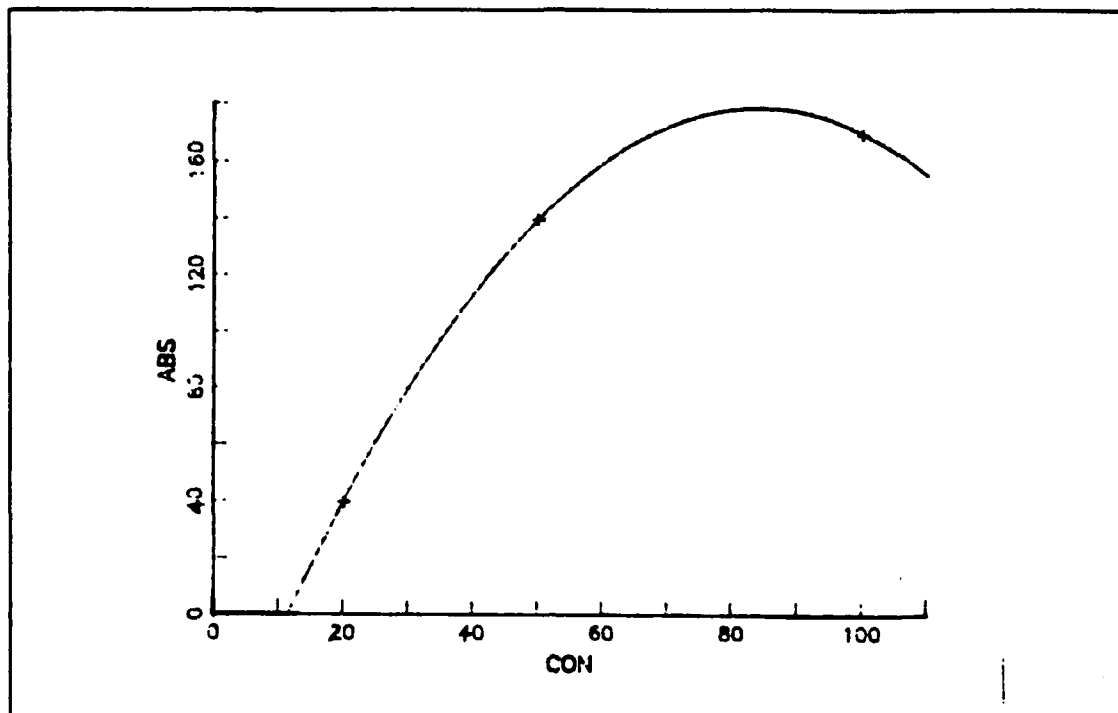


Figure 3.5 The Curve Bending Downward at High Concentration by Algorithm 5.

$$C = (20, 50, 100),$$

$$A = (40, 50, 170),$$

then

$$K = (59.2, -1.5, 0.03),$$

giving the curve in Figure 3.6. With A and C as defined earlier, the coefficients K are given by

$$K \leftarrow A \cdot C^{-1} \cdot \begin{bmatrix} 1 \\ 0 \\ 1 \end{bmatrix}.$$

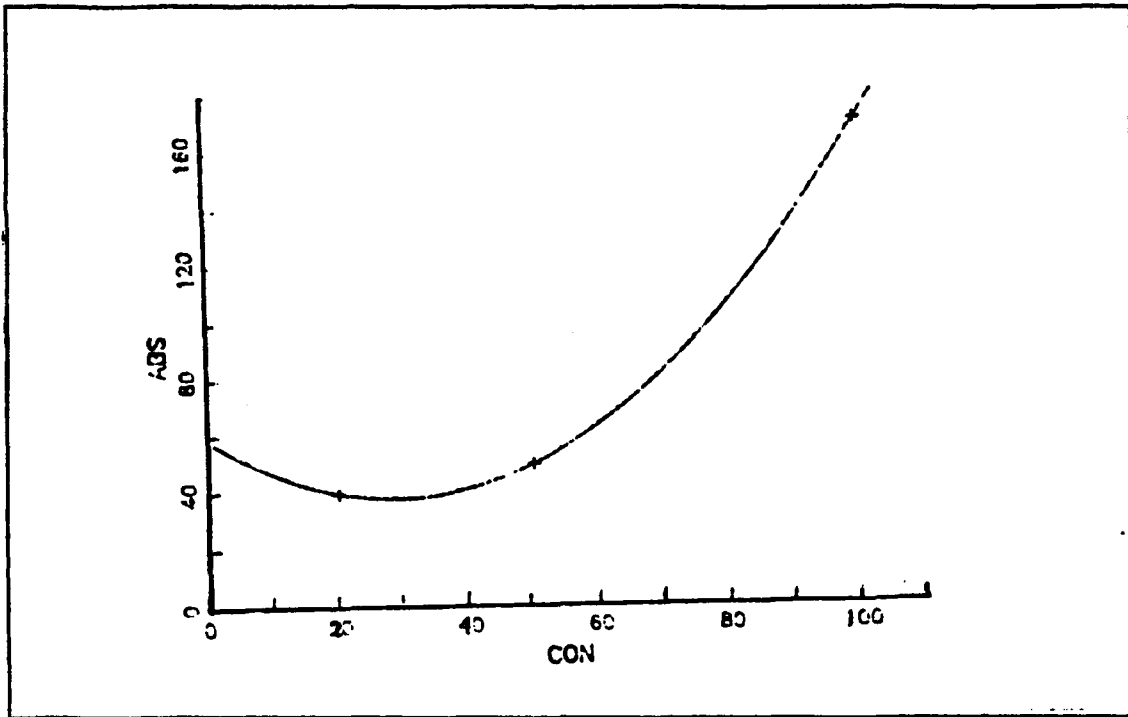


Figure 3.6 The Curve Bending Upward at Lower Concentration by Algorithm 5.

IV. ALGORITHM COMPARISON

A. GENERAL

We have mentioned how algorithms are used to fit the curve and each algorithm has defects for specific A values. Recall that the field test program included the concentrations: 10%, 20%, 40%, 50%, 70% and 100% of full scale for each element. Calibrating the tube consisted of two analyses each using the 20%, 50% and 100% samples; these samples were used to fix the absorbance - concentration curve for all analyses performed until the next calibration. During the planned sequence of burns, each of the 10%, 40%, and 70% verification samples were analyzed twice, once just after the calibration curve was established and then a second time about 30 analyses later. Ideally, the established calibration curve should give 10%, 40% and 70% of full scale respectively for these samples; the different possible algorithms can be compared by using these 10%, 40% and 70% samples, in terms of how close the resulting calibration curve result is to the known concentration, for the various tubes used.

Specifically, each algorithm can be used to establish a calibration curve using the observed absorbance numbers for the 20%, 50% and 100% samples. Since these curves are not, in general, exactly the same for all algorithms, the difference between the returned concentration for a 10% sample, e_{10} say, can be compared with the expected result c_{10} . We will introduce the sum of squares method to compare the algorithms. Thus, for a given algorithm, we will have two e_{10} , two e_{40} , and two e_{70} values from which we will compute the sum of the 6 squared differences. The algorithm which has the smallest mean sum of squares will be selected as the rank 1, with ranks 2, 3, ..., 5 assigned the larger mean sum of squares. The number of rank 1 and rank 2 for each element throughout all tubes will be considered as the comparing tool for that rank. The algorithm which produces the most rank 1 and rank 2 scores will be designated the best one.

B. SUM OF SQUARES METHOD

We will explain the sum of squares method by an example. For tube 401, for element Fe, the fitted curve is shown in Figure 4.1 where the coefficients are computed by algorithm 2.

The residuals E are defined by

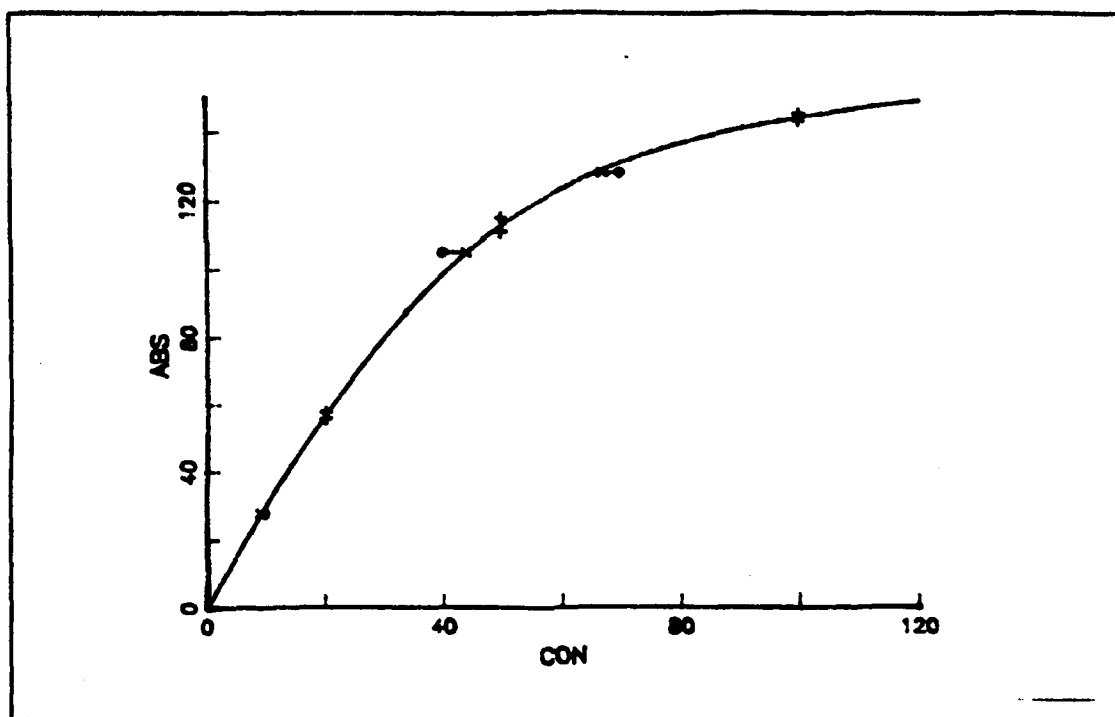


Figure 4.1 The Horizontal Deviation of Residual by Algorithm 2.

$$E = C - C'$$

where C is a known concentration and C' is the estimated concentration produced by the algorithm.

For tube 401, for Fe, the absorbance data of the verification sample A are,

$$A = (28, 105, 129)$$

and the known concentrations are,

$$C = (20, 50, 100)$$

Then we find K by algorithm 2,

$$K = (-0.3211, 0.00602, 0.00159)$$

We can compute the estimated concentration C' from equation 3.4 with (28, 105, 129) in place of A. That is

$$C' = (9.315, 44.004, 67.040).$$

Suppose $E = (e_1, e_2, e_3)$. The residual E is

$$E = C - C' = (0.685, -4.004, 2.96)$$

and the sum of squares SS is defined as the sum of the squares of these residuals,

i. e.,

$$SS = \sum_{i=1}^3 e_i^2 = 0.685^2 + 4.004^2 + 2.96^2 = 25.263. \quad (4.1)$$

Then the mean sum of squares MSS is defined as,

$$MSS = \frac{SS}{3}, \quad (4.2)$$

giving

$$MSS = 25.263 / 3 = 8.42.$$

Residuals at each point are shown in Figure 4.1 where the residual is indicated by the horizontal line between symbol 'x' and 'o'. Estimated concentration C' is indicated by the symbol 'x' and concentration is indicated by the symbol 'o'.

When we are using algorithm 4 with the same data, then

$$K = (79.237, -1.775, 0.013)$$

This curve is shown in Figure 4.2. Horizontal deviation is the distance between symbol 'x' and 'o', which is shown as a horizontal line. The estimated concentrations are

$$C' = (39.75, 36.57, 67.184).$$

As we see, the estimated concentration at absorbance 28 is larger than at absorbance 105, because the curve is bent downward at the lower concentration with this algorithm. Therefore, the residuals are larger. Also, algorithm 2 is considerably better than algorithm 4 for this set of data.

Accordingly,

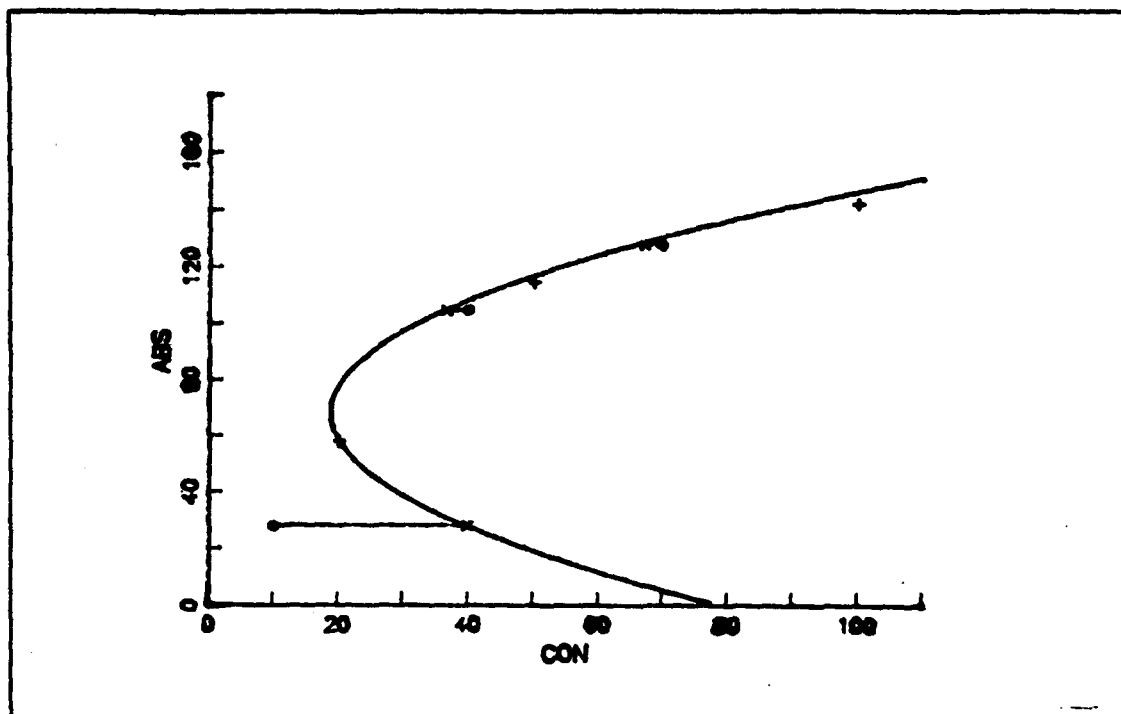


Figure 4.2 The Horizontal Deviation of Residual by Algorithm 4.

$$E = (-29.75, 3.43, 2.816),$$

$$SS = (-29.75)^2 + 3.43^2 + 2.816^2 = 735.257,$$

and

$$MSS = 735.257/3 = 245.086.$$

For these data with these two algorithms,

$$MSS \text{ by algorithm 2} = 8.42,$$

$$MSS \text{ by algorithm 4} = 245.086,$$

and algorithm 2 is better than algorithm 4.

Estimated concentrations C' by 6 algorithms for Fe are summarized in Table 4. Known concentrations C , shown in the first column, were used 4 times 10%, 40% and 70% of full scale. After the calibration curve was established, each of the 10%, 40%, and 70% verification sample was analyzed twice. For tube 401 this calibration was

done twice, so that 12 verification data are gathered in Table 4. Notice the differences between the estimated concentrations in column 2 and 3 in Table 4. The estimated concentrations by PWMA and by algorithm 1 are very close. Since algorithm PWMA and algorithm 1 are exactly the same as noted earlier; the PWMA concentration values are those printed out by the microprocessor, while the algorithm values are computed from the curve established from the rounded absorbance values. Since algorithms 2 through 5 must be employed with the rounded absorbance values, rather than the single precision values available to the microprocessor the rank comparisons will be restricted to only algorithms 1 through 5. The mean square residuals will be computed for the PWMA, but they will not be used in comparison with the others; cases in which the PWMA actually gives the smallest mean square will be denoted by an asterisk(*). We might expect that algorithm PWMA and algorithm 1 should perform very similarly. In the last column, .00 means we can not obtain acceptable estimated concentrations with this algorithm. If $K_2^2 - 4 K_3 (K_1 - A)$ was negative, the estimated concentration was forced to be .00 for all verification samples with this calibration. The estimated concentration in the first three rows of algorithm 2 and 4 are the same as our previous example.

From the data in Table 4, we calculate the sum of squares and the mean sum of squares, which are displayed in Table 5. It says that the sum of squares by PWMA is 100, and the mean sum of squares is 8.33, so the rank of the mean sum of squares is labeled with * as it is the smallest. If we don't count the PWMA in comparison, algorithm 3 is the best. It has the smallest mean sum of squares, 9.09; thus algorithm 3 suggested by Limbeck is the best for the tube PPM401 for element Fe. The number of acceptable data points for each algorithm is given in the row labeled N; if the test plan was followed exactly, and the algorithm performed satisfactorily for the calibration samples, N would equal 12 for each case. If the algorithm performed totally unsatisfactorily, its N value is 0 and its mean sum of squares was set equal to 888 (see algorithm 5 in Table 5).

The sum of squares, the mean sum of squares, and the rank of the algorithm for tube 401 for 9 elements are summarized in Table 6. The equivalent data for the other tubes are attached in Appendix D.

Table 6 shows the best algorithm for each element. Algorithm 3 is the best algorithm for element Fe. Algorithm 1 has the smallest mean sum of squares for Ag, and so on. As can be seen, the best algorithm differs according to the element. For tube 401, no single algorithm is the best for all the elements.

TABLE 4
ESTIMATED CONCENTRATION BY 6 ALGORITHMS

Standard	Algorithm					
	PWMA	1	2	3	4	5
10.00	9.00	8.87	9.31	13.44	39.75	.00
40.00	44.00	44.79	44.00	45.66	36.57	.00
70.00	70.00	70.23	67.04	67.96	67.18	.00
10.00	13.00	13.21	13.57	17.64	29.09	.00
40.00	42.00	42.32	41.36	43.08	32.93	.00
70.00	70.00	70.23	67.04	67.96	67.18	.00
10.00	8.00	8.00	8.34	17.04	53.22	.00
40.00	43.00	43.05	42.64	60.82	36.86	.00
70.00	74.00	75.32	74.43	88.77	83.90	.00
10.00	6.00	6.07	6.39	13.58	66.29	.00
40.00	36.00	35.69	35.19	52.89	26.16	.00
70.00	73.00	73.52	72.64	87.41	81.58	.00

TABLE 5
ALGORITHM COMPARISON FOR FE

	Algorithm					
	PWMA	1	2	3	4	5
SS(Fe)	100.00	128.08	121.12	109.09	3119.08	.00
N	12.00	12.00	12.00	12.00	12.00	.00
MSS	8.33	10.67	10.09	9.09	259.92	888.00
RANK	*	3	2	1	4	5

In Chapter II we mentioned that Langley used 6 tubes, Myrtle Beach used 5 tubes, Pensacola used 11 tubes, and Elmendorf used 6 tubes. Due to missing data, we excluded several tubes in our analyses. For example, tube PPM406 doesn't have verification sample 70 data. For similar reasons we discard 1 tube from Langley, 1

TABLE 6
SUMMARY OF THE RANKS FOR TUBE 401

	Algorithm					
	PWMA	1	2	3	4	5
SS(Fe) n MSS RANK	100.00 12.00 8.33 *	128.03 12.00 10.67 3	121.08 12.00 10.09 2	109.09 12.00 9.09 1	3119.08 12.00 259.92 4	.00 .00 888.00 5
SS(Ag) n MSS RANK	2.00 12.00 .17	12.85 12.00 .07 1	1.70 12.00 .14 4	1.48 12.00 .12 3	12.34 12.00 1.03 5	.67 6.00 .11 2
SS(Al) n MSS RANK	268.00 12.00 22.33 *	317.88 12.00 26.49 3	300.21 12.00 25.02 2	297.57 12.00 24.80 1	429.29 12.00 35.77 4	.00 .00 888.00 5
SS(Cr) n MSS RANK	16.00 12.00 1.33 *	17.29 12.00 1.44 3	17.16 12.00 1.43 2	17.10 12.00 1.42 1	39.32 12.00 3.28 4	.00 .00 888.00 5
SS(Cu) n MSS RANK	106.00 12.00 8.83	102.05 12.00 8.50 4	100.06 12.00 8.34 2	99.74 12.00 8.31 1	100.09 12.00 8.34 3	107.97 12.00 9.00 5
SS(Mg) n MSS RANK	91.00 12.00 7.58 *	170.50 12.00 14.21 3	156.21 12.00 13.02 2	148.12 12.00 12.34 1	179.10 12.00 14.92 4	.00 .00 888.00 5
SS(Ni) n MSS RANK	38.00 12.00 3.17	34.87 12.00 2.91 3	26.02 12.00 2.17 1	31.70 12.00 2.64 2	192.41 12.00 16.03 4	.00 .00 888.00 5
SS(Si) n MSS RANK	162.00 12.00 13.50	149.09 12.00 12.42 2	141.29 12.00 11.77 1	166.29 12.00 13.86 3	171.89 12.00 14.32 4	103.04 6.00 17.17 5
SS(Ti) n MSS RANK	135.00 12.00 11.25	142.77 12.00 11.90 4	364.12 12.00 30.34 5	141.42 12.00 11.79 2	134.41 12.00 11.20 1	141.47 12.00 11.79 3

tube from Myrtle Beach, 2 tubes from Pensacola, leaving 24 tubes in all. The summary of 24 tubes for each element is shown in Table 7 through Table 15. We may say that algorithms which have the most rank 1 or rank 2 scores but the least rank 4 or rank 5

TABLE 7
RANKS OF 24 TUBES FOR FE

Algo.	rank					
	*	1	2	3	4	5
PWMA	10					
1		7	1	11	5	0
2		7	8	3	0	5
3		5	12	6	1	0
4		1	0	0	12	11
5		4	3	3	6	8

TABLE 8
RANKS OF 24 TUBES FOR AG

Algo.	rank					
	*	1	2	3	4	5
PWMA	6					
1		7	3	4	8	2
2		7	8	5	2	2
3		0	7	11	6	0
4		1	3	1	5	13
5		9	3	3	3	6

scores are good ones, and algorithms which have the least rank 1 or rank 2 scores but the most rank 4 or rank 5 scores are not good ones. The entries in these tables are the counts of the ranks. For example, in Table 7 algorithm 1 for Fe has 7 rank 1's, 1 rank 2, 11 rank 3's, 5 rank 4's, and 0 rank 5 out of 24. For Fe, the PWMA internal algorithm, with rounded concentrations, had the smallest mean sum of squares 10 time out of 24. In Table 7, we conclude that algorithm 1 is better than algorithms 2, 3, 4, and 5. Algorithm 2 has 7 rank 1's but also 5 rank 5's. Therefore, algorithm 2 is not as good as algorithm 1. If algorithm 2 did not have those undesirable cases with

TABLE 9
RANKS OF 24 TUBES FOR AL

Algo.	rank					
	*	1	2	3	4	5
PWMA	12					
1		2	8	6	4	4
2		10	4	7	3	0
3		6	8	9	1	0
4		3	1	1	10	9
5		3	3	1	6	11

TABLE 10
RANKS OF 24 TUBES FOR CR

Algo.	rank					
	*	1	2	3	4	5
PWMA	9					
1		9	2	7	5	1
2		4	6	10	2	2
3		3	15	4	2	0
4		0	0	2	11	11
5		8	1	0	4	10

denominator zero, then it would be the best. Also algorithm 4 has 11 rank 5's because it has a lot of deviations for two extreme cases as we saw in the previous Chapter. Algorithm 5 is almost the same as algorithm 4. Algorithm 3, with 5 rank 1's and 12 rank 2's, is acceptable but not as good as algorithm 1. We conclude that algorithm 1 is the best for element Fe.

For element Ag, even though algorithm 5 has 9 rank 1's, it also has 6 rank 5's. Therefore, we can not select algorithm 5 as the best one. Algorithm 2, or 1 are acceptable for element Ag.

TABLE 11
RANKS OF 24 TUBES FOR CU

Algo.	rank					
	*	1	2	3	4	5
PWMA	15					
1		12	1	1	9	1
2		4	8	5	6	1
3		1	7	9	0	2
4		2	2	3	4	13
5		5	6	1	5	7

TABLE 12
RANKS OF 24 TUBES FOR MG

Algo.	rank					
	*	1	2	3	4	5
PWMA	12					
1		12	5	6	0	1
2		4	7	11	2	0
3		6	9	4	5	0
4		0	2	3	14	5
5		2	1	0	3	18

For Al, algorithm 2 dominates the other algorithms, since it has 10 rank 1's. Algorithm 3 is the second algorithm. Algorithm 3 could be useable even though it has fewer rank 1 scores.

For Cr, algorithm 4 is not acceptable, for it has 22 tubes rank 4's or 5's. Algorithm 1 is the best for element Cr.

For Cu, Mg, and Ni, algorithm 1 dominates the others. For Si, algorithm 2 or 3 are the best. For Ti, algorithm 2 is good. Algorithm 1 and 4 are acceptable too.

TABLE 13
RANKS OF 24 TUBES FOR NI

Algo.	rank					
	*	1	2	3	4	5
PWMA	7					
1		13	2	8	1	0
2		4	2	2	5	11
3		4	18	2	0	0
4		1	1	3	16	3
5		2	1	9	2	10

TABLE 14
RANKS OF 24 TUBES FOR SI

Algo.	rank					
	*	1	2	3	4	5
PWMA	11					
1		3	7	8	3	2
2		8	7	2	3	4
3		7	2	3	11	1
4		1	7	9	4	3
5		4	1	2	3	13

Table 16 summarizes the results of the previous tables by listing the best and alternate algorithm for each element. We can say that algorithm 1 provides the best fit for elements Fe, Cr, Cu, Mg, and Ni, and algorithm 2 provides the best fit for elements Ag, Al, Si, and Ti.

TABLE 15
RANKS OF 24 TUBES FOR TI

Algo.	rank					
	*	1	2	3	4	5
PWMA	11					
1		5	9	3	5	2
2		7	3	2	8	4
3		4	1	6	7	6
4		5	5	11	2	1
5		3	5	3	2	11

TABLE 16
THE BEST ALGORITHM FOR EACH ELEMENT

	Element								
	Fe	Ag	Al	Cr	Cu	Mg	Ni	Si	Ti
First	1	2	2	1	1	1	1	2	2
Second	3	1	3	3	2	3	3	3	1,4

V. CONCLUSIONS AND RECOMMENDATIONS

There is no single best algorithm for all nine elements, since the degree of curvature is different according to the element. The PWMA didn't consider different algorithms for different elements. The computation in the PWMA used only by one algorithm for all nine elements. If we employ different algorithms for different elements in an analyzer, it will be better because each element has different characteristics. We recommend that one may use algorithm 1 for elements Fe, Cr, Cu, Mg, and Ni and algorithm 2 for elements Ag, Al, Si, and Ti to obtain better estimation in computing concentration. The microprocessor can easily check for the existence of an asymptote. If one occurs, use algorithm 1.

APPENDIX A

PWMA FIELD TEST BURN SEQUENCE

Day 1

1. Tube Condition
2. Auto Zero - Operate in AZ mode. If the AZ failed, repeat the AZ and number the run as 2A. If the AZ failed twice in a row, replace the tube.
3. CAL Ia - Operate in CAL I mode.
4. CAL Ib - Operate in CAL I mode.
5. CAL IIa - Operate in CAL II mode.
6. CAL IIb - Operate in CAL II mode.
7. CAL IIIa - Operate in CAL III mode.
8. CAL IIIb - Operate in CAL III mode. If the calibration failed, repeat the calibration and number the calibration runs as 3A, ..., 8A, accordingly.
9. Blank - Run a burn with no sample; operate in Analysis mode. From now on, all the samples will be run in Analysis mode except Reslope.
10. Verification Sample # 1: 10%
11. Verification Sample # 2: 40%
12. Verification Sample # 3: 70%
13. Blank
- 14-18. Correlation sample # 1 - Repeat the same sample five times, this procedure applies for all the correlation sample.
- 19-23. Correlation sample # 2
- 24-28. Correlation sample # 3
29. Blank
- 30-37. Random samples - These may be real engine oil sample. If there are not enough real engine oil samples, repeat one sample as many times as necessary. (8 samples total)
38. 10% sample
39. 40% sample
40. 70% sample

Day 2

41. Blank
42. Reslope A - Operate in RSLP mode. If the reslope A failed, disregard and continue the sequence.
43. Reslope B - Operate in RSLP mode. If the reslope B failed, disregard and continue the sequence.
44. 10% sample
45. 40% sample
46. 70% sample
- 47-51. Correlation sample # 4
- 52-56. Correlation sample # 5
57. Blank
- 58-63. Random samples (6 samples total)
64. CAL I check sample - Operate in Analysis mode.
65. CAL II check sample - Operate in Analysis mode.
66. CAL III check sample - Operate in Analysis mode.
67. Blank
- 68-77. Random samples (10 samples total)
78. 10% sample
79. 40% sample
80. 70% sample

Day 3

81. Blank
82. CAL Ia.
83. CAL Ib.

84. CAL IIa.
85. CAL IIb.
86. CAL IIIa.
87. CAL IIIb.
88. Blank
89. 10% sample
90. 40% sample
91. 70% sample
92-96. Correlation sample # 6
97-101. Correlation sample # 7
102. Blank
103-117. Random samples (15 samples total)
118. 10% sample
119. 40% sample
120. 70% sample

Day 4

121. Blank
122. Reslope A
123. Reslope B
124. 10% sample
125. 40% sample
126. 70% sample
127-131. Correlation sample # 8
132-136. Correlation sample # 9
137. Blank
138-157. Random samples (20 samples total)
158. CAL I check sample
159. CAL II check sample
160. CAL III check sample

APPENDIX B DATA PPM401

There are several data, we list just one of them for references.

PPM401											
PS23 401											
CAL IA	P	3	58	53	10	38	23	44	40	10	10
CAL IB	P	4	56	52	10	34	22	46	38	8	9
CAL IIA	P	5	115	98	19	83	47	60	92	28	30
CAL IIB	P	6	111	100	21	84	43	61	96	18	29
CAL IIIA	P	7	143	129	26	104	70	72	119	33	42
CAL IIIB	P	8	145	131	24	104	71	74	120	35	44
VERIF 10	P	10	9	1	2	1	4	3	3	2	5
VERIF 10	ABSORBANCES		28	28	5	17	12	32	19	4	11
VERIF 40	P	11	44	4	10	4	18	9	12	9	7
VERIF 40	ABSORBANCES		105	90	17	70	42	56	79	20	20
VERIF 70	P	12	70	7	9	6	28	17	20	12	10
VERIF 70	ABSORBANCES		129	114	16	88	56	66	105	26	29
VERIF 10	P	38	13	1	8	1	8	2	4	9	8
VERIF 10	ABSORBANCES		40	28	15	25	22	30	25	21	21
VERIF 40	P	39	42	4	7	5	16	8	12	15	7
VERIF 40	ABSORBANCES		101	84	14	77	39	54	76	30	18
VERIF 70	P	40	70	7	11	7	26	19	21	12	12
VERIF 70	ABSORBANCES		129	115	18	94	54	68	107	26	34
PS24 401											
RESLOPE 1	P	42	117	101	12	72	51	55	80	37	19
RESLOPE 2	P	43	114	96	13	69	51	54	80	34	17
VERIF 10	P	44	12	1	1	1	4	2	3	4	3
VERIF 10	ABSORBANCES		36	29	3	17	13	28	19	10	7
VERIF 40	SI>LIMIT P	45	46	4	8	4	20	8	11	25	8
VERIF 40	ABSORBANCES		107	91	15	74	45	53	70	42	21
VERIF 70	SI>LIMIT P	46	74	9	10	7	31	14	17	39	10
VERIF 70	ABSORBANCES		131	124	16	95	61	63	99	67	30
CAL I	P	64	23	2	3	2	8	4	5	8	4
CAL I	ABSORBANCES		65	52	7	38	23	42	34	17	10
CAL II	P	65	50	5	9	6	20	14	14	14	9
CAL II	ABSORBANCES		113	98	16	88	44	61	89	28	25
CAL III	CR,NI,SI,P	66	103	10	28	11	35	25	31	30	23
CAL III	ABSORBANCES		149	130	28	110	65	74	124	50	50
VERIF 10	SI>LIMIT?P	78	9	1	2	1	4	2	2	0	3
VERIF 10	ABSORBANCES		28	24	5	21	12	30	15	117	6
VERIF 40	P	79	41	4	11	5	15	11	12	17	8
VERIF 40	ABSORBANCES		99	80	18	79	37	59	76	31	23
VERIF 70	SI>LIMIT P	80	73	7	22	7	25	23	21	23	13
VERIF 70	ABSORBANCES		131	114	24	95	53	72	107	39	34

**APPENDIX C
PROGRAMMING LIST**

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V COEF1
[1] A 8/12/86 PROGRAM TO FIT BILL'S 2 POINT RATIONAL EQUATION, CHECK OUT
[2] A VERIFICATION READINGS AND THE VALUES AFTER THE RESLOPE AS WELL.
[3] H+ 10.10.250.10.40.250.30.20.2
[4] E+20+5*19*I+1
[5] T1+ 83#61p0
[6] R1+ 30#61p0
[7] Q1+ 60#61p0
[8] 'FOR TUBE 401'
[9] 'MODEL A/(K1+K2A)!'
[10] LOOP:A++/ 3 2 p (A,PPM401[1+16;E[I]+15])+2
[11] K+((A[1 2]+20 50 *H[I])#A[1 2]#. * 0 1),[0.1](A[2 3]+ 50 100
    #H[I]) A[2 3]#. * 0 1
[12] A ORIGINAL COEFFICIENTS FOR THE TWO CURVES.
[13] D+ A,PPM401[7+112;E[I]+15]
[14] O+D[2*16]+(,(D[2*16]#. * 0 1)+. *K)[1 3 6 7 9 12]
[15] U1+ 83#61p0
[16] Q+ 60#61p0D[-1+2*16]
[17] Q1+Q1,[1] Q
[18] U+(60#61p0D[-1+2*16]),U1
[19] T1+T1,[1] U1
[20] R+ 30#61p 10 40 70 *H[(E[I]+5)-4]
[21] R1+R1,[1] R
[22] 'FOR ELEMENT: '(EL[I:]),' PRIOR TO RESLOPE'
[23] U+U+(30#61p 10 40 70 *H[(E[I]+5)-4]),U
[24] C++/(A,PPM401[20+12;E[I]+15])+2
[25] C+50*H[I]+(2pC)+(K+. *C* 0 1)
[26] A GIVES MULTIPLICATIVE CORRECTION FACTORS FOR THE TWO LINES
[27] D+ A,PPM401[(22+16),34+16;E[I]+15]
[28] O+C[6p 1 1 2]*D[2*16]+(,(D[2*16]#. * 0 1)+. *K)[1 3 6 7 9 12]
[29] 'AFTER THE RESLOPE'
[30] U+(60#61p0D[-1+2*16]), 83#61p0
[31] Q+ A,PPM401[(28+16);E[I]+15]
[32] O+C[1 1 2 2]*D[2*1 2 2 3]+(,(D[2*1 2 2 3]#. * 0 1)+. *K)[1 3 4 6]
[33] U+( 30+U),[1](((60#4 1 pD[1 3 3 5]), 83#4 1 pQ),[1] 30+U)
[34] U+U+(30#10 1 p 10 40 70 20 50 50 100 10 40 70 *H[(E[I]+5)-4]),U
[35] +((I+I+1)<=9)/LOOP
[36] I+1
[37] LOOT:A++/ 3 2 p (A,PPM401[41+16;E[I]+15])+2
[38] K+((A[1 2]+20 50 *H[I])#A[1 2]#. * 0 1),[0.1](A[2 3]+ 50 100
    #H[I]) A[2 3]#. * 0 1
[39] D+ A,PPM401[47+112;E[I]+15]
[40] O+D[2*16]+(,(D[2*16]#. * 0 1)+. *K)[1 3 6 7 9 12]
[41] U1+ 83#61p0
[42] Q+ 60#61p0D[-1+2*16]
[43] Q1+Q1,[1] Q
[44] U+(60#61p0D[-1+2*16]),U1
[45] T1+T1,[1] U1
[46] R+ 30#61p 10 40 70 *H[(E[I]+5)-4]
[47] R1+R1,[1] R
[48] 'FOR ELEMENT: '(EL[I:]),' PRIOR TO RESLOPE'
[49] U+U+(30#61p 10 40 70 *H[(E[I]+5)-4]),U
[50] +((I+I+1)<=9)/LOOT

```

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V COEFR
[1] A 8/12/86 PROGRAM TO FIT BILL'S 2 POINT RATIONAL EQUATION, CHECK OUT
[2] A VERIFICATION READINGS AND THE VALUES AFTER THE RESLOPE AS WELL.
[3] H+ 10.10.250.10.40.250.30.20.2
[4] E+20+5*19*I+1
[5] T2+ 83#61p0
[6] 'FOR TUBE 401'

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[7] LOOP:A++/ 3 2 p (a, PPM401[1+16;E[I]+15])+2
[8] K+((A°.x 1 -1),A*2)±(20 50 100 xH[I])°. * 1 0 1
[9] □+K+3 1 p (BK)+.x 1 1 1
[10] ORIGINAL COEFFICIENTS FOR THE TWO CURVES.
[11] D+a, PPM401[7+12;E[I]+15]
[12] O+((K[1;]×D[2×16])+K[3;]×D[2×16]×D[2×16])+((K[2;]×D[2×16])-1)
[13] U2+8 3 # 6 1 pO
[14] U+(7 1 # 6 1 pD[-1+2×16]),U2
[15] T2+T2, [1] U2
[16] 'FOR ELEMENT: ', (EL[I;]), ' PRIOR TO RESLOPE'
[17] □+U+(4 1 # 6 1 p 10 40 70 xH[(E[I]+5)-4]),U
[18] C+/(a, PPM401[20+12;E[I]+15])+2
[19] C+50×H[I]+(K[1;]×C+K[3;]×C×C)+((K[2;]×C)-1)
[20] GIVES MULTIPLICATIVE CORRECTION FACTORS FOR THE TWO LINES
[21] D+a, PPM401[(22+16),34+16;E[I]+15]
[22] O+C[6p1]×(K[1;]×D[2×16]+K[3;]×D[2×16]×D[2×16])+((K[2;]×D[2
6]) -1)
[23] 'AFTER THE RESLOPE'
[24] U+(7 1 # 6 1 pD[-1+2×16]), 8 3 # 6 1 pO
[25] D+a, PPM401[(28+16);E[I]+15]
[26] O+C[4p1]×(K[1;]×D[2 4 4 6]+K[3;]×D[2 4 4 6]×D[2 4 4 6])+((K[
2;] D[2 4 4 6]) -1)
[27] →((1/O≥2000)=1)/ERR
[28] U+(3 0 +U), [1] ((7 1 # 4 1 pD[1 3 3 5]), 8 3 # 4 1 pO), [1] 3 0 +U)
[29] □+U+(3 0 # 10 1 p 10 40 70 20 50 50 100 10 40 70 xH[(E[I]+5)-4]),U
[30] →((I+I+1)≤9)/LOOP
[31] I+1
[32] LOOT:A++/ 3 2 p (a, PPM401[41+16;E[I]+15])+2
[33] K+((A°.x 1 -1),A*2)±(20 50 100 xH[I])°. * 1 0 1
[34] □+K+3 1 p (BK)+.x 1 1 1
[35] ORIGINAL COEFFICIENTS FOR THE TWO CURVES.
[36] D+a, PPM401[47+12;E[I]+15]
[37] O+((K[1;]×D[2×16])+K[3;]×D[2×16]×D[2×16])+((K[2;]×D[2×16])-1)
[38] U2+8 3 # 6 1 pO
[39] U+(7 1 # 6 1 pD[-1+2×16]),U2
[40] T2+T2, [1] U2
[41] 'FOR ELEMENT: ', (EL[I;]), ' PRIOR TO RESLOPE'
[42] □+U+(4 1 # 6 1 p 10 40 70 xH[(E[I]+5)-4]),U
[43] →((I+I+1)≤9)/LOOT
[44] →0
[45] ERR:O+4pO
[46] U+(3 0 +U), [1] ((7 1 # 4 1 pD[1 3 3 5]), 8 3 # 4 1 pO), [1] 3 0 +U)
[47] □+U+(4 1 # 10 1 p 10 40 70 20 50 50 100 10 40 70 xH[(E[I]+5)-4]),U
[48] →((I+I+1)≤9)/LOOP

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▽ COEF3
[1] A 9/06/86 PROGRAM TO FIT BILL'S 2 POINT RATIONAL EQUATION, CHECK OUT
[2] A VERIFICATION READINGS AND THE VALUES AFTER THE RESLOPE AS WELL.
[3] A MODEL C = A+(C1+C2×A+C3×A×A)
[4] 'MODEL C = A+(C1+C2×A+C3×A×A)'
[5] H+1 0.1 0.25 0.1 0.4 0.25 0.3 0.2 0.2
[6] T3+8 3 # 6 1 pO
[7] E+20+5×19×I+1
[8] 'FOR TUBE 401'
[9] LOOP:A++/ 3 2 p (a, PPM201[1+16;E[I]+15])+2
[10] □+K+(A+C+ 20 50 100 xH[I])A°. * 0 1 2
[11] ORIGINAL COEFFICIENTS FOR THE TWO CURVES.
[12] D+a, PPM401[7+12;E[I]+15]
[13] O+D[2×16]+((D[2×16]°. * 0 1 2)+.x BK)
[14] U3+8 3 # 6 1 pO
[15] U+(6 0 # 6 1 pD[-1+2×16]),U3
[16] T3+T3, [1] U3
[17] 'FOR ELEMENT: ', (EL[I;]), ' PRIOR TO RESLOPE'
[18] □+U+(3 0 # 6 1 p 10 40 70 xH[(E[I]+5)-4]),U
[19] C+/(a, PPM401[20+12;E[I]+15])+2
[20] C+50+(1pC)+(K+.x C* 0 1 2)
[21] GIVES MULTIPLICATIVE CORRECTION FACTORS FOR THE TWO LINES
[22] D+a, PPM401[(22+16),34+16;E[I]+15]
[23] O+C[6p1]×D[2×16]+((D[2×16]°. * 0 1 2)+.x BK)

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[24] 'AFTER THE RESLOPE'
[25] U+(6 0 # 6 1 p D[ -1+2x16]), 8 3 # 6 1 p 0
[26] D+# PPM401[(28+16);E[I]+15]
[27] O+C[4p1]x D[2x1 2 2 3]+((D[2x1 2 2 3]o.* 0 1 2)+.x#K)
[28] U+(3 0 +U),[1](((6 0 # 4 1 p D[1 3 3 5]), 8 3 # 4 1 p 0),[1] 3 0 +U)
[29] #+U+(3 0 # 10 1 p 10 40 70 20 50 50 100 10 40 70 xH[(E[I]+5)-4]),U
[30] #((I+I+1)≤9)/LOOP
[31] I+1
[32] LOOP:A++/ 3 2 p (#,PPM201[41+16;E[I]+15])+2
[33] #+K+(A+C+ 20 50 100 xH[I])#A.* 0 1 2
[34] # ORIGINAL COEFFICIENTS FOR THE TWO CURVES.
[35] D+# PPM401[47+112;E[I]+15]
[36] O+D[2x16]+((D[2x16]o.* 0 1 2)+.x#K)
[37] U3+ 8 3 # 6 1 p 0
[38] U+(6 0 # 6 1 p D[ -1+2x16]),U3
[39] T3+T3,[1] U3
[40] 'FOR ELEMENT:',(EL[I;]),'. PRIOR TO RESLOPE'
[41] #+U+(3 0 # 6 1 p 10 40 70 xH[(E[I]+5)-4]),U
[42] #((I+I+1)≤9)/LOOP

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▽ COEF4
[1] # 9/09/86 PROGRAM TO FIT BILL'S 2 POINT RATIONAL EQUATION, CHECK OUT
[2] # VERIFICATION READINGS AND THE VALUES AFTER THE RESLOPE AS WELL.
[3] # MODEL C = C1+C2xA+C3xAxA
[4] 'MODEL C = C1+C2xA+C3xAxA'
[5] H+ 1 0. 1 0. 25 0. 1 0. 4 0. 25 0. 3 0. 2 0. 2
[6] E+20+5x19xI+1
[7] T4+ 8 3 # 6 1 p 0
[8] 'FOR TUBE 401'
[9] LOOP:A++/ 3 2 p (#,PPM501[1+16;E[I]+15])+2
[10] #+K+(C+ 20 50 100 xH[I])#A.* 0 1 2
[11] # ORIGINAL COEFFICIENTS FOR THE TWO CURVES.
[12] D+# PPM401[7+112;E[I]+15]
[13] O+((D[2x16]o.* 0 1 2)+.xK)
[14] U4+ 8 3 # 6 1 p 0
[15] T4+T4,[1] U4
[16] U+(6 0 # 6 1 p D[ -1+2x16]),U4
[17] 'FOR ELEMENT:',(EL[I;]),'. PRIOR TO RESLOPE'
[18] #+U+(3 0 # 6 1 p 10 40 70 xH[(E[I]+5)-4]),U
[19] C++/(#,PPM401[20+12;E[I]+15])+2
[20] C+(50xH[I])+(K+.xC* 0 1 2)
[21] # GIVES MULTIPLICATIVE CORRECTION FACTORS FOR THE TWO LINES
[22] D+# PPM401[(22+16),34+16;E[I]+15]
[23] O+Cx((D[2x16]o.* 0 1 2)+.xK)
[24] 'AFTER THE RESLOPE'
[25] U+(6 0 # 6 1 p D[ -1+2x16]), 8 3 # 6 1 p 0
[26] D+# PPM401[(28+16);E[I]+15]
[27] O+Cx((D[2x1 2 2 3]o.* 0 1 2)+.xK)
[28] U+(3 0 +U),[1](((6 0 # 4 1 p D[1 3 3 5]), 8 3 # 4 1 p 0),[1] 3 0 +U)
[29] #+U+(3 0 # 10 1 p 10 40 70 20 50 50 100 10 40 70 xH[(E[I]+5)-4]),U
[30] #((I+I+1)≤9)/LOOP
[31] I+1
[32] LOOP:A++/ 3 2 p (#,PPM501[41+16;E[I]+15])+2
[33] #+K+(C+ 20 50 100 xH[I])#A.* 0 1 2
[34] # ORIGINAL COEFFICIENTS FOR THE TWO CURVES.
[35] D+# PPM401[47+112;E[I]+15]
[36] O+((D[2x16]o.* 0 1 2)+.xK)
[37] U4+ 8 3 # 6 1 p 0
[38] T4+T4,[1] U4
[39] U+(6 0 # 6 1 p D[ -1+2x13]),U4
[40] 'FOR ELEMENT:',(EL[I;]),'. PRIOR TO RESLOPE'
[41] #+U+(3 0 # 6 1 p 10 40 70 xH[(E[I]+5)-4]),U
[42] #((I+I+1)≤9)/LOOP

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▽ COEF5
[1] # 9/09/86 PROGRAM TO FIT BILL'S 2 POINT RATIONAL EQUATION, CHECK OUT
[2] # VERIFICATION READINGS AND THE VALUES AFTER THE RESLOPE AS WELL.
[3] # MODEL A = B1+B2xC+B3xCxC
[4] 'MODEL A = B1+B2xC+B3xCxC'

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[5] R+ 10.10.250.10.40.250.30.20.2
[6] K+20+5*18*I+1
[7] T5+83#61p0
[8] 'FOR TUBE 401'
[9] LOOP:A++/32p(a,PPM401[1+16;E[I]+15])+2
[10] D+K+(a)(2050100xH[I]).*012
[11] * ORIGINAL COEFFICIENTS FOR THE TWO CURVES.
[12] D+a,PPM401[7+12;E[I]+15]
[13] +((-K[2]+2xK[3])<100xH[I])/OTHER
[14] O+((-K[2])+((K[2]*2)-4xK[3]x(K[1]-D[2x16]))*0.5)+2xK[3]
[15] U5+83#61p0
[16] T5+T5,[1]U5
[17] U+(60#61pD[-1+2x16]),83#61p0
[18] 'FOR ELEMENT: '(EL[I;]),' PRIOR TO RESLOPE'
[19] D+U+(30#61p104070xH[(E[I]+5)-4]),U
[20] C+/(a,PPM401[20+12;E[I]+15])+2
[21] C+C+(K+.x(50xH[I])*012)
[22] * GIVES MULTIPLICATIVE CORRECTION FACTORS FOR THE TWO LINES
[23] D+a,PPM401[(22+16),34+16;E[I]+15]
[24] II+1
[25] DOQ:DD+(K[2]*2)-4xK[3]x(K[1]-D[2xII])
[26] +((DD<0)/ERRORQ
[27] O[II]+Cx((-K[2])+DD*0.5)+2xK[3]
[28] LQ:+((II+II+1)<=6)/DOQ
[29] 'AFTER THE RESLOPE'
[30] U+(60#61pD[-1+2x16]),83#61p0
[31] D+a,PPM401[(28+16);E[I]+15]
[32] II+1
[33] KK+2446
[34] DO1:DD+(K[2]*2)-4xK[3]x(K[1]-D[KK[II]])
[35] +((DD<0)/ERROR1
[36] O[II]+Cx((-K[2])+DD*0.5)+2xK[3]
[37] L1:+((II+II+1)<=4)/DO1
[38] AO+Cx((-K[2])+((K[2]*2)-4xK[3]x(K[1]-D[2446]))*0.5)+2xK[3]
[39] U+(30+U),[1]((60#41pD[1335]),83#41p0),[1]30+U)
[40] D+U+(30#101p104070205050100104070xH[(E[I]+5)-4]),U
[41] +((I+I+1)<=9)/LOOP
[42] I+1
[43] LOOT:A++/32p(a,PPM401[41+16;E[I]+15])+2
[44] D+K+(a)(2050100xH[I]).*012
[45] D+a,PPM401[47+12;E[I]+15]
[46] +((-K[2]+2xK[3])<100xH[I])/OTHER1
[47] * ORIGINAL COEFFICIENTS FOR THE TWO CURVES.
[48] +((-K[2]*2)-4xK[3]x(K[1]-D[2x16]))<=0)/OTHER1
[49] O+((-K[2])+((K[2]*2)-4xK[3]x(K[1]-D[2x16]))*0.5)+2xK[3]
[50] U5+83#61p0
[51] T5+T5,[1]U5
[52] U+(60#61pD[-1+2x13]),83#61p0
[53] 'FOR ELEMENT: '(EL[I;]),' PRIOR TO RESLOPE'
[54] D+U+(30#61p104070xH[(E[I]+5)-4]),U
[55] +((I+I+1)<=9)/LOOT
[56] +0
[57] OTHER1:U5+83#61p0
[58] U+(60#61pD[-1+2x16]),83#61p0
[59] T5+T5,[1]U5
[60] 'FOR ELEMENT: '(EL[I;]),' PRIOR TO RESLOPE'
[61] D+U+(30#61p104070xH[(E[I]+5)-4]),U
[62] +((I+I+1)<=9)/LOOT
[63] ERROR1:O[II]+8888
[64] +L1
[65] ERRORQ:O[II]+8888
[66] +LQ
[67] AERROR1:O[II]+8888
[68] A+L1
[69] AERRORQ:O[II]+8888
[70] A+LQ
[71] OTHER:J+I
[72] U+(60#61pD[-1+2x16]),83#61p0
[73] U5+83#61p0

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[74] T5←T5,[1] U5
[75] 'FOR ELEMENT:' (EL[J:]), ' PRIOR TO RESLOPE'
[76] □←U+(3 0 * 6 1 p 10 40 70 *H[(E[J]+5)-4]),U
[77] D←PPM401[(22+16),34+16;E[J]+15]
[78] 'AFTER THE RESLOPE'
[79] U←(6 0 * 6 1 p D[1+2*16]), 8 3 * 6 1 p 0
[80] D←PPM401[(28+16);E[J]+15]
[81] U←(3 0 +U),[1](((6 0 * 4 1 p D[1 3 3 5]), 8 3 * 4 1 p 0),[1] 3 0 +U)
[82] □←U+(3 0 * 10 1 p 10 40 70 20 50 50 100 10 40 70 *H[(E[J]+5)-4]),U
[83] →((I+I+1)≤9)/LOOP
[84] I+1
[85] →LOOT
      ▽ SUMOFSQ
[1] AA←Q1,T1,T2,T3,T4,T5
[2] '-----'
[3] E←6*19*I+1
[4] LOOP:S←12 6 p, AA[E[I]+16:],[1] AA[(54+I*6)+16:]
[5] R←12 1 p, R1[E[I]+16:]
[6] □←R2←9 2 *R,[2] S
[7] NN←7 p 0
[8] SSUM←7 p 0
[9] V←7 p 0
[10] IN←1
[11] NEST:KO←1
[12] DO:SS+S[IN;KO]
[13] RR←R[IN:]
[14] →((SS>0)/COUNT
[15] →((KO+KO+1)≤6)/DO
[16] COUNT:NN[KO]+NN[KO]+1
[17] SSUM[KO]+SSUM[KO]+(SS-RR)*2
[18] →((KO+KO+1)≤6)/DO
[19] JK←1
[20] RATIO:→(NN[JK]=0)/GO
[21] V[JK]+SSUM[JK]+NN[JK]
[22] →((JK+JK+1)≤6)/RATIO
[23] →((IN+IN+1)≤12)/NEST
[24] GO:V[JK]+888
[25] →((JK+JK+1)≤6)/RATIO
[26] →((IN+IN+1)≤12)/NEST
[27] '-----'
[28] □←TT←EL[I:], ' ', 10 2 * 1 +SSUM
[29] □←UU←'SIZE', 10 2 * 1 +NN
[30] □←VV←'RATIO', 10 2 * 1 +V
[31] □←'RANK', 10 0 * AA1 + 1 +V
[32] '-----'
[33] →((I+I+1)≤9)/LOOP

```

APPENDIX D SUMMARY OF RANK

DATAAL PPM201

FE	175.00	190.49	68.89	119.33	972.52	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	29.17	31.75	10.93	19.89	162.09	888.00
RANK		3	1	2	4	5
AG	20.00	22.11	33.74	22.85	37.23	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	3.33	3.69	5.62	3.81	6.20	888.00
RANK	*	1	3	2	4	5
AL	36.00	49.29	47.87	47.76	47.82	48.75
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	6.00	8.22	7.98	7.96	7.97	8.12
RANK	*	5	3	1	2	4
CR	10.00	10.04	9.46	9.58	21.59	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	1.67	1.67	1.58	1.60	3.60	888.00
RANK		3	1	2	4	5
CU	28.00	29.44	33.00	31.86	42.77	33.90
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	4.67	4.91	5.50	5.31	7.13	5.65
RANK	*	1	3	2	5	4
MG	32.00	18.54	20.96	21.13	117.57	21.28
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	5.33	3.09	3.49	3.52	19.59	3.55
RANK		1	2	3	5	4
NI	19.00	25.62	370.96	14.67	89.70	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	3.17	4.27	61.83	2.44	14.95	888.00
RANK		2	4	1	3	5
SI	283.00	.00	383.29	193.42	.00	.00
SIZE	6.00	.00	6.00	6.00	.00	.00
RATIO	47.17	888.00	63.88	32.24	888.00	888.00
RANK		3	2	1	4	5
TI	22.00	21.31	167.05	18.08	21.07	19.45
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	3.67	3.55	27.84	3.01	3.51	3.24
RANK		4	5	1	3	2

DATAAL PPM202

FE	87.00	90.85	69.46	62.50	387.26	47.50
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	14.50	15.14	11.58	10.42	64.54	7.92
RANK		4	3	2	5	1
AG	5.00	4.67	3.31	4.80	35.16	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	.83	.78	.55	.80	5.86	888.00
RANK		2	1	3	4	5
AL	11.00	14.13	16.47	16.22	20.50	18.22

SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	1.83	2.36	2.75	2.70	3.42	3.04
RANK	*	1	3	2	5	4
CR	2.00	1.33	2.03	1.78	4.37	2.28
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	.33	.22	.34	.30	.73	.38
RANK		1	3	2	5	4
CU	42.00	38.05	32.38	34.15	59.89	25.46
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	7.00	6.34	5.40	5.69	9.98	4.24
RANK		4	2	3	5	1
MG	52.00	46.76	47.40	47.26	89.28	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	8.67	7.79	7.90	7.88	14.88	888.00
RANK		1	3	2	4	5
NI	7.00	4.04	6486.57	9.90	40.81	15.54
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	1.17	.67	1081.10	1.65	6.80	2.59
RANK		1	5	2	4	3
SI	304.00	302.19	293.80	289.35	296.75	298.05
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	50.67	50.37	48.97	48.23	49.46	49.67
RANK		5	2	1	3	4
TI	8.00	8.79	10.58	9.53	9.41	9.47
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	1.33	1.46	1.76	1.59	1.57	1.58
RANK	*	1	5	4	2	3

DATAAL PPM203

FE	230.00	240.99	96.31	179.67	396.15	123.58
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	38.33	40.17	16.05	29.95	66.02	20.60
RANK		4	1	3	5	2
AG	2.00	1.59	.53	1.14	17.87	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	.33	.26	.09	.19	2.98	888.00
RANK		3	1	2	4	5
AL	52.00	45.26	44.56	46.30	58.88	41.04
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	8.67	7.54	7.43	7.72	9.81	6.84
RANK		3	2	4	5	1
CR	1.00	.53	.48	.25	4.27	.29
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	.17	.09	.08	.04	.71	.05
RANK		4	3	1	5	2
CU	245.00	344.41	386.51	367.05	354.90	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	40.83	57.40	64.42	61.18	59.15	888.00
RANK	*	1	4	3	2	5
MG	9.00	5.04	7.18	7.31	29.65	12.93
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	1.50	.84	1.20	1.22	4.94	2.15
RANK		1	2	3	5	4
NI	9.00	9.17	120.78	17.14	22.59	19.61
SIZE	6.00	6.00	6.00	6.00	6.00	6.00

RATIO	1.50	1.53	20.13	2.86	3.77	3.27
RANK	*	1	5	2	4	3
SI	305.00	757.69	552.89	1687.68	753.81	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	50.83	126.28	92.15	281.28	125.63	888.00
RANK	*	3	1	4	2	5
TI	13.00	13.54	14.30	14.49	14.32	14.36
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	2.17	2.26	2.38	2.42	2.39	2.39
RANK	*	1	2	5	3	4

DATAAL PPM206

FE	725.00	823.52	739.28	807.14	979.47	798.93
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	60.42	68.63	61.61	67.26	81.62	66.58
RANK	*	4	1	3	5	2
AG	3.00	19.31	17.25	17.95	47.74	15.83
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	.25	1.61	1.44	1.50	3.98	2.64
RANK	*	3	1	2	5	4
AL	317.00	111.67	265.15	104.02	6031.69	.00
SIZE	12.00	6.00	6.00	6.00	6.00	.00
RATIO	26.42	18.61	44.19	17.34	1005.28	888.00
RANK	*	2	3	1	5	4
CR	7.00	19.88	20.78	20.50	33.57	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	.58	1.66	1.73	1.71	2.80	888.00
RANK	*	1	3	2	4	5
CU	114.00	170.55	169.35	169.37	183.95	160.21
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	9.50	14.21	14.11	14.11	15.33	13.35
RANK	*	4	2	3	5	1
MG	77.00	163.79	172.05	163.16	179.81	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	6.42	13.65	14.34	13.60	14.98	888.00
RANK	*	2	3	1	4	5
NI	3.00	93.12	333.76	93.53	162.60	73.45
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	.25	7.76	27.81	7.79	13.55	12.24
RANK	*	1	5	2	4	3
SI	748.00	6691.70	1407.42	345.57	1493.15	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	62.33	557.64	117.28	28.80	124.43	888.00
RANK	*	4	2	1	3	5
TI	27.00	87.04	78.46	102.10	89.66	52.04
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	2.25	7.25	6.54	8.51	7.47	8.67
RANK	*	2	1	4	3	5

DATAAL PPM207

FE	2898.00	2914.98	19613.44	3395.53	1457.12	.00
SIZE	12.00	12.00	12.00	12.00	6.00	.00
RATIO	241.50	242.92	1634.45	282.96	242.85	888.00
RANK	*	2	5	3	1	4

AG	14.00	154.71	.26	52.41	1.33	.27
SIZE	12.00	12.00	6.00	12.00	6.00	6.00
RATIO	1.17	12.89	.04	4.37	.22	.04
RANK		5	1	4	3	2

AL	320.00	.00	234.39	.00	.00	.00
SIZE	11.00	.00	12.00	.00	.00	.00
RATIO	29.09	888.00	19.53	888.00	888.00	888.00
RANK		2	1	3	4	5

CR	11.00	60.80	32.88	52.55	28.17	.00
SIZE	12.00	12.00	12.00	12.00	6.00	.00
RATIO	.92	5.07	2.74	4.38	4.69	888.00
RANK	*	4	1	2	3	5

CU	29.00	290.09	135.19	943.80	20.49	23.40
SIZE	12.00	12.00	12.00	12.00	6.00	6.00
RATIO	2.42	24.17	11.27	78.65	3.41	3.90
RANK	*	4	3	5	1	2

MG	89.00	395.16	97.60	794.64	107.73	.00
SIZE	12.00	12.00	6.00	12.00	6.00	.00
RATIO	7.42	32.93	16.27	66.22	17.95	888.00
RANK	*	3	1	4	2	5

NI	15.00	226.95	271.83	221.93	201.59	4.00
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	1.25	18.91	22.65	18.49	16.80	.67
RANK		4	5	3	2	1

SI	443.00	382.24	703.85	528.94	405.50	.00
SIZE	12.00	6.00	12.00	6.00	6.00	.00
RATIO	36.92	63.71	58.65	88.16	67.58	888.00
RANK	*	2	1	4	3	5

TI	67.00	230.65	67.83	358.10	171.87	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	5.58	19.22	5.65	29.84	14.32	888.00
RANK	*	3	1	4	2	5

DATAAL PPM302

FE	179.00	295.18	284.07	279.17	1040.62	327.22
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	14.92	24.60	23.67	23.26	86.72	27.27
RANK	*	3	2	1	5	4

AG	12.00	13.72	11.42	12.52	27.56	.24
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	1.00	1.14	.95	1.04	2.30	.04
RANK		4	2	3	5	1

AL	191.00	27.67	184.49	37.31	.00	.00
SIZE	10.00	6.00	6.00	6.00	.00	.00
RATIO	19.10	4.61	30.75	6.22	888.00	888.00
RANK		1	3	2	4	5

CR	6.00	6.31	6.63	6.36	13.35	5.10
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	.50	.53	.55	.53	1.11	.43
RANK		2	4	3	5	1

CU	32.00	56.40	64.10	63.24	85.96	65.02
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	2.67	4.70	5.34	5.27	7.16	5.42
RANK	*	1	3	2	5	4

MG	85.00	118.39	118.44	114.24	191.42	.00

SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	7.08	9.87	9.87	9.52	15.95	888.00
RANK	*	2	3	1	4	5
NI	20.00	21.67	97.51	14.37	100.50	10.37
SIZE	12.00	12.00	6.00	12.00	12.00	6.00
RATIO	1.67	1.81	16.25	1.20	8.37	1.73
RANK		3	5	1	4	2
SI	65.00	97.88	92.17	93.20	92.57	87.42
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	5.42	8.16	7.68	7.77	7.71	7.29
RANK	*	5	2	4	3	1
TI	39.00	48.27	49.77	51.86	49.15	48.74
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	3.25	4.02	4.15	4.32	4.10	4.06
RANK	*	1	4	5	3	2

DATAAL PPM303

FE	108.00	107.97	111.92	112.21	347.53	123.26
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	18.00	18.00	18.65	18.70	57.92	20.54
RANK		1	2	3	5	4
AG	.00	.21	.36	.32	7.19	.84
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	.00	.03	.06	.05	1.20	.14
RANK	*	1	3	2	5	4
AL	6.00	11.92	14.39	11.16	23.00	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	1.00	1.99	2.40	1.86	3.83	888.00
RANK	*	2	3	1	4	5
CR	1.00	2.29	3.35	2.85	12.25	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	.17	.38	.56	.47	2.04	888.00
RANK	*	1	3	2	4	5
CU	40.00	39.34	44.48	53.20	43.62	42.99
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	6.67	6.56	7.41	8.87	7.27	7.17
RANK		1	4	5	3	2
MG	32.00	24.80	41.08	45.75	34.14	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	5.33	4.13	6.85	7.63	5.69	888.00
RANK		1	3	4	2	5
NI	6.00	6.50	.00	15.41	64.70	.00
SIZE	6.00	6.00	.00	6.00	6.00	.00
RATIO	1.00	1.08	888.00	2.57	10.78	888.00
RANK	*	1	4	2	3	5
SI	53.00	64.80	505.64	66.91	60.38	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	8.83	10.80	84.27	11.15	10.06	888.00
RANK	*	2	4	3	1	5
TI	11.00	11.60	6.72	12.54	12.18	12.21
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	1.83	1.93	1.12	2.09	2.03	2.04
RANK		2	1	5	3	4

DATAAL1 PPM304

FE	64.00	2180.05	3250.60	7185.16	50624.67	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	10.67	363.34	541.77	1197.53	8437.45	888.00
RANK	*	1	2	4	5	3
AG	10.00	10.03	9.97	9.97	13.44	8.62
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	1.67	1.67	1.66	1.66	2.24	1.44
RANK		4	2	3	5	1
AL	72.00	88.29	88.09	87.99	103.49	90.89
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	12.00	14.71	14.68	14.66	17.25	15.15
RANK	*	3	2	1	5	4
CR	4.00	3.97	3.07	3.32	13.88	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	.67	.66	.51	.55	2.31	888.00
RANK		3	1	2	4	5
CU	23.00	25.91	24.46	24.46	24.99	26.71
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	3.83	4.32	4.08	4.08	4.16	4.45
RANK	*	4	1	2	3	5
MG	78.00	82.36	.00	92.85	104.65	.00
SIZE	6.00	6.00	.00	6.00	6.00	.00
RATIO	13.00	13.73	888.00	15.48	17.44	888.00
RANK	*	1	4	2	3	5
NI	3.00	2.32	22.76	5.26	197.67	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	.50	.39	3.79	.88	32.95	888.00
RANK		1	3	2	4	5
SI	17.00	16.86	22.34	12.79	16.26	16.24
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	2.83	2.81	3.72	2.13	2.71	2.71
RANK		4	5	1	3	2
TI	24.00	26.31	27.28	27.70	27.08	26.45
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	4.00	4.39	4.55	4.62	4.51	4.41
RANK	*	1	4	5	3	2

DATAAL PPM305

FE	402.00	460.36	421.86	427.68	3416.65	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	67.00	76.73	70.31	71.28	569.44	888.00
RANK	*	3	1	2	4	5
AG	2.00	3.41	3.68	3.63	6.13	4.19
SIZE	5.00	6.00	6.00	6.00	6.00	6.00
RATIO	.40	.57	.61	.60	1.02	.70
RANK	*	1	3	2	5	4
AL	35.00	44.39	21.72	71.60	.00	.00
SIZE	6.00	6.00	6.00	6.00	.00	.00
RATIO	5.83	7.40	3.62	11.93	888.00	888.00
RANK		2	1	3	4	5
CR	2.00	1.87	2.05	1.91	3.59	2.19
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	.33	.31	.34	.32	.60	.36
RANK		1	3	2	5	4

CU	165.00	179.86	176.91	179.08	212.23	176.33
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	27.50	29.98	29.49	29.85	35.37	29.39
RANK	*	4	2	3	5	1

MG	118.00	118.06	119.86	119.84	467.95	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	19.67	19.68	19.98	19.97	77.99	888.00
RANK	*	1	3	2	4	5

NI	35.00	32.50	27.84	21.81	81.19	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	5.83	5.42	4.64	3.64	13.53	888.00
RANK		3	2	1	4	5

SI	19.00	14.18	17.28	16.45	19.92	16.95
SIZE	6.00	6.00	6.00	6.00	6.00	6.00
RATIO	3.17	2.36	2.88	2.74	3.32	2.83
RANK		1	4	2	5	3

TI	38.00	51.08	1687.75	61.65	51.55	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	6.33	8.51	281.29	10.28	8.59	888.00
RANK	*	1	4	3	2	5

DATAAL PPM401

FE	100.00	128.03	121.08	109.09	3119.08	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	8.33	10.67	10.09	9.09	259.92	888.00
RANK	*	3	2	1	4	5

AG	2.00	.85	1.70	1.48	12.34	.67
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	.17	.07	.14	.12	1.03	.11
RANK		1	4	3	5	2

AL	268.00	317.88	300.21	297.57	429.29	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	22.33	26.49	25.02	24.80	35.77	888.00
RANK	*	3	2	1	4	5

CR	16.00	17.29	17.16	17.10	39.32	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	1.33	1.44	1.43	1.42	3.28	888.00
RANK	*	3	2	1	4	5

CU	106.00	102.05	100.06	99.74	100.09	107.97
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	8.83	8.50	8.34	8.31	8.34	9.00
RANK		4	2	1	3	5

MG	91.00	170.50	156.21	148.12	179.10	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	7.58	14.21	13.02	12.34	14.92	888.00
RANK	*	3	2	1	4	5

NI	38.00	34.87	26.02	31.70	192.41	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	3.17	2.91	2.17	2.64	16.03	888.00
RANK		3	1	2	4	5

SI	162.00	149.09	141.29	166.29	171.89	103.04
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	13.50	12.42	11.77	13.86	14.32	17.17
RANK		2	1	3	4	5

TI	135.00	142.77	364.12	141.42	134.41	141.47

SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	11.25	11.90	30.34	11.79	11.20	11.79
RANK		4	5	2	1	3

DATAAL PPM404

FE	30.00	38.60	92.84	73.23	1614.32	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	5.00	6.43	15.47	12.21	269.05	888.00
RANK	*	1	3	2	4	5

AG	9.00	7.92	6.83	8.49	98.37	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	1.50	1.32	1.14	1.42	16.39	888.00
RANK		2	1	3	4	5

AL	231.00	427.14	240.35	327.70	966.05	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	38.50	71.19	40.06	54.62	161.01	888.00
RANK	*	3	1	2	4	5

CR	1.00	.97	2.22	1.10	10.77	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	.17	.16	.37	.18	1.80	888.00
RANK		1	3	2	4	5

CU	113.00	116.47	125.29	127.72	238.24	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	18.83	19.41	20.88	21.29	39.71	888.00
RANK	*	1	2	3	4	5

MG	47.00	44.59	45.87	46.35	46.32	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	7.83	7.43	7.64	7.73	7.72	888.00
RANK		1	2	4	3	5

NI	11.00	8.71	5.84	6.41	145.33	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	1.83	1.45	.97	1.07	24.22	888.00
RANK		3	1	2	4	5

SI	36.00	53.14	42.51	69.86	820.92	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	6.00	8.86	7.08	11.64	136.82	888.00
RANK	*	2	1	3	4	5

TI	32.00	26.54	25.41	26.86	25.17	.00
SIZE	6.00	6.00	6.00	6.00	6.00	.00
RATIO	5.33	4.42	4.23	4.48	4.20	888.00
RANK		3	2	4	1	5

DATAAL PPM405

FE	239.00	223.84	175.23	193.22	4758.95	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	19.92	18.65	14.60	16.10	396.58	888.00
RANK		3	1	2	4	5

AG	24.00	19.51	18.68	19.07	23.69	14.31
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	2.00	1.63	1.56	1.59	1.97	1.19
RANK		4	2	3	5	1

AL	120.00	136.21	133.36	135.06	138.79	133.60
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	10.00	11.35	11.11	11.25	11.57	11.13
RANK	*	4	1	3	5	2

CR	3.00	6.97	4.91	6.10	14.73	3.51
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	.25	.58	.41	.51	1.23	.59
RANK	*	3	1	2	5	4
CU	47.00	41.02	47.51	46.43	48.57	49.51
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	3.92	3.42	3.96	3.87	4.05	4.13
RANK		1	3	2	4	5
MG	266.00	254.16	220.37	194.30	213.89	74.36
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	22.17	21.18	18.36	16.19	17.82	12.39
RANK		5	4	2	3	1
NI	26.00	32.24	8.76	20.00	167.12	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	2.17	2.69	.73	1.67	13.93	888.00
RANK		3	1	2	4	5
SI	65.00	71.11	71.03	71.21	85.72	60.57
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	5.42	5.93	5.92	5.93	7.14	5.05
RANK		3	2	4	5	1
TI	40.00	44.50	45.33	43.82	44.29	25.18
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	3.33	3.71	3.78	3.65	3.69	4.20
RANK	*	3	4	1	2	5

DATAAL PPM408

FE	469.00	547.44	363.16	415.18	17220.25	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	39.08	45.62	30.26	34.60	1435.02	888.00
RANK		3	1	2	5	4
AG	33.00	28.08	26.34	26.91	37.04	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	2.75	2.34	2.19	2.24	3.09	888.00
RANK		3	1	2	4	5
AL	74.00	63.13	58.35	61.43	149.71	10.28
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	6.17	5.26	4.86	5.12	12.48	1.71
RANK		4	2	3	5	1
CR	2.00	3.48	3.37	3.36	96.86	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	.17	.29	.28	.28	8.07	888.00
RANK	*	3	2	1	4	5
CU	32.00	35.41	41.01	42.99	43.21	45.57
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	2.67	2.95	3.42	3.58	3.60	3.80
RANK	*	1	2	3	4	5
MG	91.00	83.41	95.59	94.42	280.57	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	7.58	6.95	7.97	7.87	23.38	888.00
RANK		1	3	2	4	5
NI	30.00	28.84	8.25	16.98	478.03	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	2.50	2.40	.69	1.41	39.84	888.00
RANK		3	1	2	4	5

SI	132.00	329.10	306.22	308.16	316.09	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	11.00	27.42	25.52	25.43	26.34	888.00
RANK	*	4	2	1	3	5

TI	35.00	32.77	27.75	29.39	27.98	28.11
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	2.92	2.73	2.31	2.45	2.33	2.34
RANK		5	1	4	2	3

DATAAL PPM409

FE	915.00	896.53	673.04	788.52	2643.86	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	76.25	74.71	56.09	65.71	220.32	888.00
RANK		3	1	2	4	5

AG	41.00	147.67	5523.28	292.54	90.84	7.25
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	3.42	12.31	460.27	24.38	7.57	1.21
RANK		3	5	4	2	1

AL	277.00	283.60	265.18	264.74	255.95	265.68
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	23.08	23.63	22.10	22.06	21.33	22.14
RANK		5	3	2	1	4

CR	3.00	4.47	5.43	5.10	26.16	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	.25	.37	.45	.42	2.18	888.00
RANK	*	1	3	2	4	5

CU	134.00	146.62	157.18	154.99	218.25	153.89
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	11.17	12.22	13.10	12.92	18.19	12.82
RANK	*	1	4	3	5	2

MG	73.00	77.80	78.61	76.53	95.99	66.93
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	6.08	6.48	6.55	6.38	8.00	11.15
RANK	*	2	3	1	4	5

NI	3.00	2.85	73.03	6.58	97.48	7.02
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	.25	.24	6.09	.55	8.12	1.17
RANK		1	4	2	5	3

SI	525.00	.00	216.03	.00	5087.06	.00
SIZE	12.00	.00	12.00	.00	12.00	.00
RATIO	43.75	888.00	18.00	888.00	423.92	888.00
RANK		3	1	4	2	5

TI	29.00	25.84	26.30	23.93	26.04	21.23
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	2.42	2.15	2.19	1.99	2.17	3.54
RANK		2	4	1	3	5

DATAAL PPM412

FE	41.00	50.60	142.39	76.18	1098.40	85.99
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	3.42	4.22	11.87	6.35	91.53	14.33
RANK	*	1	3	2	5	4

AG	6.00	5.50	3.93	4.56	11.75	1.68
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	.50	.46	.33	.38	.98	.14

RANK		4	2	3	5	1
AL	85.00	241.77	357.39	280.80	53.34	.00
SIZE	12.00	12.00	12.00	12.00	6.00	.00
RATIO	7.08	20.15	29.78	23.40	8.89	888.00
RANK	*	2	4	3	1	5
CR	3.00	2.53	1.94	2.47	15.28	.25
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	.25	.21	.16	.21	1.27	.04
RANK		4	2	3	5	1
CU	33.00	35.82	47.30	44.36	51.29	47.38
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	2.75	2.99	3.94	3.70	4.27	3.95
RANK	*	1	3	2	5	4
MG	53.00	42.21	43.66	44.11	163.60	18.44
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	4.42	3.52	3.64	3.68	13.63	3.07
RANK		2	3	4	5	1
NI	23.00	16.25	809.13	19.34	96.26	22.57
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	1.92	1.35	67.43	1.61	8.02	3.76
RANK		1	5	2	4	3
SI	880.00	1541.92	.00	.00	3279.27	.00
SIZE	12.00	12.00	.00	.00	12.00	.00
RATIO	73.33	128.49	888.00	888.00	273.27	888.00
RANK	*	1	3	4	2	5
TI	21.00	21.31	20.92	20.66	21.39	20.88
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	1.75	1.78	1.74	1.72	1.78	1.74
RANK		4	3	1	5	2

DATAAL PPM413

FE	320.00	298.68	2303.91	231.51	986.67	147.91
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	26.67	24.89	191.99	19.29	82.22	12.33
RANK		3	5	2	4	1
AG	17.00	17.38	15.35	16.32	21.01	14.15
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	1.42	1.45	1.28	1.36	1.75	1.18
RANK		4	2	3	5	1
AL	124.00	267.30	138.11	14.47	16.94	16.80
SIZE	12.00	12.00	12.00	6.00	6.00	6.00
RATIO	10.33	22.27	11.51	2.41	2.82	2.80
RANK		5	4	1	3	2
CR	7.00	4.48	20.26	4.47	15.25	1.18
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	.58	.37	1.69	.37	1.27	.20
RANK		3	5	2	4	1
CU	81.00	65.38	68.28	61.92	99.51	57.06
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	6.75	5.45	5.69	5.16	8.29	4.75
RANK		3	4	2	5	1
MG	72.00	66.45	66.15	65.62	110.91	55.46
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	6.00	5.54	5.51	5.47	9.24	9.24
RANK		3	2	1	4	5

NI	16.00	16.21	306.24	16.84	87.17	14.39
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	1.33	1.35	25.52	1.40	7.26	2.40
RANK	*	1	5	2	4	3
SI	548.00	3253.60	363.43	37162.64	1317.74	.00
SIZE	12.00	12.00	6.00	12.00	12.00	.00
RATIO	45.67	271.13	60.57	3096.89	109.81	888.00
RANK	*	3	1	5	2	4
TI	25.00	30.57	26.31	31.13	29.47	29.45
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	2.08	2.55	2.19	2.59	2.46	2.45
RANK	*	4	1	5	3	2

DATAAL PPM415

FE	95.00	98.18	97.72	76.81	768.92	110.22
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	7.92	8.18	8.14	6.40	64.08	9.18
RANK		3	2	1	5	4
AG	8.00	7.57	7.88	7.93	8.29	7.53
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	.67	.63	.66	.66	.69	.63
RANK		2	3	4	5	1
AL	67.00	88.27	109.28	88.64	113.79	85.55
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	5.58	7.36	9.11	7.39	9.48	7.13
RANK	*	2	4	3	5	1
CR	2.00	1.43	.93	.86	5.70	.51
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	.17	.12	.08	.07	.48	.04
RANK		4	3	2	5	1
CU	23.00	25.87	27.27	28.05	65.71	32.73
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	1.92	2.16	2.27	2.34	5.48	2.73
RANK	*	1	2	3	5	4
MG	31.00	24.13	29.92	30.08	124.74	12.40
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	2.58	2.01	2.49	2.51	10.39	2.07
RANK		1	3	4	5	2
NI	13.00	12.73	423.45	16.17	39.13	19.71
SIZE	12.00	12.00	6.00	12.00	12.00	12.00
RATIO	1.08	1.06	70.58	1.35	3.26	1.64
RANK		1	5	2	4	3
SI	288.00	197.67	344.15	222.31	174.00	.00
SIZE	11.00	6.00	12.00	6.00	6.00	.00
RATIO	26.18	32.94	28.68	37.05	29.00	888.00
RANK	*	3	1	4	2	5
TI	20.00	14.81	9.61	14.86	14.62	.00
SIZE	12.00	12.00	6.00	12.00	12.00	.00
RATIO	1.67	1.23	1.60	1.24	1.22	888.00
RANK		2	4	3	1	5

DATAAL PPM416

FE	372.00	1112.62	906.69	971.94	1775.91	93.76
SIZE	12.00	12.00	12.00	12.00	12.00	6.00

RATIO	31.00	92.72	75.56	80.99	147.99	15.63
RANK		4	2	3	5	1
AG	11.00	23.19	22.85	22.93	18.28	6.51
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	.92	1.93	1.90	1.91	1.52	1.09
RANK	*	5	3	4	2	1
AL	81.00	114.20	106.24	108.97	117.53	110.48
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	6.75	9.52	8.85	9.08	9.79	9.21
RANK	*	4	1	2	5	3
CR	5.00	19.94	17.17	18.37	18.20	.89
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	.42	1.66	1.43	1.53	1.52	.15
RANK		5	2	4	3	1
CU	26.00	141.62	145.05	140.81	127.41	138.40
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	2.17	11.80	12.09	11.73	10.62	11.53
RANK	*	4	5	3	1	2
MG	29.00	106.64	115.62	112.50	121.80	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	2.42	8.89	9.64	9.37	10.15	888.00
RANK	*	1	3	2	4	5
NI	15.00	96.20	153.68	90.07	118.51	83.23
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	1.25	8.02	12.81	7.51	9.88	13.87
RANK	*	2	4	1	3	5
SI	148.00	928.14	1098.13	1078.96	425.30	40.81
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	12.33	77.35	91.51	89.91	35.44	6.80
RANK		3	5	4	2	1
TI	35.00	53.25	14.04	59.28	57.97	.00
SIZE	12.00	12.00	6.00	12.00	12.00	.00
RATIO	2.92	4.44	2.34	4.94	4.83	888.00
RANK		2	1	4	3	5

DATAAL PPM501

FE	180.00	78.30	179.09	131.65	7817.24	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	15.00	6.52	14.92	10.97	651.44	888.00
RANK		1	3	2	4	5
AG	10.00	9.19	6.30	7.81	31.30	1.95
SIZE	11.00	12.00	12.00	12.00	12.00	6.00
RATIO	.91	.77	.53	.65	2.61	.32
RANK		4	2	3	5	1
AL	104.00	139.48	60.48	45.26	46.08	43.34
SIZE	10.00	12.00	12.00	6.00	6.00	6.00
RATIO	10.40	11.62	5.04	7.54	7.68	7.22
RANK		5	1	3	4	2
CR	3.00	5.68	23.18	6.15	31.26	.00
SIZE	11.00	12.00	12.00	12.00	12.00	.00
RATIO	.27	.47	1.93	.51	2.61	888.00
RANK	*	1	3	2	4	5
CU	16.00	11.63	9.58	10.06	19.84	9.96
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	1.33	.97	.80	.84	1.65	.83

RANK	4	1	3	5	2	
MG	183.00	187.93	190.79	190.92	1091.13	.00
SIZE	11.00	12.00	12.00	12.00	12.00	.00
RATIO	16.64	15.66	15.90	15.91	90.93	888.00
RANK		1	2	3	4	5
NI	21.00	19.30	141.79	24.83	185.73	.00
SIZE	11.00	12.00	12.00	12.00	12.00	.00
RATIO	1.91	1.61	11.82	2.07	15.48	888.00
RANK		1	3	2	4	5
SI	408.00	898.40	499.81	232.03	537.23	564.70
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	34.00	74.87	41.65	19.34	44.77	47.06
RANK		5	2	1	3	4
TI	23.00	19.37	18.96	18.74	18.88	8.07
SIZE	11.00	12.00	12.00	12.00	12.00	6.00
RATIO	2.09	1.61	1.58	1.56	1.57	1.34
RANK		5	4	2	3	1

DATAAL PPM502

FE	310.00	299.60	267.45	281.80	402.95	260.74
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	25.83	24.97	22.29	23.48	33.58	21.73
RANK		4	2	3	5	1
AG	4.00	3.14	3.51	4.14	10.61	4.13
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	.33	.26	.29	.34	.88	.34
RANK		1	2	4	5	3
AL	438.00	488.21	453.86	484.93	4541.83	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	36.50	40.68	37.82	37.91	378.49	888.00
RANK	*	3	1	2	4	5
CR	6.00	7.42	6.21	7.13	21.56	1.12
SIZE	11.00	12.00	12.00	12.00	12.00	6.00
RATIO	.55	.62	.52	.59	1.80	.19
RANK		4	2	3	5	1
CU	36.00	40.26	41.72	44.04	48.01	51.05
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	3.00	3.35	3.48	3.67	4.00	4.25
RANK	*	1	2	3	4	5
MG	182.00	169.20	191.42	189.20	2219.15	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	15.17	14.10	15.95	15.77	184.93	888.00
RANK		1	3	2	4	5
NI	25.00	11.97	185.09	7.33	37.72	1.79
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	2.08	1.00	15.42	.61	3.14	.30
RANK		3	5	2	4	1
SI	230.00	136.61	53.51	143.39	58.97	.00
SIZE	11.00	12.00	6.00	12.00	6.00	.00
RATIO	20.91	11.38	8.92	11.95	9.83	888.00
RANK		3	1	4	2	5
TI	30.00	34.99	36.46	64.72	42.50	15.75
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	2.50	2.92	3.04	5.39	3.54	2.62
RANK	*	2	3	5	4	1

DATAAL PPM503

FE	1648.00	2124.98	2092.60	2073.75	2461.83	1363.09
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	139.00	177.08	174.38	172.81	205.15	227.18
RANK	*	3	2	1	4	5
AG	27.00	50.22	43.14	43.26	13.81	.00
SIZE	11.00	12.00	12.00	12.00	6.00	.00
RATIO	2.45	4.18	3.60	3.60	2.30	888.00
RANK		4	2	3	1	5
AL	141.00	215.83	209.43	67.76	67.44	.00
SIZE	12.00	12.00	12.00	6.00	6.00	.00
RATIO	11.75	17.99	17.45	11.29	11.24	888.00
RANK		4	3	2	1	5
CR	9.00	23.01	23.92	22.41	25.39	7.38
SIZE	11.00	12.00	12.00	12.00	12.00	6.00
RATIO	.82	1.92	1.99	1.87	2.12	1.23
RANK	*	3	4	2	5	1
CU	447.00	652.44	650.39	637.09	611.21	281.30
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	37.25	54.37	54.20	53.09	50.93	46.88
RANK	*	5	4	3	2	1
MG	79.00	183.36	182.30	181.26	13744.22	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	6.58	15.28	15.19	15.10	1145.35	888.00
RANK	*	3	2	1	5	4
NI	149.00	252.63	5861.19	250.95	220.62	156.12
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	12.42	21.05	488.43	20.91	18.39	26.02
RANK	*	3	5	2	1	4
SI	1411.00	768.20	4163.73	880.75	3977.22	.00
SIZE	12.00	6.00	12.00	12.00	6.00	.00
RATIO	117.58	128.03	346.98	73.40	662.87	888.00
RANK		2	3	1	4	5
TI	94.00	192.46	117.78	124.93	36.04	.00
SIZE	12.00	12.00	12.00	12.00	6.00	.00
RATIO	7.83	16.04	9.82	10.41	6.01	888.00
RANK		4	2	3	1	5

DATAAL PPM504

FE	1047.00	1171.91	1653.12	1159.36	1525.00	1168.39
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	87.25	97.66	137.76	96.61	127.08	97.37
RANK	*	3	5	1	4	2
AG	12.00	14.46	12.87	13.81	15.14	13.80
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	1.00	1.21	1.07	1.15	1.26	1.15
RANK	*	4	1	3	5	2
AL	317.00	760.40	361.65	187.29	676.56	.00
SIZE	12.00	12.00	12.00	6.00	6.00	.00
RATIO	26.42	63.37	30.14	31.21	112.76	888.00
RANK	*	3	1	2	4	5
CR	11.00	9.05	13.12	9.30	11.83	9.38
SIZE	12.00	12.00	12.00	12.00	12.00	12.00

RATIO	.92	.75	1.09	.77	.99	.78
RANK		1	5	2	4	3
CU	222.00	240.31	193.48	226.03	284.58	214.89
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	18.50	20.03	16.12	18.84	23.71	17.91
RANK		4	1	3	5	2
MG	132.00	160.42	157.35	157.49	1412.08	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	11.00	13.37	13.11	13.12	117.67	888.00
RANK	*	3	1	2	4	5
NI	74.00	71.55	77.34	71.63	83.46	72.91
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	6.17	5.96	6.44	5.97	6.95	6.08
RANK		1	4	2	5	3
SI	108.00	203.42	263.66	238.05	222.00	80.49
SIZE	11.00	11.00	11.00	11.00	12.00	6.00
RATIO	9.82	18.49	23.97	21.64	18.50	13.42
RANK	*	2	5	4	3	1
TI	21.00	25.83	93.57	33.93	3.09	.00
SIZE	12.00	12.00	12.00	12.00	6.00	.00
RATIO	1.75	2.15	7.80	2.83	.52	888.00
RANK		2	4	3	1	5

DATAAL PPM505

FE	440.00	424.89	63786.81	578.67	669.23	628.68
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	36.67	35.41	5315.57	48.22	55.77	52.39
RANK		1	5	2	4	3
AG	2.00	1.50	3.23	1.87	1.65	1.84
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	.17	.13	.27	.16	.14	.15
RANK		1	5	4	2	3
AL	1239.00	454.57	360.05	.00	36003.11	.00
SIZE	12.00	6.00	12.00	.00	6.00	.00
RATIO	103.25	75.76	30.00	888.00	6000.52	888.00
RANK		2	1	3	5	4
CR	2.00	1.86	2.26	2.99	3.07	3.04
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	.17	.16	.19	.25	.26	.25
RANK		1	2	3	5	4
CU	71.00	65.44	37.30	78.48	81.53	81.01
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	5.92	5.45	3.11	6.54	6.79	6.75
RANK		2	1	3	5	4
MG	144.00	154.07	153.27	153.39	481.96	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	12.00	12.84	12.77	12.78	40.16	888.00
RANK	*	3	1	2	4	5
NI	10.00	12.28	47.80	21.40	21.93	21.83
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	.83	1.02	3.98	1.78	1.83	1.82
RANK	*	1	5	2	4	3
SI	71.00	69.86	210.06	71.19	70.94	28.93
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	5.92	5.82	17.50	5.93	5.91	4.82

RANK	2	5	4	3	1	
TI	21.00	25.80	23.69	39.51	29.40	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	1.75	2.15	1.97	3.29	2.45	888.00
RANK	*	2	1	4	3	5

DATAAL PPM507

FE	598.00	560.43	2731.83	738.05	976.95	793.70
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	49.83	46.70	227.65	61.50	81.41	66.14
RANK		1	5	2	4	3

AG	9.00	6.41	7.98	7.16	11.86	7.30
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	.75	.53	.66	.60	.99	.61
RANK		1	4	2	5	3

AL	324.00	557.96	509.22	.00	.00	.00
SIZE	6.00	6.00	12.00	.00	.00	.00
RATIO	54.00	92.99	42.44	888.00	888.00	888.00
RANK		2	1	3	4	5

CR	4.00	4.57	2.92	6.39	11.64	1.38
SIZE	12.00	12.00	6.00	12.00	12.00	6.00
RATIO	.33	.38	.49	.53	.97	.23
RANK		2	3	4	5	1

CU	61.00	57.13	69.86	65.67	80.20	69.83
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	5.08	4.76	5.82	5.47	6.68	5.82
RANK		1	4	2	5	3

MG	180.00	231.80	206.02	243.13	2856.46	.00
SIZE	12.00	12.00	12.00	12.00	12.00	.00
RATIO	15.00	19.32	17.17	20.26	238.04	888.00
RANK	*	2	1	3	4	5

NI	35.00	33.17	39.69	44.62	56.64	48.32
SIZE	12.00	12.00	12.00	12.00	12.00	12.00
RATIO	2.92	2.76	3.31	3.72	4.72	4.03
RANK		1	2	3	5	4

SI	194.00	258.31	75630.78	432.132415930.14		.00
SIZE	6.00	6.00	11.00	6.00	12.00	.00
RATIO	32.33	43.05	6875.53	72.02	201327.51	888.00
RANK	*	1	4	2	5	3

TI	45.00	43.50	66.30	44.06	45.37	20.91
SIZE	12.00	12.00	12.00	12.00	12.00	6.00
RATIO	3.75	3.62	5.52	3.67	3.78	3.49
RANK		2	5	3	4	1

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7-87

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