

THE OPTIMUM QUANTIZATION PROCESS

by

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requirements for the degree

MASTER OF SCIENCE

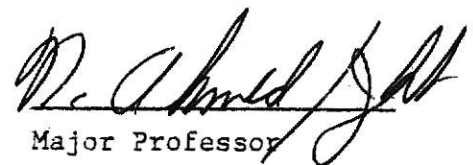
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CHAPTER I

INTRODUCTION

In practice it is impossible to encode a given data sequence exactly, since data words with infinite lengths would be required. However, very accurate representations of data sequences are achieved in digital computers by resorting to floating-point arithmetic. In contrast, accuracy presents serious problems when one is restricted to using fixed-point arithmetic, as is the case in special purpose hardware. Since fixed-point arithmetic offers faster execution time and more economical hardware, methods have been sought to encode data sequences in an efficient manner using fixed-point arithmetic. One such method is attributed to Max [1]. This method is optimum when the distortion measure is the mean square error criterion, and the data sequence to be encoded can be described by a gaussian random variable.

The purpose of this report is to develop a set of computer programs that implement Max's method on the computer system in the Electrical Engineering Department's Signal Processing Laboratory. The motivation for doing so is that the computer programs can be readily used for encoding purposes in applications such as the compression of image and speech data.

Chapter II is devoted to an introduction of Max's method, while its relevance to encoding gaussian random variables is addressed in Chapter III. Illustrative examples are included along with the computer program documentation presented in Chapter IV. Chapter V supplements the documentation with a user's guide. Some conclusions and recommendations for future work are summarized in Chapter VI.

CHAPTER II

THE OPTIMUM QUANTIZER

A quantizer enables one to encode a given discrete-time sequence of numbers, $x(n)$, in binary form. Why would anyone want to quantize? It is because only a finite number of bits are available to encode a given sequence. Thus, in practice, it is impossible to encode $x(n)$ exactly. As a consequence, efficient quantization techniques are sought to use a finite number of bits to encode $x(n)$ in such a way that the resulting digital sequence $\hat{x}(n)$ is a faithful reconstruction of $x(n)$. An optimum quantizer is one which produces minimum distortion with respect to some criterion--e.g., the mean square error. The objective is that the output sequence from the quantizer is as close a representation of the input sequence as possible.

For convenience, we let the input to the quantizer be a continuous random variable X ; i.e., X takes a continuum of values which we wish to quantize into a finite set of discrete values Y_i . In practice, the entire range of possible input values is broken up into smaller subranges. Each subrange has a lower limit, X_i and upper limit X_{i+1} . Any input X between these two limits is mapped into corresponding output Y_i . We consider the cost function

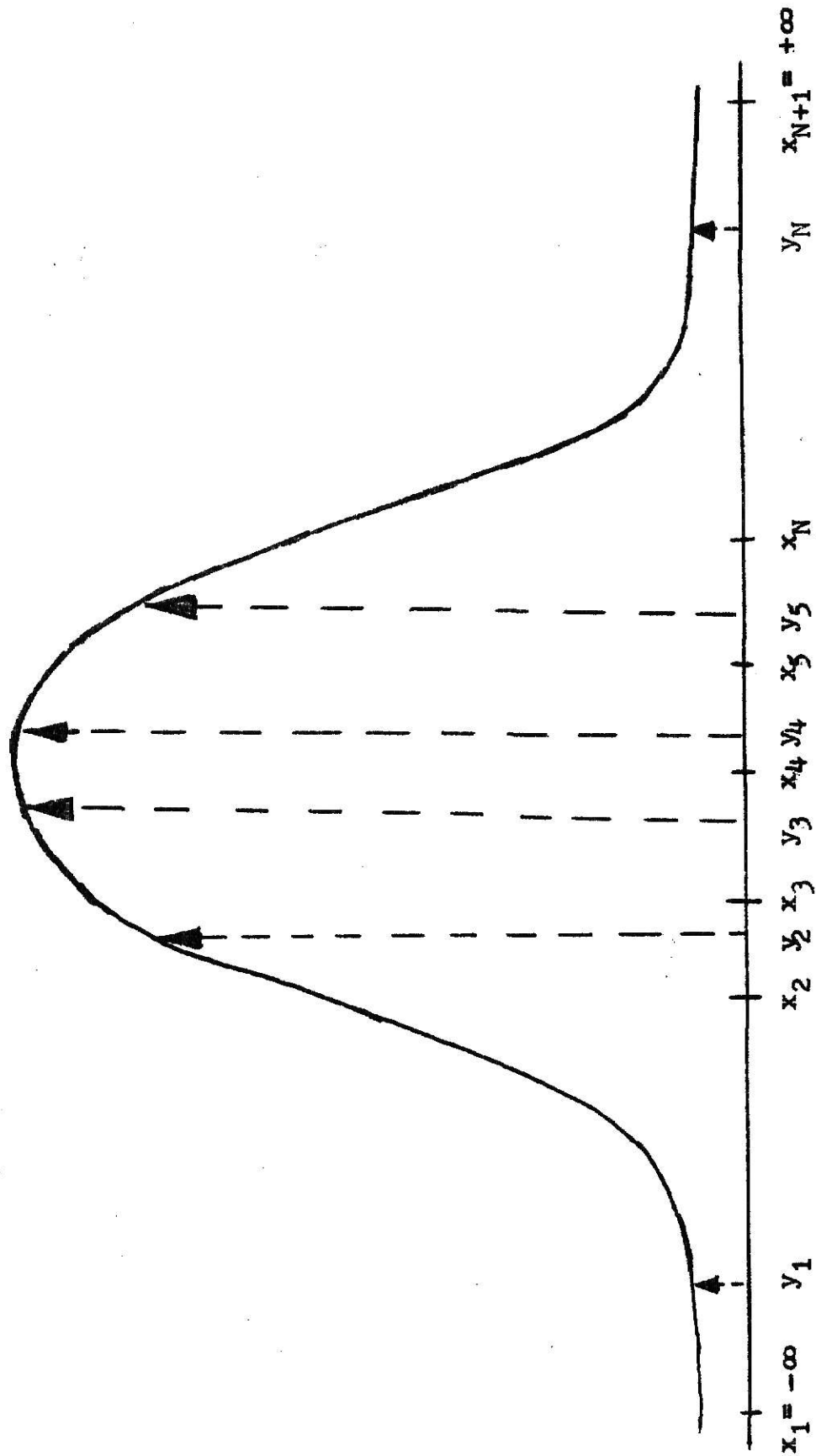
$$E((x - y)^2) = \sum_{i=1}^N \int_{x_i}^{x_{i+1}} (x - y_i)^2 p_X(x) dx \quad (2-1)$$

where $p_X(x)$ denotes the probability density function of X , and N is the number of finite output levels. In the most general case $x_1 = -\infty$ and $x_{N+1} = +\infty$, as illustrated in Figure 2.1. Simplifying for a symmetric distribution, $p_X(x)$, such as the gaussian distribution, x_{N+1} still equals $+\infty$ but

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Note that the input abscissa (x_i) is continuous, while the output abscissa (y_i) is discrete.



OUTPUT LEVELS, GENERAL DISTRIBUTION

FIGURE 2.1

$$\begin{aligned}
 &\text{For N} \\
 &\text{EVEN} \left\{ \begin{aligned} x_N &= 0 \\ \frac{x_N}{2} + 1 &= 0 \\ y_N - i &= -\frac{y_N}{2} + (i + 1) \\ \frac{x_N}{2} - (i - 1) &= -\frac{x_N}{2} + (i + 1) \end{aligned} \right. \\
 &\text{For N} \\
 &\text{ODD} \left\{ \begin{aligned} \frac{y_{(N+1)}}{2} &= 0 \\ \frac{x_{(N+1)}}{2} + i &= -\frac{x_{(N+1)}}{2} - (i - 1) \\ \frac{y_{(N+1)}}{2} + i &= -\frac{y_{(N+1)}}{2} - i \end{aligned} \right.
 \end{aligned}$$

as depicted in Figure 2.2. Equation (2-1) can also be written as

$$\begin{aligned}
 E((X - Y)^2) &= \sum_{j=2}^{N+1} \int_{x_{j-1}}^{x_j} (x - y_{j-1})^2 p_X(x) dx \\
 &= \int_{x_1}^{x_2} (x - y_1)^2 p_X(x) dx + \int_{x_2}^{x_3} (x - y_2)^2 p_X(x) dx + \dots \\
 &+ \int_{x_{j-1}}^{x_j} (x - y_{j-1})^2 p_X(x) dx + \int_{x_j}^{x_{j+1}} (x - y_j)^2 p_X(x) dx + \dots \\
 &+ \int_{x_N}^{x_{N+1}} (x - y_N)^2 p_X(x) dx \tag{2-2}
 \end{aligned}$$

The x_i 's and y_i 's are determined when (2-2) is minimized. Conditions for a minimum are discussed in Appendix 2.1. Taking the partial with respect to x_j , only two terms survive since the other terms are constant with respect to x_j and the partial of a constant is zero. Thus (2-2) yields

OUTPUT LEVELS, SYMMETRIC DISTRIBUTION

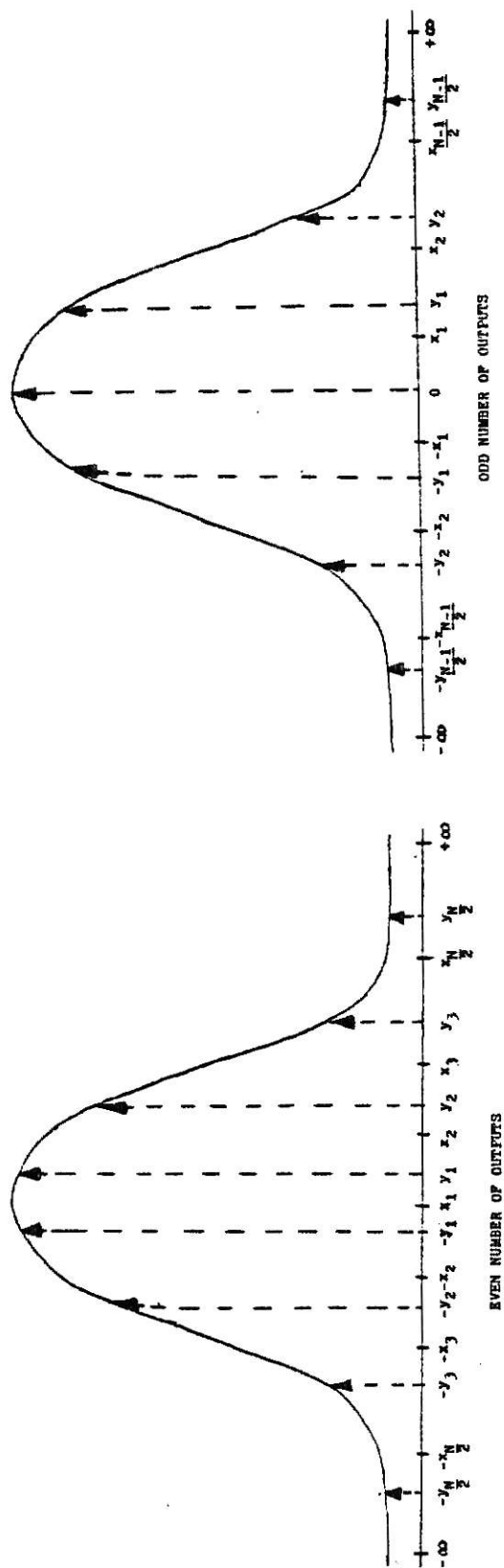


FIGURE 2.2

$$\frac{\partial E((X-Y)^2)}{\partial x_j} = \frac{\partial}{\partial x_j} \int_{x_{j-1}}^{x_j} (x-y_{j-1})^2 p_X(x) dx + \frac{\partial}{\partial x_j} \int_{x_j}^{x_{j+1}} (x-y_j)^2 p_X(x) dx \quad (2-3)$$

To evaluate the partial derivative in (2-3), we utilize Liebnitz's rule, which states that

$$\frac{d}{dx} \int_{a(x)}^{b(x)} g(z) dz = g(b(x)) \frac{db(x)}{dx} - g(a(x)) \frac{da(x)}{dx} \quad (2-4)$$

Noting that $\frac{\partial x_{j-1}}{\partial x_j} = 0$, and applying (2-4) to (2-3) results in

$$\begin{aligned} \frac{\partial E((X-Y)^2)}{\partial x_j} &= (x_j - y_{j-1})^2 p_X(x_j) (1) - (x_{j-1} - y_{j-1})^2 p_X(x_{j-1}) (0) \\ &\quad + (x_{j+1} - y_j)^2 p_X(x_{j+1}) (0) - (x_j - y_j)^2 p_X(x_j) (1) \end{aligned}$$

or

$$\frac{\partial E((X-Y)^2)}{\partial x_j} = (x_j - y_{j-1})^2 p_X(x_j) - (x_j - y_j)^2 p_X(x_j) \quad (2-5)$$

From (2-5) it is clear that one condition for minimizing the mean square error is

$$(x_j - y_{j-1})^2 = (x_j - y_j)^2$$

or

$$y_j = 2x_j - y_{j-1} \quad \text{for } j=2, \dots, N \quad (2-6)$$

Now for the partial with respect to y_j . Only one term y_j varies from y_1 to y_N . Thus (2-1) yields

$$\frac{\partial E((X-Y)^2)}{\partial y_j} = \int_{x_j}^{x_{j+1}} 2(x-y_j)(-1) p_X(x) dx \quad (2-7)$$

From (2-7) the second condition for the minimization of the mean square error is obtained as

$$\int_{x_j}^{x_{j+1}} (x-y_j) p_X(x) dx = 0 \quad \text{for } j=1, \dots, N \quad (2-8)$$

Thus to summarize, (2-6) and (2-8) completely describe the optimum quantizer, since their simultaneous solution yields all subranges of the possible input values, X_i and X_{i+1} , and their corresponding assigned output value, Y_i . A general solution is difficult; i.e., for any $p_X(x)$. However, an iterative procedure [1] can produce solutions for a specified $p_X(x)$. The procedure is to choose initial values x_1 and y_1 , then compute the set of x_i and y_i using (2-6) and (2-8). Once the last values, x_N and y_N , have been computed, y_N is compared to the average abscissa between x_N and infinity since (2-8) can be written

$$y_j = \frac{\int_{x_j}^{x_{j+1}} x p_X(x) dx}{\int_{x_j}^{x_{j+1}} p_X(x) dx} \quad \text{for } j=1, \dots, N \quad (2-9)$$

where y_j is recognized as the average abscissa corresponding to the centroid of the area between x_j and x_{j+1} . If the two values of y_N are within a specified accuracy, then the optimum quantizer has been characterized. Otherwise the initial values, x_1 and y_1 , are chosen again based on the difference between these two values for y_N . The remaining x_i 's and y_i 's are then recalculated. The entire procedure continues until the two values for y_N are within a specified accuracy.

APPENDIX 2.1

This appendix contains the conditions necessary to ensure minimization of an equation containing two variables. Conditions to ensure minimization of (2-1). A minimum occurs for the critical points where

$$\frac{\partial E}{\partial x_i} = \frac{\partial E}{\partial y_i} = 0$$

if

$$\begin{vmatrix} \frac{\partial^2 E}{\partial x_i^2} & \frac{\partial^2 E}{\partial y_i \partial x_i} \\ \frac{\partial^2 E}{\partial x_i \partial y_i} & \frac{\partial^2 E}{\partial y_i^2} \end{vmatrix} > 0$$

and if

$$\frac{\partial^2 E}{\partial x_i^2} > 0, \quad \frac{\partial^2 E}{\partial y_i^2} > 0$$

Applying the minimization criteria to (2-1) we obtain

$$E = \sum_{i=1}^N \int_{x_i}^{x_{i+1}} (y_i - x)^2 p_X(x) dx$$

$$\frac{\partial^2 E}{\partial y_i^2} = \int_{x_i}^{x_{i+1}} 2 p_X(x) dx$$

$$\therefore \frac{\partial^2 E}{\partial y_i^2} > 0 \text{ ALWAYS}$$

$$\frac{\partial^2 E}{\partial x_i^2} = 2p_X(x_i) [y_i - y_{i-1}]$$

$$+ \left[\frac{d}{dx_i} p(x_i) \right] [y_i - y_{i-1}]$$

$$[(-1) (y_{i-1} + y_i) + 2x_i]$$

$$\frac{\partial^2 E}{\partial x_i \partial y_i} = \frac{\partial^2 E}{\partial y_i \partial x_i} = \underbrace{2p(x_i)[y_{i-1} - y_i]}_{\text{ALWAYS NEGATIVE}}$$

CHAPTER III

THE OPTIMUM GAUSSIAN QUANTIZER

The gaussian distribution is frequently assumed for $p_X(x)$ in (2-6) and (2-8) that yield the optimum quantizer. A set of tables containing the x_i and y_i parameters for the gaussian case were generated by Max [1]. In this chapter we discuss some details related to generating Max's tables.

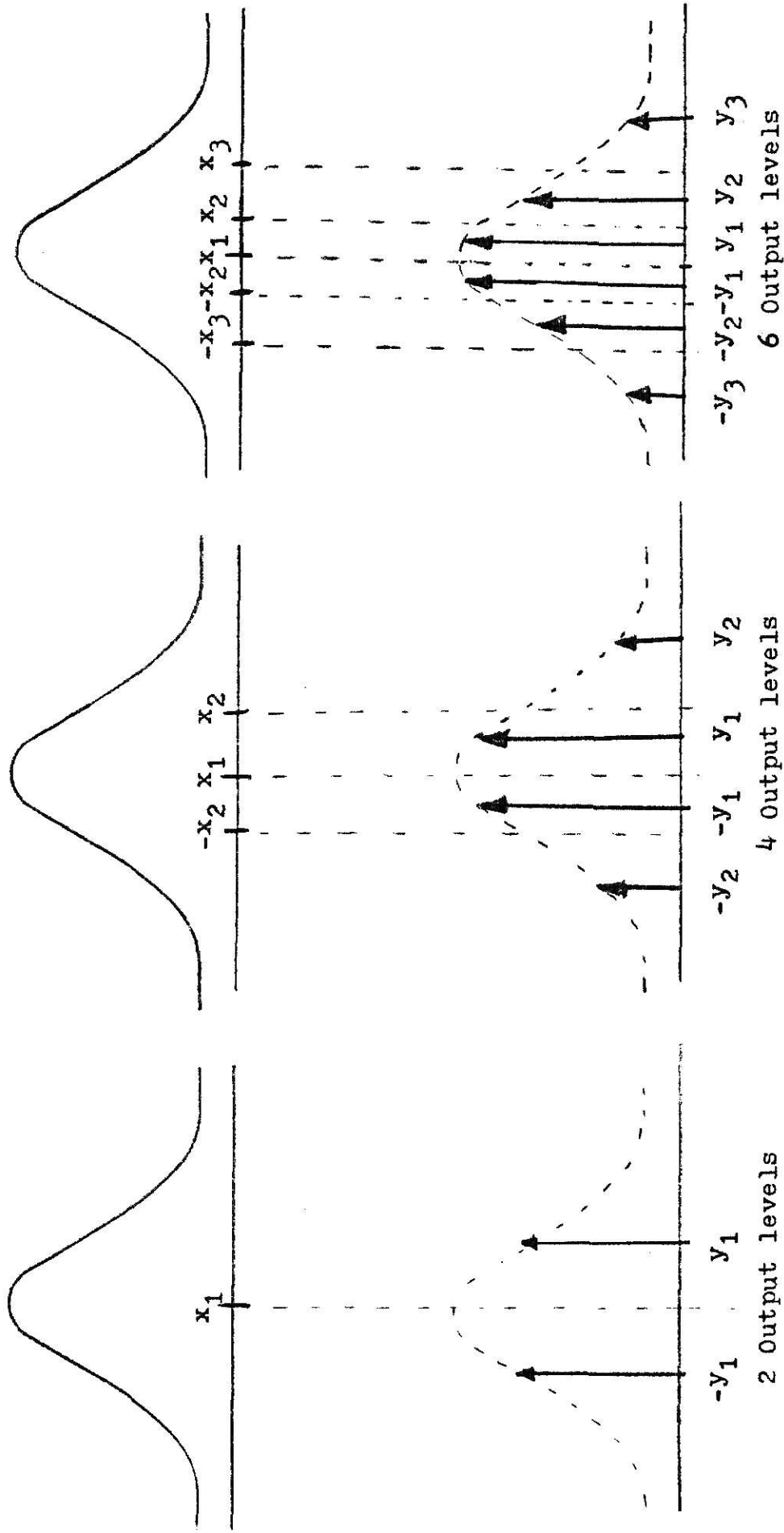
The critical portion of the iterative procedure to obtain the optimum quantizer for a given number of output levels N , is the initial guess for x_1 and y_1 . It should be pointed out that only one of these values is chosen. The other value is set equal to zero, forcing the output of the quantizer to be symmetric about the origin, as illustrated in Figure 3.1. What about the value to be chosen? To this end, one can refer to Max's tables which are available for values of N from 2 through 36 [1]. It is convenient to obtain an empirical relationship between the initial values published by Max [1] and the number of output levels N . Hein [3] has obtained a handy relationship via a least squares fit of the initial values published by Max [1]. This relationship is given by [3]

$$P_1 = \frac{2.1613}{N} - \frac{1.653}{N^2} \quad (3-1)$$

where P_1 is x_1 if N is odd, or y_1 if N is even. Values for P_1 are tabulated and compared to Max's initial values in Table 3.1. Clearly the results compare closely.

The remaining x_i 's and y_i 's are obtained from (2-6) and (2-8); i.e.,

EVEN NUMBER OF OUTPUTS (x_1 is always zero)

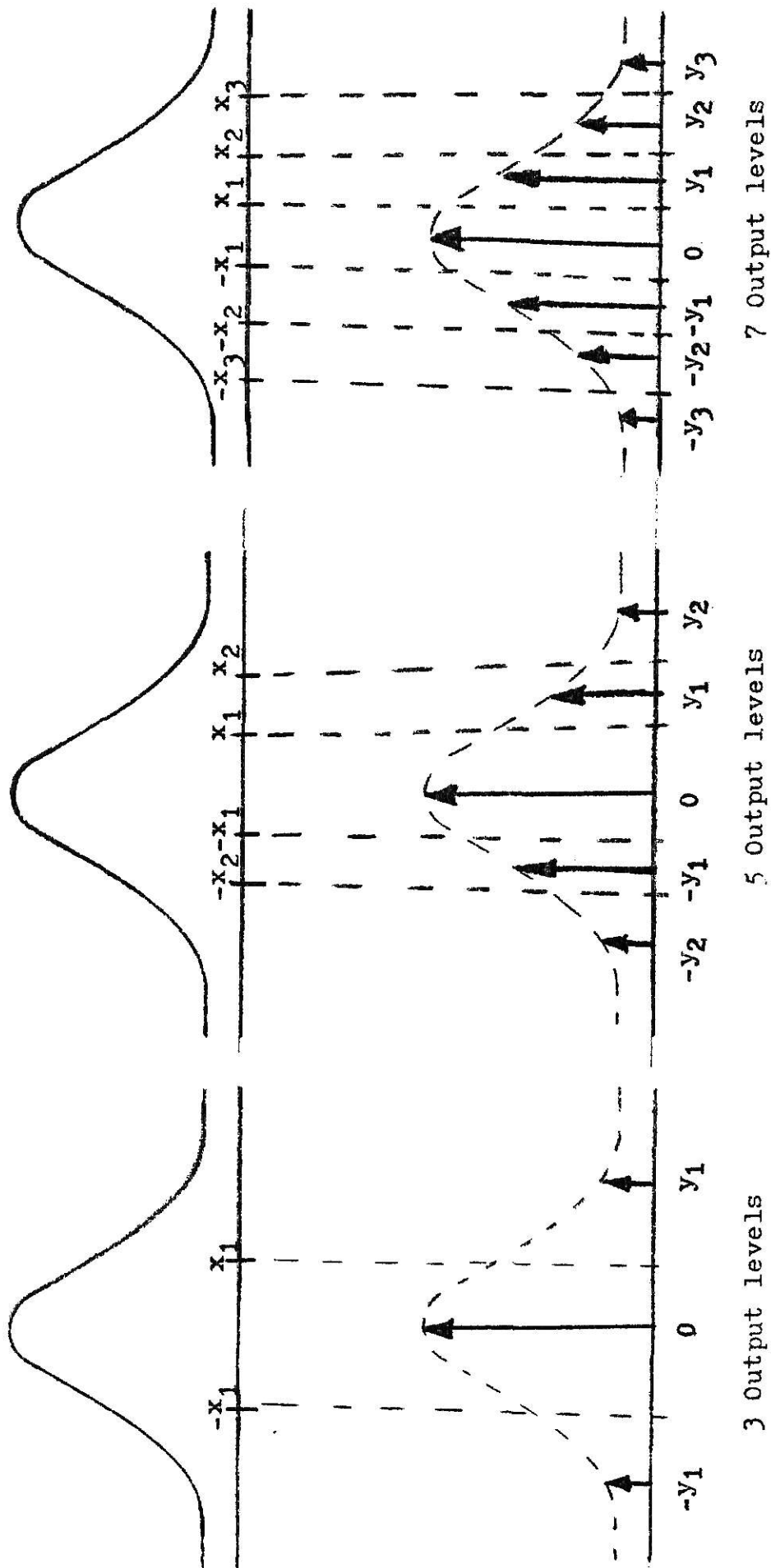


Note Y_i is the centroid of the area between X_i and X_{i+1} .

The positive values of X_i and Y_i are tabulated in Table 3.2, these values are symmetric about the origin.

FIGURE 3.1

ODD NUMBER OF OUTPUTS (Y_1 is always zero)



Note Y_i is the centroid of the area between X_i and X_{i+1} .
The positive values of X_i and Y_i are tabulated in Table 3.2, these values are symmetric about the origin

FIGURE 3.1 (CONT.)

TABLE 3.1
INITIAL VALUES

NUMBER OF OUTPUT LEVELS	INITIAL VALUES EQ. (3-1) $P_1 = \frac{2.1613}{N} - \frac{1.653}{N^2}$	INITIAL VALUES PUBLISHED BY MAX
4	0.43701	0.4528
5	0.36614	0.3823
6	0.31430	0.3177
7	0.27502	0.2803
8	0.24433	0.2451
9	0.21974	0.2218
10	0.19960	0.1996
11	0.18282	0.1837
12	0.16863	0.1684
13	0.15647	0.1569
14	0.14594	0.1457
15	0.13674	0.1369
16	0.12862	0.1284
17	0.12142	0.1215
18	0.11497	0.1148
19	0.10917	0.1092
20	0.10393	0.1038
21	0.09917	0.09918
22	0.09483	0.09469
23	0.09084	0.09085
24	0.08718	0.08708
25	0.08381	0.08381
26	0.08068	0.08060
27	0.07778	0.07779
28	0.07508	0.07502
29	0.07256	0.07257
30	0.07021	0.07016
31	0.06800	0.06802
32	0.06593	0.06500
33	0.06398	0.06400
34	0.06214	0.06212
35	0.06040	0.06043
36	0.05876	0.05876
64	0.03337	0.03342
128	0.01678	0.01685

$$y_j = 2x_j - y_{j-1} \quad \text{for } j=2, \dots, M$$

and

$$\int_{x_j}^{x_{j+1}} (x - y_j) p_X(x) dx = 0 \quad \text{for } j=1, \dots, M$$

where

$$M = \begin{cases} \frac{N}{2} & , N \text{ EVEN} \\ \frac{(N-1)}{2} & , N \text{ ODD} \end{cases}$$

due to symmetry, i.e, $x_j = -x_j$ and $y_j = -y_j$, as shown in Figure 3.2.

Let $p_X(x)$ be gaussianly distributed

$$p_X(x) = \frac{1}{\sqrt{2\pi}} e^{-\frac{x^2}{2}} \quad , \quad |x| < \infty \quad (3-2)$$

The value of x_{j+1} in (2-8) is determined via an iterative procedure for finding the zeroes of functions, Newton's method [5] which is described in Appendix 3.1. The actual integration in (2-8) is accomplished by expressing (2-8) as

$$\int_{x_j}^{x_{j+1}} x p_X(x) dx - y_j \int_{x_j}^{x_{j+1}} p_X(x) dx = 0 \quad (3-3)$$

where the first part is easily integrable, once it is observed that for our particular choice of $p_X(x)$

$$\frac{dp_X(x)}{dx} = -x p_X(x) \quad (3-4)$$

The second part is evaluated by replacing the integral with its equivalent infinite series representation [4] which converges very rapidly as shown in Figure 3.2. Thus if

$$S(x) = \frac{1}{\sqrt{\pi}} e^{-\frac{x^2}{2}} \sum_{k=0}^{\infty} \frac{(x/\sqrt{2})^{2k+1} 2^k}{(1)(3)(5)\cdots(2k+1)}$$

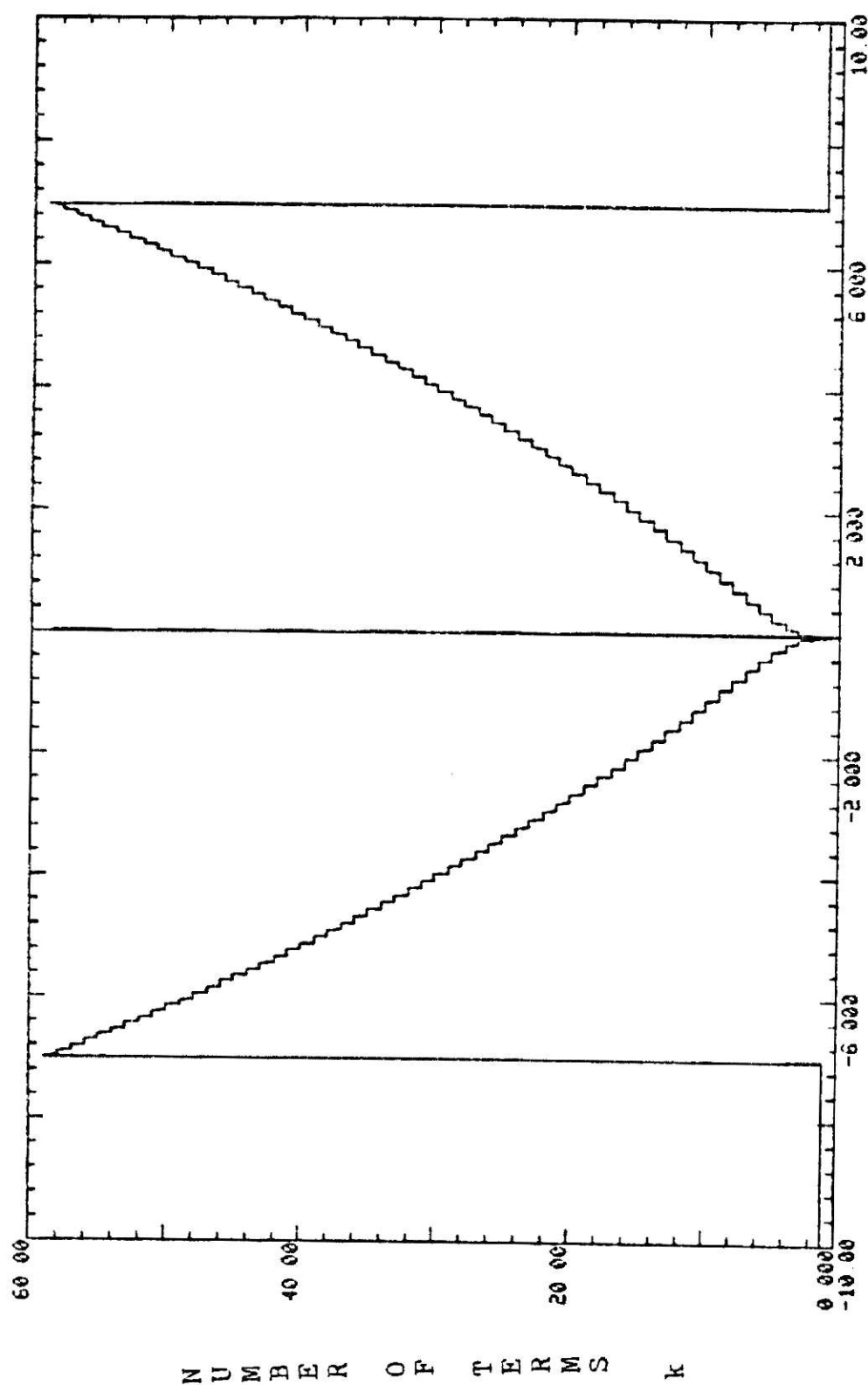


FIGURE 3.2

$$S(x) = \frac{1}{\sqrt{\pi}} e^{-\frac{x^2}{2}} \sum_{k=0}^{\infty} \frac{(x/\sqrt{2})^{2k+1} 2^k}{(1)(3)(5) \dots (2k+1)} \quad (3-5)$$

then

$$\int_{x_j}^{x_{j+1}} p_X(x) dx = S(x_{j+1}) - S(x_j) \quad (3-6)$$

Also, we note that

$$\int_{x_j}^{\infty} p_X(x) dx = S(+\infty) - S(x_j) = 0.5 - S(x_j) \quad (3-7)$$

since

$$\frac{1}{\sqrt{2\pi}} \int_0^{+\infty} e^{-\frac{x^2}{2}} dx = 0.5 \quad (3-8)$$

After the last x_i and y_i are calculated, y_M is also determined by an alternate method. The abscissa of the centroid of the area between x_M and infinity should also equal y_M , as seen in (2-9)--i.e.,

$$y_{M_c} = \frac{\int_{x_M}^{\infty} x p_X(x) dx}{\int_{x_M}^{\infty} p_X(x) dx}$$

The two values for y_M are compared. If they are close enough, then the optimum quantizer has been determined. Otherwise, the initial value must be updated and the x_i 's and y_i 's again computed. The initial value is updated using an iterative procedure for finding the zeroes of functions, ie force $y_M - y_{M_c}$ toward zero. Newton's method can't be used in this instance because an explicit expression for the first derivative of the function to be zeroed is not available. Thus a

similar iterative procedure must be used, the secant method [5] which is described in Appendix 3.2

$$P_1^{**} = \frac{P_1(y_M^* - y_{M_c}^*) - P_1^*(y_M - y_{M_c})}{(y_M^* - y_{M_c}^*) - (y_M - y_{M_c})} \quad (3-9)$$

where y_M is calculated from (2-6) and (2-8), while y_{M_c} is obtained using (2-9). In (3-9),

$$P_1 = \begin{cases} x_1 & \text{if } N \text{ is odd} \\ y_1 & \text{if } N \text{ is even} \end{cases}$$

Two asterisks denote the updated value, one asterisk denotes the current value, while no asterisk denotes the previous value. The secant method converges only if the initial values (current and last) are sufficiently close to the actual initial value of the optimum quantizer. However, when it does converge it converges very quickly [5]. For example, with $N=128$ output levels, the entire iterative process took only six iterations for an accuracy to six decimal places on the Data General Nova 4X computer.

The program to generate the parameters X_i and Y_i , which completely describe the optimum gaussian quantizer for a finite number of output levels, N , is listed in Appendix 3.3. Tables generated using this program are summarized in Table 3.2. A graphical representation of the input and output values for the optimum quantizer, tabulated for a specific N in Table 3.2, is perhaps more enlightening. Figures 3.4 thru 3.10 show these input/output transfer functions for $N=2$ thru $N=128$ respectively. Figure 3.11 shows the input/output transfer function for $N=5$.

THE OPTIMUM QUANTIZER

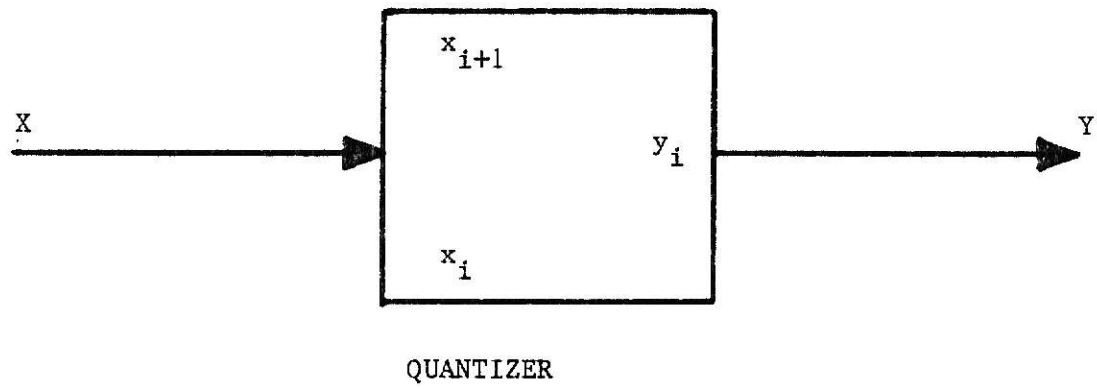


FIGURE 3.3

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TABLE 3.2

N= 2			N= 13		
J	X(J)	Y(J)	J	X(J)	Y(J)
1	.000000	.797885	1	.156887	.313773
N= 3			2	.476013	.638252
J	X(J)	Y(J)	3	.812602	.986953
1	.612003	1.224007	4	1.184109	1.381266
N= 4			5	1.622901	1.864535
J	X(J)	Y(J)	6	2.214547	2.564559
1	.000000	.452781	N= 14		
2	.981600	1.510419	J	X(J)	Y(J)
N= 5			1	.000000	.145708
J	X(J)	Y(J)	2	.293516	.441323
1	.382284	.764568	3	.595885	.750446
2	1.244357	1.724147	4	.913045	1.065643
N= 6			5	1.276590	1.467538
J	X(J)	Y(J)	6	1.703080	1.938623
1	.000000	.317717	7	2.281868	2.625112
2	.658912	1.000107	N= 15		
3	1.446853	1.893598	J	X(J)	Y(J)
N= 7			1	.136929	.273857
J	X(J)	Y(J)	2	.414311	.554765
1	.280289	.560577	3	.702951	.851138
2	.874363	1.188149	4	1.013013	1.174888
3	1.610763	2.033378	5	1.360479	1.546069
N= 8			6	1.776280	2.006491
J	X(J)	Y(J)	7	2.343715	2.680939
1	.000000	.245095	N= 16		
2	.500530	.756006	J	X(J)	Y(J)
3	1.049959	1.343911	1	.000000	.128397
4	1.747929	2.151946	2	.258224	.388051
N= 9			3	.522406	.656762
J	X(J)	Y(J)	4	.799554	.942347
1	.221819	.443639	5	1.099296	1.256245
2	.681218	.918797	6	1.437155	1.618065
3	1.197595	1.476392	7	1.843554	2.069043
4	1.865532	2.254671	8	2.400829	2.732615
N= 10			N= 17		
J	X(J)	Y(J)	J	X(J)	Y(J)
1	.000000	.199624	1	.121497	.242994
2	.404742	.609859	2	.366939	.490883
3	.833844	1.057829	3	.620086	.749290
4	1.324587	1.591345	4	.887446	1.025603
5	1.968230	2.345115	5	1.178253	1.330902
N= 11			6	1.507680	1.684458
J	X(J)	Y(J)	7	1.905731	2.127003
1	.183729	.367458	8	2.453907	2.780811
2	.559913	.752368	N= 18		
3	.965599	1.178830	J	X(J)	Y(J)
4	1.435738	1.692646	1	.000000	.114771
5	2.059198	2.425750	2	.230560	.346348
N= 12			3	.465327	.584306
J	X(J)	Y(J)	4	.709087	.833868
1	.000000	.168439	5	.967991	1.102113
2	.340144	.511849	6	1.250977	1.399840
3	.694316	.876784	7	1.572936	1.746631
4	1.081252	1.285721	8	1.963506	2.180981
5	1.534384	1.763047	9	2.503445	2.825908
6	2.140753	2.498459			

THE OPTIMUM GAUSSIAN QUANTIZER PARAMETERS

TABLE 3.2 (CONT.)

N=19			N=24		
J	X(J)	Y(J)	J	X(J)	Y(J)
1	.109205	.218409	1	.000000	.087075
2	.329383	.440357	2	.174590	.262105
3	.555079	.669801	3	.350982	.439859
4	.790737	.911674	4	.531118	.622378
5	1.042246	1.172817	5	.717236	.812095
6	1.315314	1.463810	6	.912102	1.012110
7	1.633597	1.803385	7	1.119376	1.226643
8	2.017416	2.231448	8	1.344255	1.461868
9	2.549850	2.868252	9	1.594795	1.727721
			10	1.884872	2.042022
			11	2.242618	2.443213
			12	2.745395	3.047577
N=20			N=25		
J	X(J)	Y(J)	J	X(J)	Y(J)
1	.000000	.103765	1	.083805	.167611
2	.208280	.312794	2	.252210	.336808
3	.418643	.528492	3	.423048	.509288
4	.637515	.748539	4	.598134	.686980
5	.864096	.983652	5	.779603	.872225
6	1.111070	1.238487	6	.970134	1.068042
7	1.380962	1.523438	7	1.173310	1.278578
8	1.690218	1.855998	8	1.394255	1.509932
9	2.067866	2.278734	9	1.640941	1.771950
10	2.593412	2.908089	10	1.927137	2.082325
			11	2.280796	2.477268
			12	2.778845	3.078423
N=21			N=26		
J	X(J)	Y(J)	J	X(J)	Y(J)
1	.099179	.198357	1	.000000	.080596
2	.298857	.399358	2	.161540	.242484
3	.502625	.605894	3	.324503	.406522
4	.713670	.821446	4	.490408	.574295
5	.936004	1.050562	5	.660970	.747645
6	1.175154	1.299746	6	.838242	.928539
7	1.439494	1.579242	7	1.024834	1.120830
8	1.743296	1.907350	8	1.224257	1.327684
9	2.115335	2.323319	9	1.441577	1.555469
10	2.634465	2.945612	10	1.684685	1.813901
			11	1.967265	2.120629
			12	2.317117	2.513606
			13	2.810743	3.107880
N=22			N=27		
J	X(J)	Y(J)	J	X(J)	Y(J)
1	.000000	.094689	1	.077781	.155562
2	.189945	.285201	2	.233977	.312391
3	.382218	.479235	3	.392110	.471828
4	.579363	.679490	4	.553600	.635372
5	.784387	.889284	5	.720083	.804793
6	1.001159	1.113034	6	.893548	.982302
7	1.235071	1.357108	7	1.076545	1.170789
8	1.494378	1.631848	8	1.272530	1.374271
9	1.793210	1.954771	9	1.486513	1.593756
10	2.160102	2.365433	10	1.726320	1.853884
11	2.673377	2.981322	11	2.005543	2.157203
			12	2.351832	2.546462
			13	2.841302	3.136142
N=23					
J	X(J)	Y(J)			
1	.090844	.181688			
2	.273545	.365401			
3	.459364	.533328			
4	.650872	.748017			
5	.850341	.952664			
6	1.062122	1.171590			
7	1.291364	1.411028			
8	1.546034	1.681040			
9	1.840308	1.999576			
10	2.202451	2.405326			
11	2.710289	3.015252			

TABLE 3.2 (CONT.)

N= 28			N= 32		
J	X(J)	Y(J)	J	X(J)	Y(J)
1	.000000	.075016	1	.000000	.065894
2	.150311	.225806	2	.131975	.198057
3	.301765	.377924	3	.264720	.331384
4	.455574	.533225	4	.399045	.466706
5	.613084	.692943	5	.535825	.604943
6	.775866	.858789	6	.676046	.747149
7	.945854	1.032919	7	.820867	.894585
8	1.125551	1.218183	8	.971699	1.048814
9	1.318363	1.418544	9	1.130332	1.211851
10	1.529252	1.639960	10	1.299120	1.386389
11	1.765982	1.892004	11	1.481341	1.578292
12	2.042030	2.192055	12	1.681793	1.787293
13	2.364871	2.577687	13	1.908057	2.028521
14	2.670376	3.163065	14	2.173340	2.317860
			15	2.504587	2.691315
			16	2.976185	3.261056

N= 29			N= 33		
J	X(J)	Y(J)	J	X(J)	Y(J)
1	.072566	.145132	1	.063990	.127980
2	.218212	.291292	2	.192322	.256663
3	.365425	.439558	3	.321722	.386781
4	.515340	.591122	4	.452951	.519121
5	.669241	.747360	5	.586844	.654567
6	.828647	.909935	6	.724355	.794143
7	.995448	1.080982	7	.866613	.939084
8	1.172105	1.263248	8	1.015004	1.090925
9	1.361981	1.460714	9	1.171293	1.251661
10	1.569974	1.679274	10	1.337808	1.423955
11	1.803854	1.928434	11	1.517301	1.611646
12	2.076956	2.225478	12	1.716040	1.820433
13	2.416638	2.607798	13	1.940066	2.059700
14	2.898229	3.188661	14	2.203026	2.346353
			15	2.531752	2.717152
			16	3.000350	3.283548

N= 30			N= 34		
J	X(J)	Y(J)	J	X(J)	Y(J)
1	.000000	.070159	1	.000000	.062118
2	.140546	.210932	2	.124393	.186667
3	.282023	.353114	3	.249426	.312189
4	.425417	.497719	4	.375782	.439375
5	.571802	.645884	5	.504176	.568977
6	.722412	.798939	6	.635414	.701850
7	.878724	.958510	7	.770421	.838992
8	1.042590	1.126670	8	.910303	.981613
9	1.216423	1.306177	9	1.058421	1.131229
10	1.403564	1.500951	10	1.210508	1.289787
11	1.608897	1.716844	11	1.374882	1.459977
12	1.840074	1.963303	12	1.552783	1.645589
13	2.110414	2.257524	13	1.748937	1.852285
14	2.447124	2.636725	14	1.970850	2.089415
15	2.925205	3.213685	15	2.231610	2.373805
			16	2.557943	2.742061
			17	3.023688	3.305295

N= 31		
J	X(J)	Y(J)
1	.068008	.136016
2	.204447	.272878
3	.342172	.411467
4	.482105	.552744
5	.625274	.697803
6	.772870	.847936
7	.926335	1.004733
8	1.087487	1.170240
9	1.258708	1.347177
10	1.443307	1.539436
11	1.646133	1.752830
12	1.874790	1.996749
13	2.142526	2.288304
14	2.476428	2.664553
15	2.951180	3.237808

N= 35			N=128		
J	X(J)	Y(J)	J	X(J)	Y(J)
1	.060421	.120842	1	.000000	.016846
2	.181558	.242275	2	.033680	.050513
3	.303594	.364912	3	.067356	.084199
4	.427161	.489409	4	.101058	.117917
5	.552749	.616490	5	.134799	.151681
6	.681736	.746983	6	.168591	.185502
7	.814424	.881865	7	.202448	.219394
8	.952094	1.022324	8	.236382	.253371
9	1.096092	1.169859	9	.270408	.287446
10	1.248128	1.326396	10	.304539	.321633
11	1.410501	1.494606	11	.338790	.355948
12	1.586430	1.678255	12	.373174	.390401
13	1.780614	1.882974	13	.407706	.425011
14	2.000525	2.118076	14	.442402	.459793
15	2.259176	2.400316	15	.477277	.494762
16	2.583252	2.766188	16	.512348	.529935
17	3.046278	3.326368	17	.547632	.565329
			18	.583145	.600961
			19	.618907	.636852
			20	.654935	.673019
			21	.691252	.709484
			22	.727876	.746268
			23	.764830	.783393
			24	.802139	.820884
			25	.839825	.858766
			26	.877916	.897066
			27	.916440	.935813
			28	.955425	.975037
			29	.994904	1.014771
			30	1.034911	1.055050
			31	1.075482	1.095913
			32	1.116657	1.137400
			33	1.158478	1.179556
			34	1.200950	1.222343
			35	1.244120	1.265897
			36	1.288086	1.310274
			37	1.332906	1.355537
			38	1.378646	1.401755
			39	1.425380	1.449005
			40	1.473197	1.497370
			41	1.522158	1.546946
			42	1.572393	1.597840
			43	1.624005	1.650171
			44	1.677036	1.703901
			45	1.731625	1.759349
			46	1.788020	1.816691
			47	1.846411	1.876132
			48	1.907021	1.937910
			49	1.970169	2.002307
			50	2.035984	2.069661
			51	2.104834	2.140007
			52	2.177086	2.214164
			53	2.253450	2.292736
			54	2.334611	2.376485
			55	2.421100	2.465715
			56	2.513999	2.562284
			57	2.615103	2.667923
			58	2.725862	2.783802
			59	2.849152	2.914503
			60	2.988918	3.063329
			61	3.150659	3.237989
			62	3.345521	3.453054
			63	3.597598	3.742741
			64	3.974031	4.205320

N= 36		
J	X(J)	Y(J)
1	.000000	.058753
2	.117636	.176520
3	.235814	.295109
4	.355105	.415100
5	.476110	.537119
6	.599493	.661867
7	.726008	.790150
8	.856538	.922926
9	.992145	1.061365
10	1.134157	1.206949
11	1.284267	1.361585
12	1.444756	1.527927
13	1.618824	1.709721
14	1.811144	1.912567
15	2.029155	2.145743
16	2.285839	2.425936
17	2.607726	2.789517
18	3.068161	3.346805

N= 64		
J	X(J)	Y(J)
1	.000000	.033419
2	.066854	.100290
3	.133800	.167310
4	.200947	.234583
5	.268398	.302212
6	.336259	.370306
7	.404641	.438977
8	.473662	.508347
9	.543445	.578542
10	.614123	.649703
11	.685841	.721980
12	.758759	.795538
13	.833050	.870562
14	.908910	.947259
15	.986561	1.025864
16	1.066254	1.106645
17	1.148279	1.189913
18	1.232951	1.275969
19	1.320668	1.365346
20	1.411925	1.458504
21	1.507301	1.556097
22	1.607507	1.658916
23	1.713393	1.767870
24	1.826106	1.884341
25	1.947202	2.010062
26	2.078662	2.147262
27	2.223400	2.299539
28	2.385699	2.471860
29	2.572499	2.673138
30	2.795894	2.918650
31	3.080666	3.242681
32	3.495806	3.748931

TABLE 3.2 (CONT.)

TRANSFER FUNCTION
(OPTIMUM GAUSSIAN QUANTIZER)

2 OUTPUT LEVELS

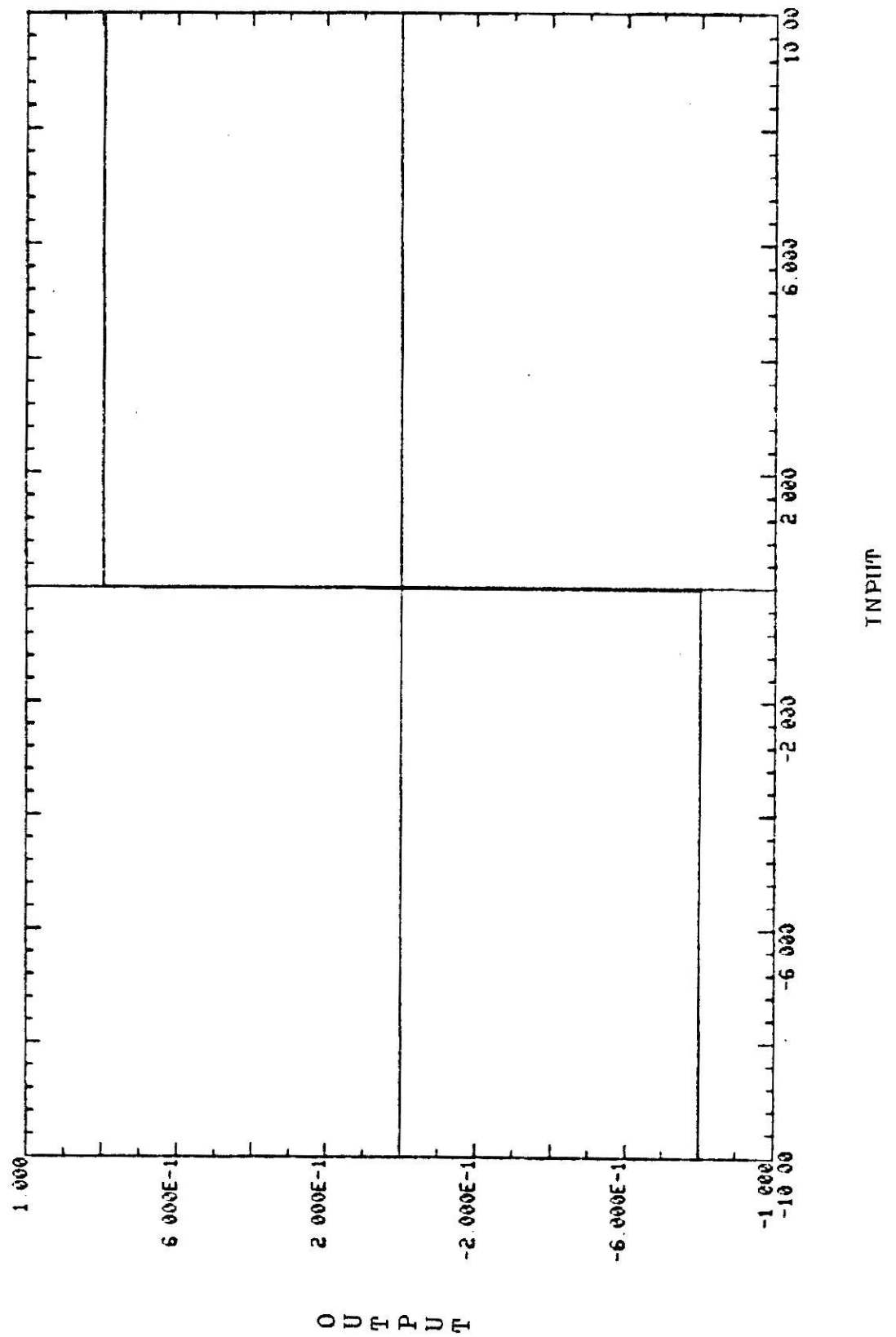


FIGURE 3.4

TRANSFER FUNCTION
(OPTIMUM GAUSSIAN QUANTIZER)

4 OUTPUT LEVELS

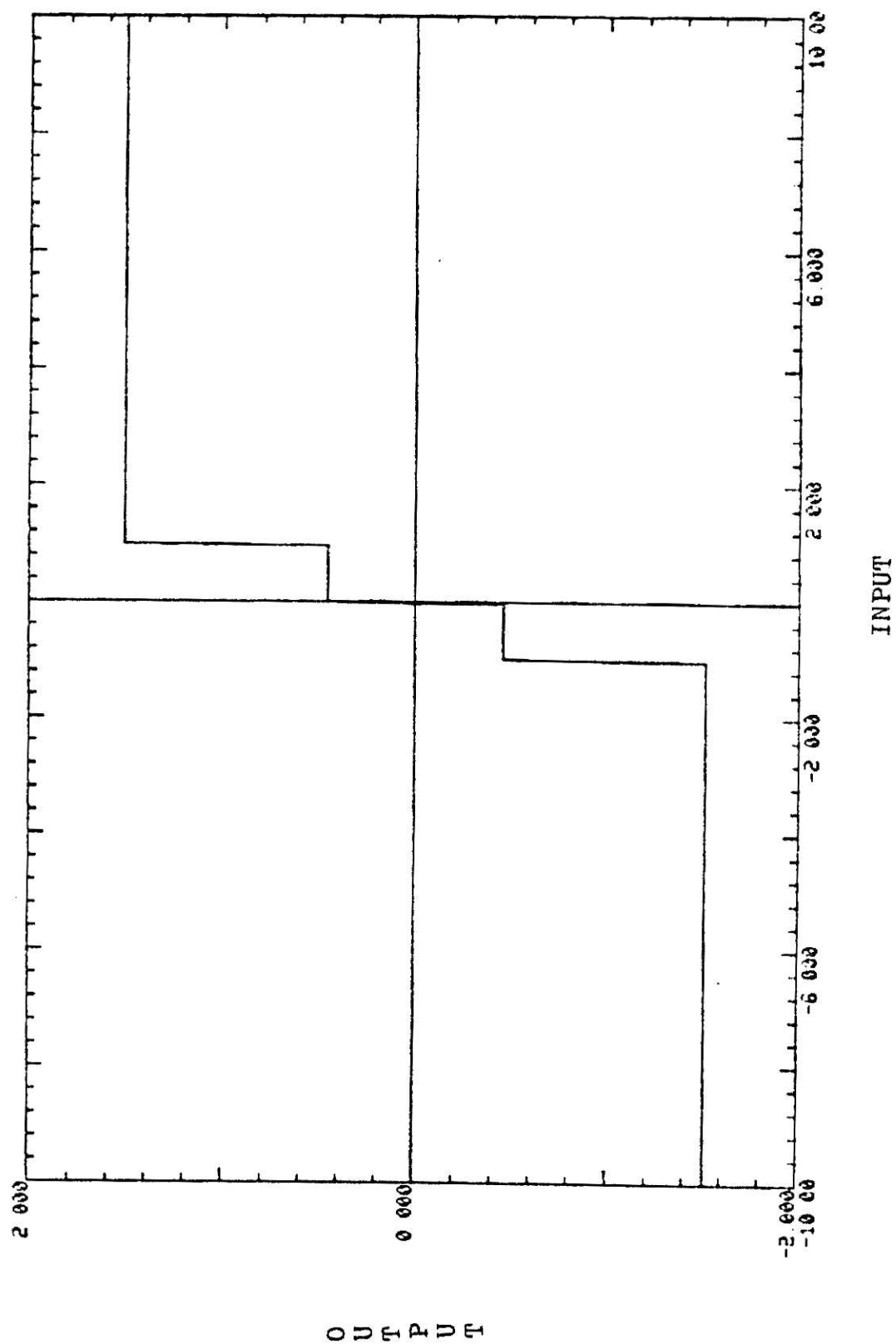


FIGURE 3.5

TRANSFER FUNCTION
(OPTIMUM GAUSSIAN QUANTIZER)

8 OUTPUT LEVELS

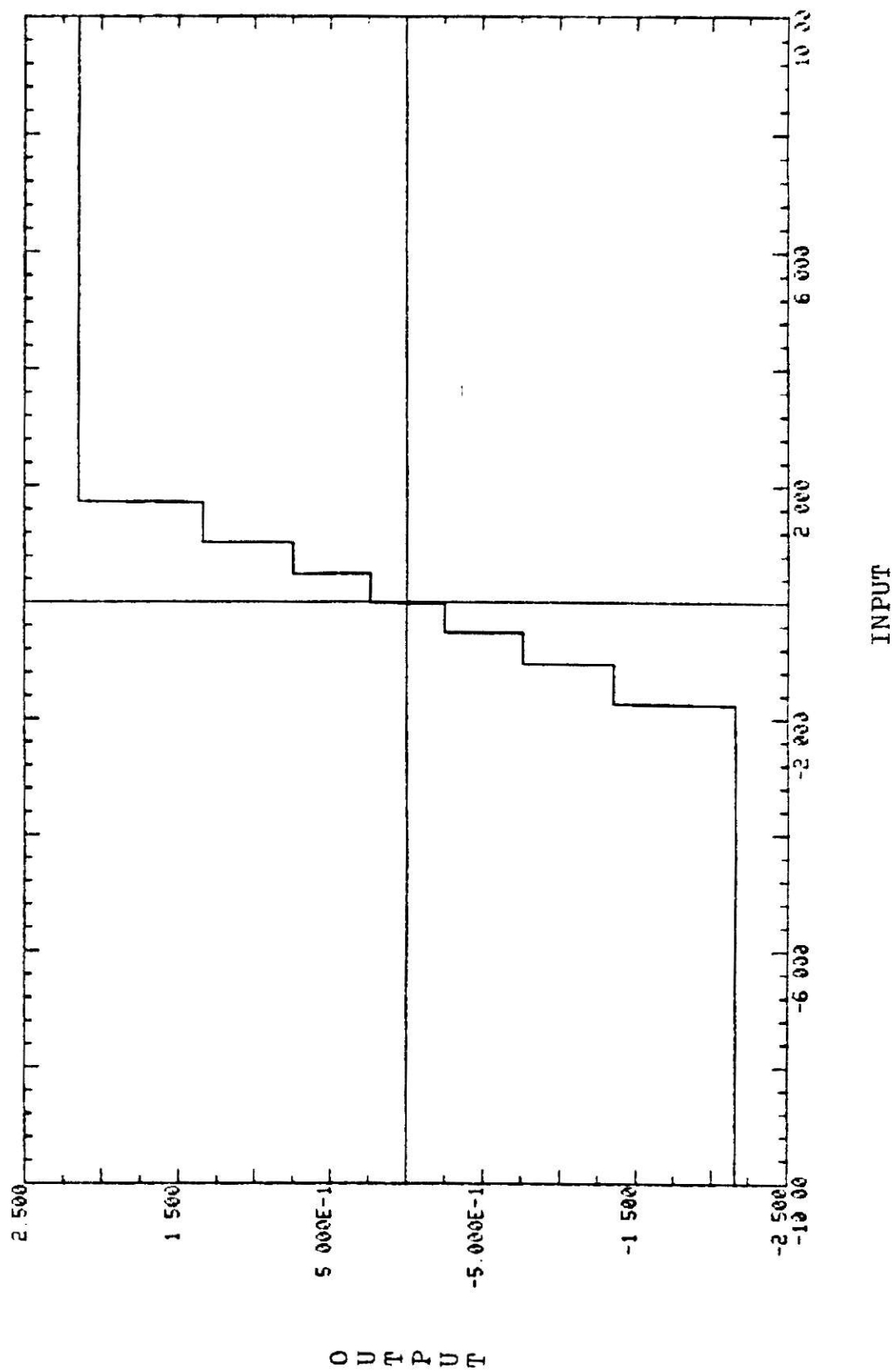


FIGURE 3.6

TRANSFER FUNCTION
(OPTIMUM GAUSSIAN QUANTIZER)

16 OUTPUT LEVELS

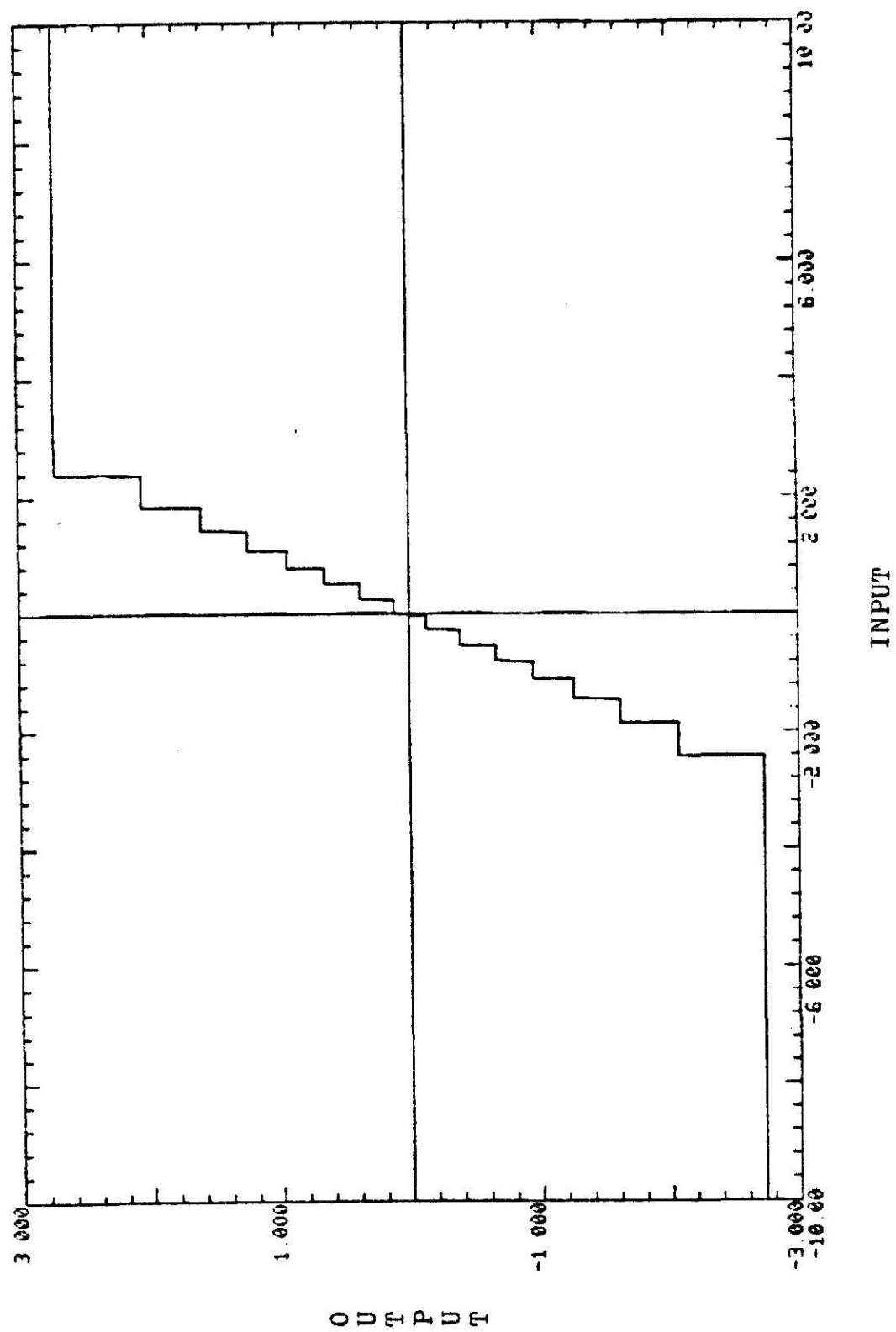


FIGURE 3.7

TRANSFER FUNCTION
(OPTIMUM GAUSSIAN QUANTIZER)

32 OUTPUT LEVELS

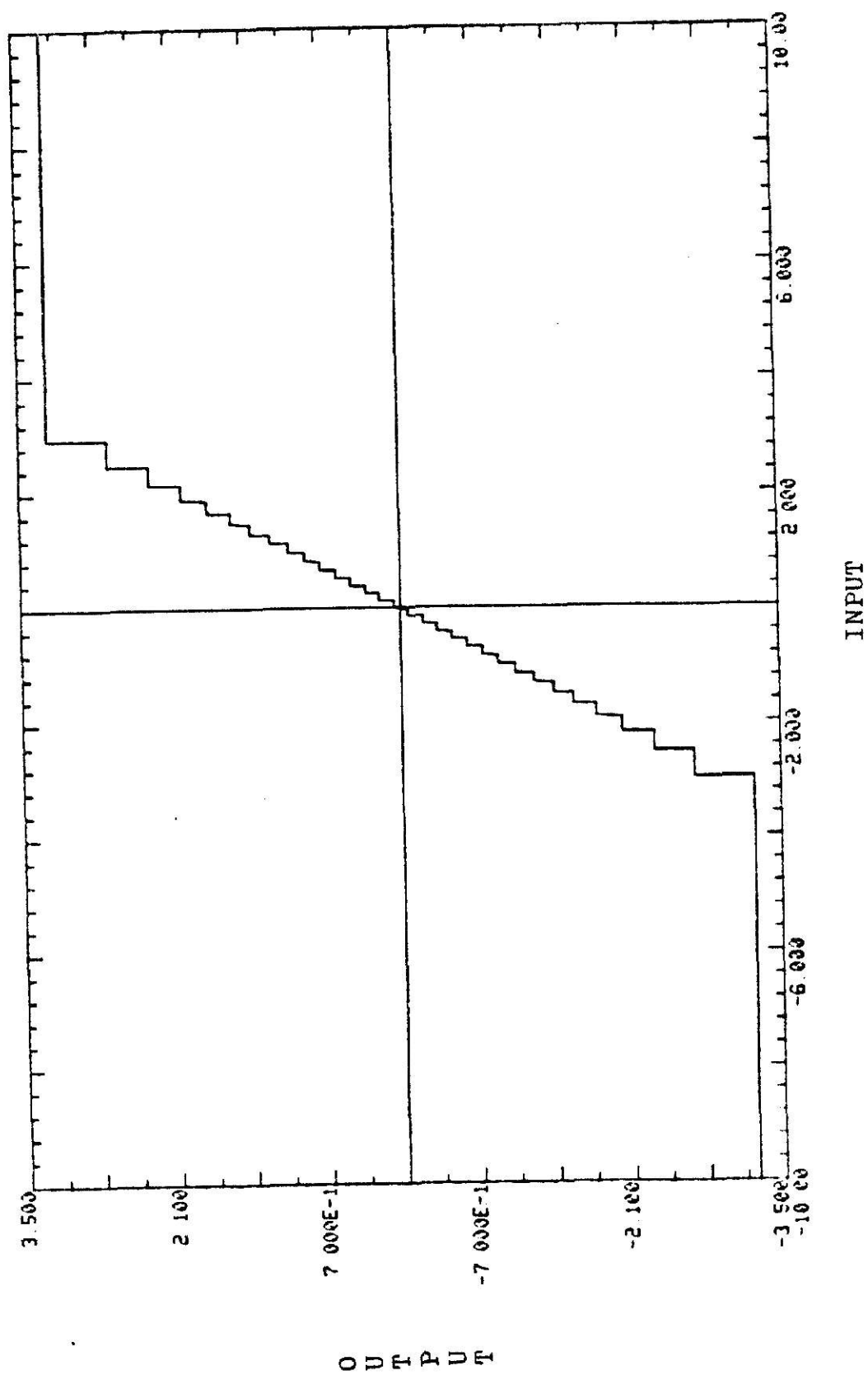


FIGURE 3.8

TRANSFER FUNCTION
(OPTIMUM GAUSSIAN QUANTIZER)

64 OUTPUT LEVELS

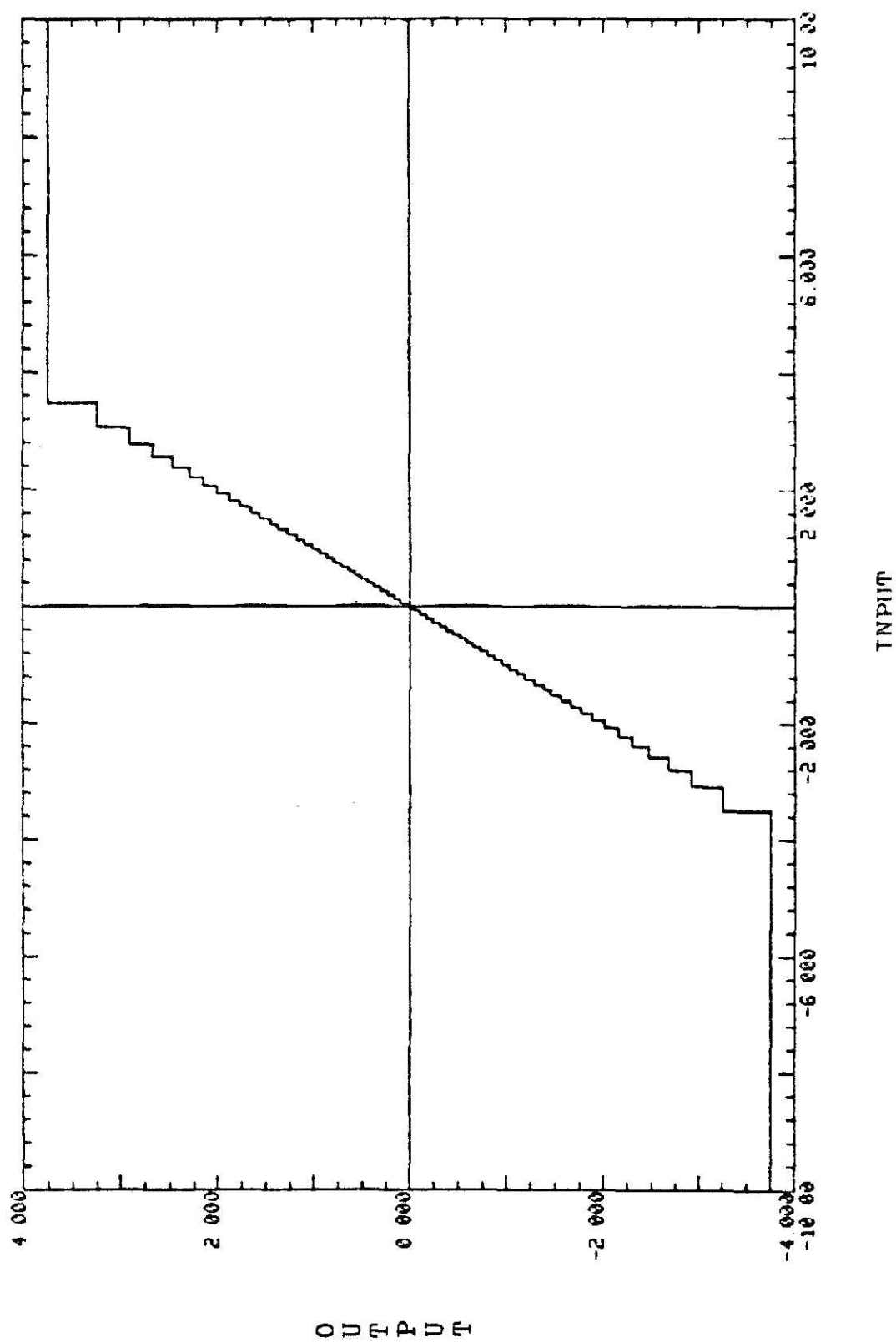


FIGURE 3.9

TRANSFER FUNCTION
(OPTIMUM GAUSSIAN QUANTIZER)

128 OUTPUT LEVELS

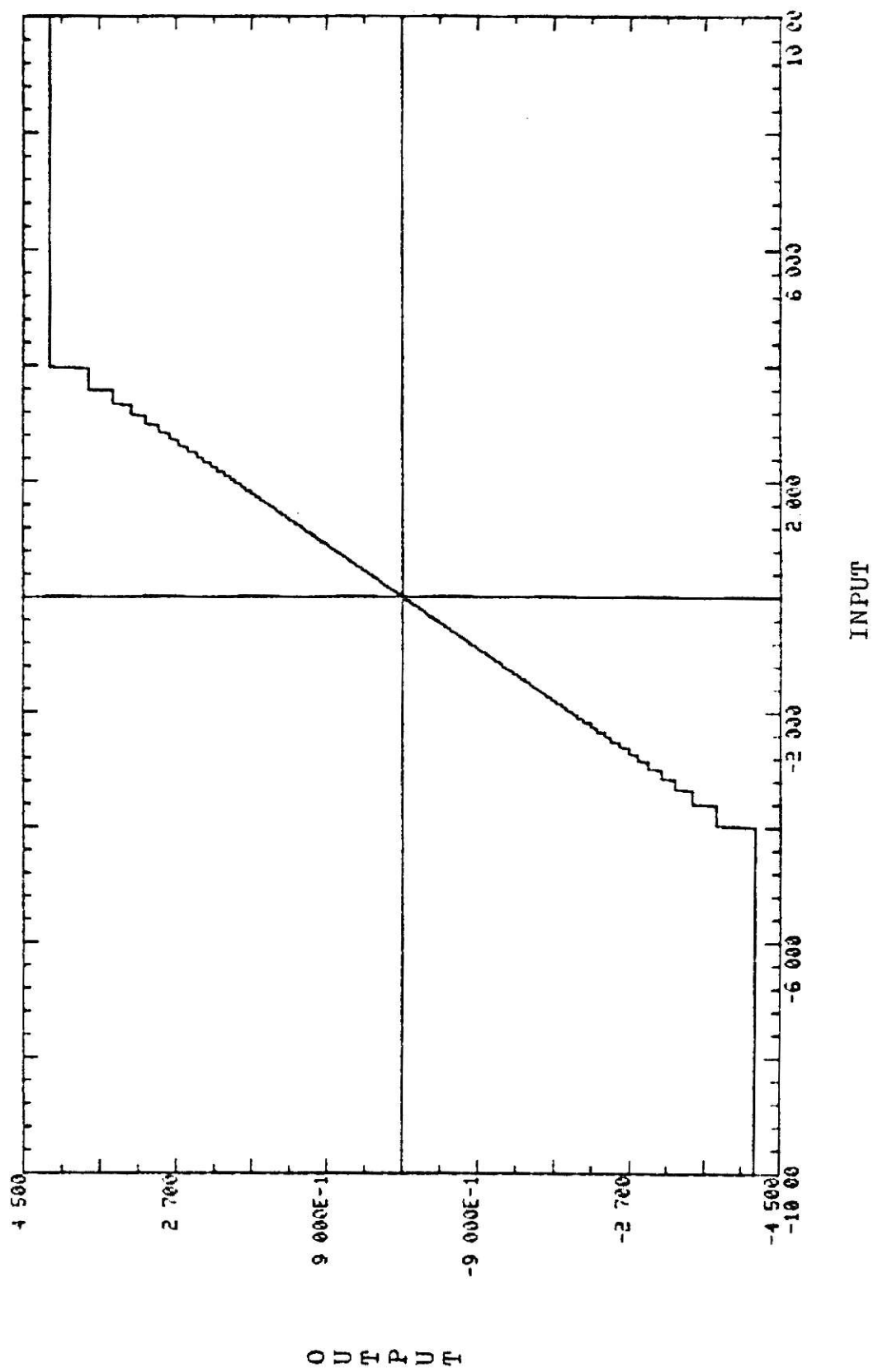


FIGURE 3.10

TRANSFER FUNCTION
(OPTIMUM GAUSSIAN QUANTIZER)

5 OUTPUT LEVELS

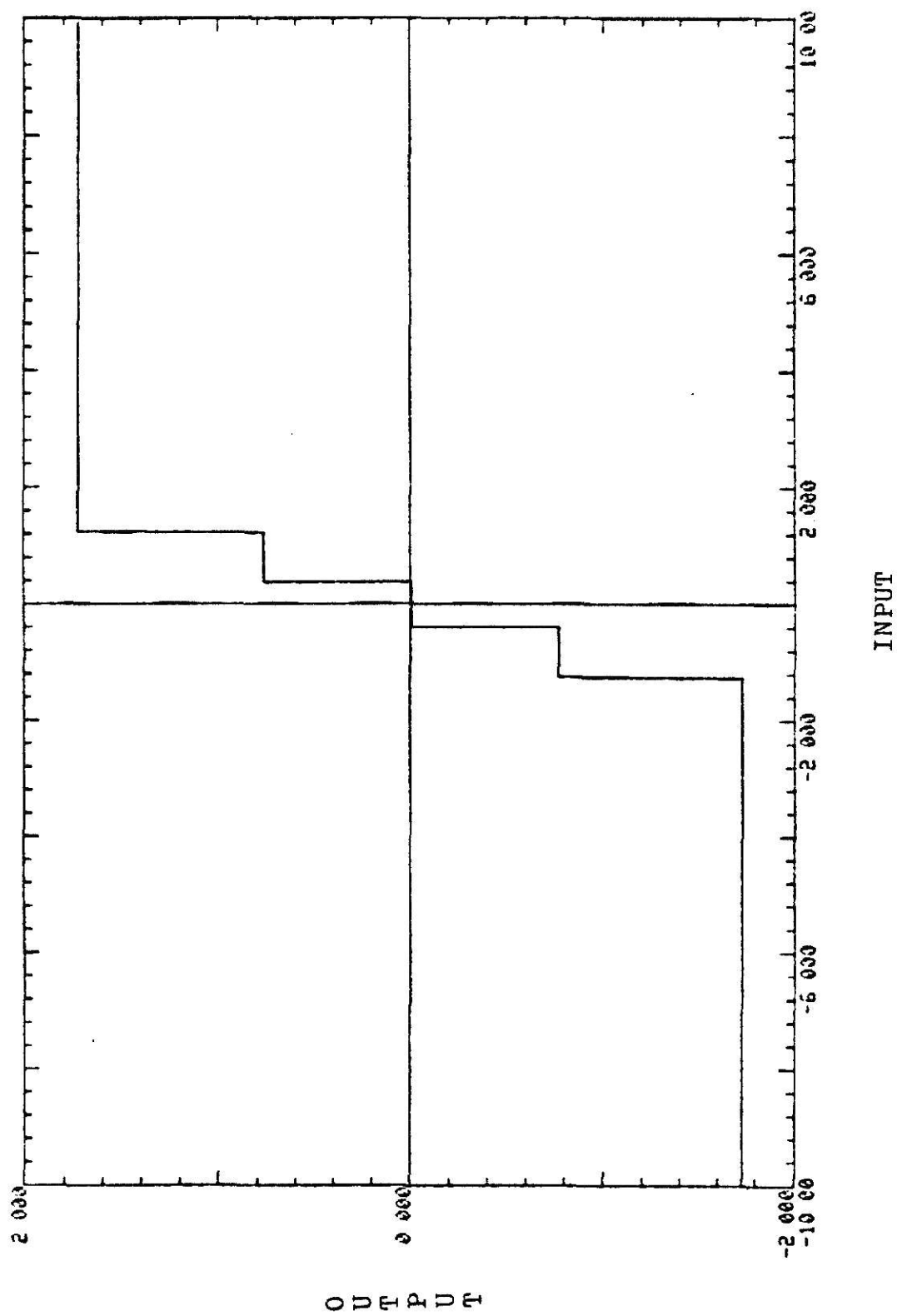


FIGURE 3.11

Entropy Considerations

The notion of entropy is useful in assessing the performance of the optimum quantizer. The entropy of the output levels is given by [6]

$$\text{ENTROPY} = \begin{cases} -2 \sum_{i=1}^M (\text{Probability of } y_i) \log_2 (\text{Probability of } y_i), & N \text{ EVEN} \\ -(2 \text{ Probability of } y_1) \log_2 (2 \text{ Probability of } y_1) \\ -2 \sum_{i=2}^M (\text{Probability of } y_i) \log_2 (\text{Probability of } y_i), & N \text{ ODD} \end{cases} \quad (3-10)$$

where entropy is defined as the average information (uncertainty) of the output levels. The smaller the uncertainty in the output, for a specified N , the better the performance. In (3-10) we note that

$$P(y_i) = \int_{x_i}^{x_{i+1}} p_X(x) dx \quad (3-11)$$

Values of the mean square error and entropy, obtained from (2-1) and (3-10) respectively, for the numbers of output levels 2 through 128 are tabulated in Table 3.3, and plotted in Figures 3.12 and 3.13. A comment about how these numerical values were obtained is now in order. The mean square error was calculated from (2-1)--i.e.,

$$E((X - Y)^2) = 2 \sum_{j=2}^{M+1} \int_{x_{j-1}}^{x_j} (x - y_{j-1})^2 p_X(x) dx$$

This integral can be expressed as the sum of three integrals since $(x - y_{j-1})^2$ can be written $y_{j-1}^2 - 2xy_{j-1} + x^2$. Thus

$$\begin{aligned} E((X - Y)^2) = & 2 \sum_{j=2}^{M+1} [y_{j-1}^2 \int_{x_{j-1}}^{x_j} p_X(x) dx - 2y_{j-1} \int_{x_{j-1}}^{x_j} x p_X(x) dx \\ & + \int_{x_{j-1}}^{x_j} x^2 p_X(x) dx] \end{aligned}$$

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POOR LEGIBILITY IN
THE ORIGINAL**

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COPY AVAILABLE**

TABLE 3.3
MEAN SQUARE ERROR AND ENTROPY

NUMBER OF OUTPUT LEVELS	MSE	ENTROPY	NUMBER OF OUTPUT LEVELS	MSE	ENTROPY
2	.363380	1.0000	65	.000823	5.7319
3	.190174	1.5358	66	.000604	5.7533
4	.117481	1.9111	67	.000587	5.7750
5	.079941	2.2029	68	.000571	5.7961
6	.057977	2.4428	69	.000554	5.8169
7	.044000	2.6469	70	.000539	5.8374
8	.034547	2.8249	71	.000524	5.8574
9	.027853	2.9827	72	.000510	5.8775
10	.022936	3.1246	73	.000496	5.8971
11	.019219	3.2533	74	.000483	5.9166
12	.016339	3.3717	75	.000470	5.9357
13	.014062	3.4807	76	.000458	5.9548
14	.012231	3.5820	77	.000446	5.9731
15	.010736	3.6766	78	.000434	5.9915
16	.009500	3.7653	79	.000424	6.0097
17	.008466	3.8488	80	.000413	6.0276
18	.007592	3.9277	81	.000403	6.0453
19	.006847	4.0025	82	.000393	6.0628
20	.006207	4.0736	83	.000383	6.0800
21	.005653	4.1411	84	.000374	6.0971
22	.005170	4.2059	85	.000366	6.1140
23	.004745	4.2673	86	.000357	6.1298
24	.004371	4.3271	87	.000349	6.1472
25	.004040	4.3840	88	.000341	6.1630
26	.003745	4.4388	89	.000333	6.1797
27	.003481	4.4916	90	.000327	6.1958
28	.003245	4.5425	91	.000320	6.2114
29	.003032	4.5916	92	.000313	6.2270
30	.002839	4.6392	93	.000307	6.2424
31	.002663	4.6852	94	.000300	6.2577
32	.002504	4.7297	95	.000294	6.2728
33	.002358	4.7730	96	.000288	6.2877
34	.002225	4.8150	97	.000282	6.3025
35	.002102	4.8557	98	.000276	6.3172
36	.001989	4.8954	99	.000271	6.3317
37	.001886	4.9340	100	.000265	6.3460
38	.001791	4.9714	101	.000260	6.3602
39	.001703	5.0082	102	.000255	6.3745
40	.001620	5.0440	103	.000250	6.3886
41	.001544	5.0788	104	.000245	6.4020
42	.001473	5.1128	105	.000240	6.4157
43	.001406	5.1461	106	.000236	6.4291
44	.001344	5.1786	107	.000231	6.4425
45	.001286	5.2104	108	.000227	6.4559
46	.001232	5.2414	109	.000223	6.4691
47	.001180	5.2719	110	.000218	6.4821
48	.001132	5.3017	111	.000214	6.4950
49	.001087	5.3309	112	.000210	6.5078
50	.001045	5.3595	113	.000207	6.5205
51	.001005	5.3876	114	.000203	6.5331
52	.000968	5.4151	115	.000200	6.5455
53	.000933	5.4421	116	.000198	6.5581
54	.000899	5.4686	117	.000194	6.5704
55	.000867	5.4947	118	.000191	6.5826
56	.000837	5.5202	119	.000188	6.5946
57	.000808	5.5453	120	.000184	6.6066
58	.000781	5.5700	121	.000181	6.6185
59	.000755	5.5943	122	.000178	6.6302
60	.000730	5.6182	123	.000175	6.6418
61	.000706	5.6416	124	.000173	6.6535
62	.000684	5.6646	125	.000170	6.6649
63	.000663	5.6875	126	.000167	6.6764
64	.000642	5.7099	127	.000165	6.6877
			128	.000163	6.6989

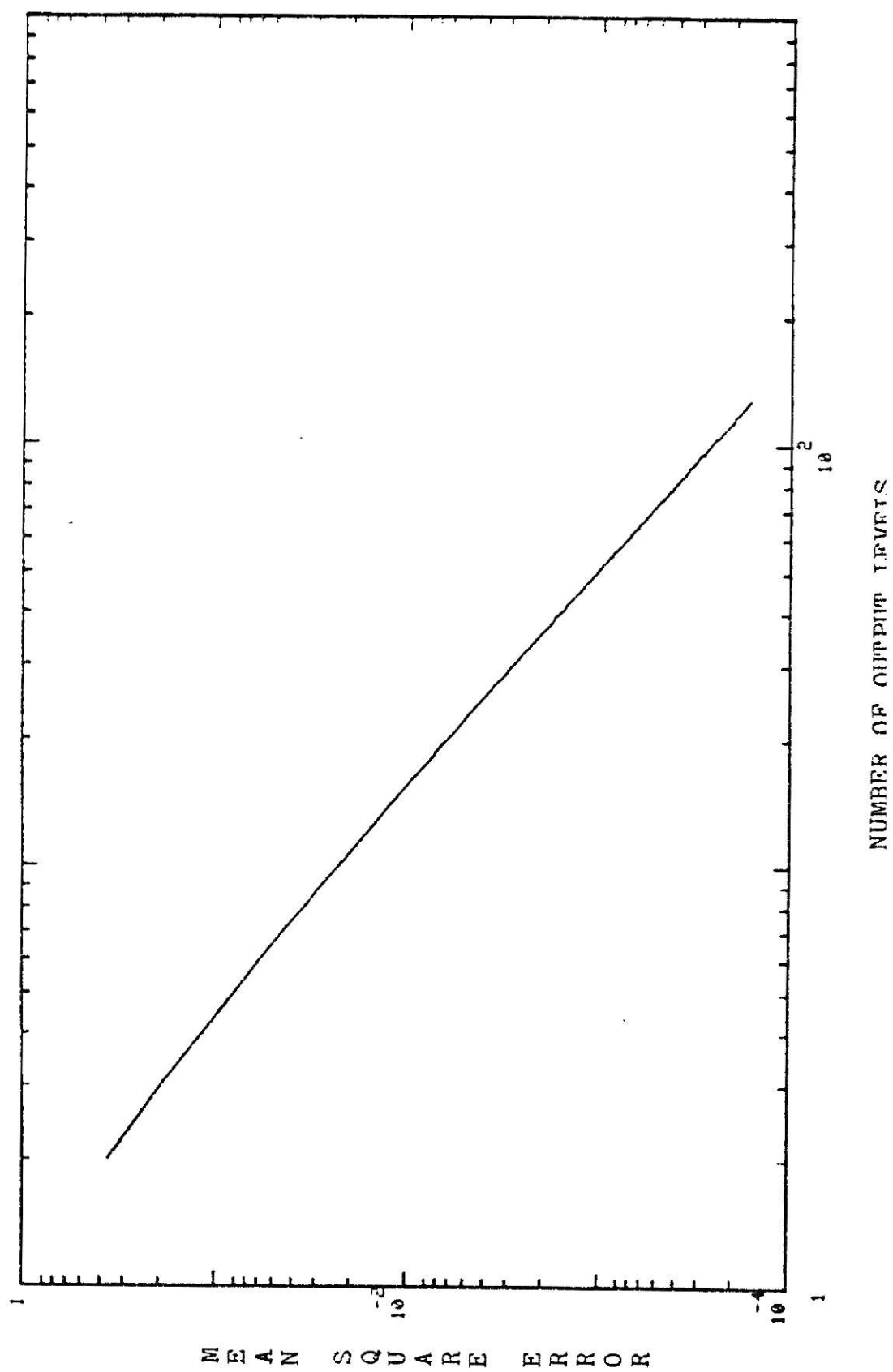


FIGURE 3.12

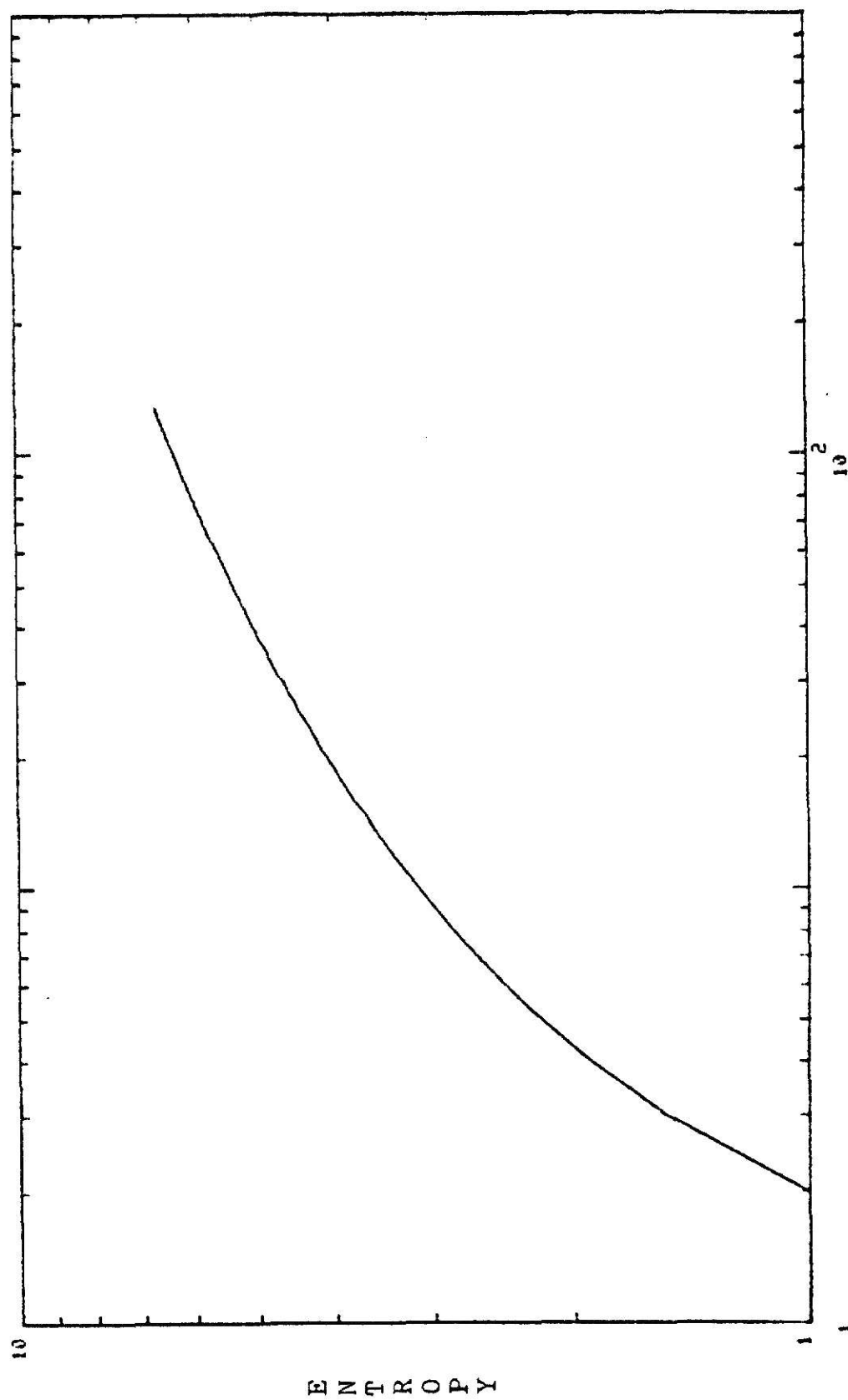


FIGURE 3.13

The forms of the first two integrals were discussed earlier in (3-3).

The last integral can be integrated by parts as follows:

$$\int u \frac{dv}{dx} dx = uv - \int v \frac{du}{dx} dx \quad (3-13)$$

where $u=x$ and $\frac{dv}{dx} = xp_X(x)$. This yields

$$\int_{x_{j-1}}^{x_j} x^2 p_X(x) dx = -xp_X(x) + \int_{x_{j-1}}^{x_j} p_X(x) dx \quad (3-14)$$

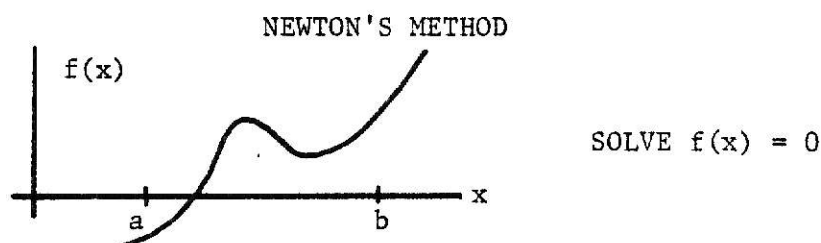
Note that (3-14) is valid only when $p_X(x)$ is gaussian, and that the last integral is of the form previously discussed in (3-3).

The tabulated values for mean square error and entropy in Table 3.3 compare very closely with those published by Max [1].

APPENDIX 3.1

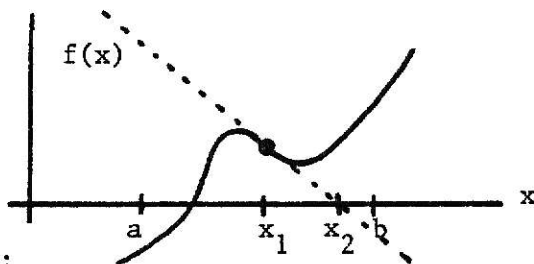
This appendix contains the iterative procedure to determine the zeroes of functions known as Newton's method.

Note if the initial value is not sufficiently close to the actual zero this method will not converge. If no prior knowledge is known about the location of the zero(es) it is suggested that another method that always converges (but is very slow) such as the bisection method be used for a few iterations to insure that the initial guess is sufficiently close to the actual zero. Newton's method may then be used for subsequent iterations to speed the process up.



- (1) Pick any x_1 in interval $\epsilon[a,b]$.
- (2) Draw tangent to graph at $f(x_1)$.
- (3) Where tangent crosses horizontal axis is x_2 .
- (4) Draw tangent to Graph at $f(x_2)$ etc.

** If tangent to graph lies outside the interval $[a,b]$ this method fails.



Derivation:

$$y - f(x_n) = [f^{(1)}(x_n)](x - x_n)$$

Next x is where $y = 0$

Thus

$$x_{n+1} = x_n - \frac{f(x_n)}{f^{(1)}(x_n)}$$

Iterative Procedure to solve (2-8)

$$\int_{x_j}^{x_{j+1}} (x - y_j) p_X(x) dx = 0$$

where x_j, y_j are known solve for x_{j+1}

* Use Newton's method for finding zeroes of functions

$$g(x) = 0 = \int_{x_j}^{x_{j+1}} (x - y_j) p_X(x) dx = 0$$

$$\hat{x}_{j+1} = x_{j+1} - \frac{g(x_{j+1})}{g^{(1)}(x_{j+1})}$$

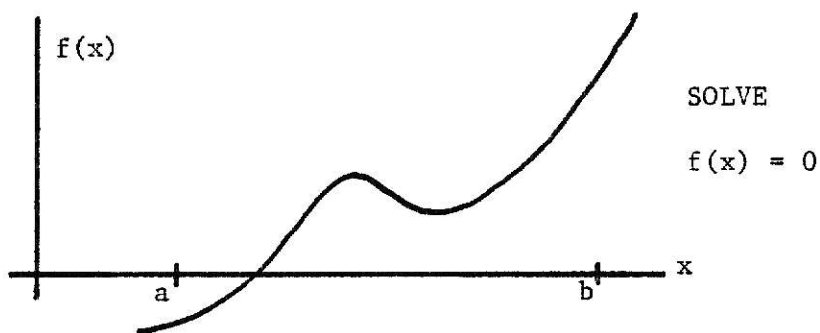
$$\hat{x}_{j+1} = x_{j+1} - \frac{\int_{x_j}^{x_{j+1}} (x - y_j) p_X(x) dx}{(x_{j+1} - y_j) p_X(x_{j+1})}$$

where $(x_{j+1})_{\text{initial}} = 2y_j - x_j$

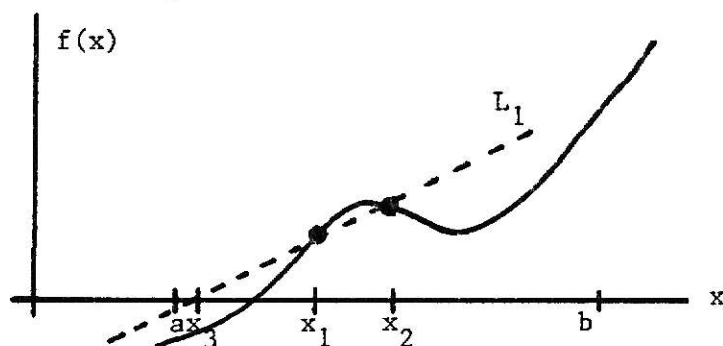
APPENDIX 3.2

This appendix contains the iterative procedure to determine the zeroes of functions known as the secant method.

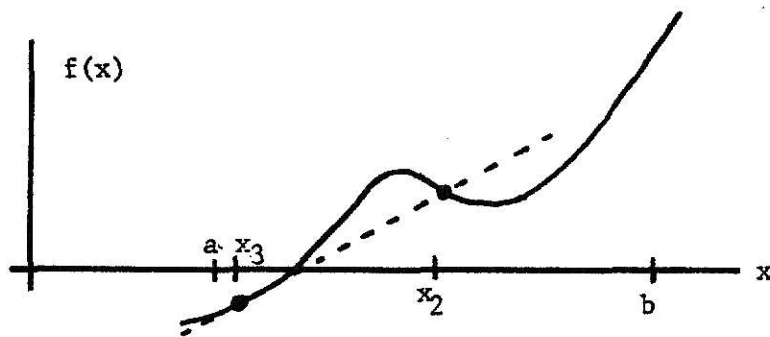
Note if the initial value is not close to the zero this method will not converge. If no prior knowledge is known about the location of the zero(es) it is suggested that another method that always converges (but very slow) such as the bisection method be used for a few iterations to insure that the initial guess is sufficiently close to the actual zero. The secant method may be used for subsequent iterations to speed the process up.



- (1) Pick any two points in the interval $\epsilon[a,b]$, ie, x_1 and x_2
- (2) Connect $(x_1, f(x_1))$ and $(x_2, f(x_2))$ by a straight line L_1



- (3) Let x_3 be the intersection of L_1 and the interval $[a,b]$ (ie, the horizontal axis)
- (4) Connect $(x_2, f(x_2))$ and $(x_3, f(x_3))$ etc.



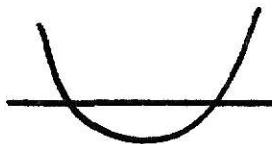
Caution: Secant line may not cross horizontal axis within the interval $[a, b]$

** The sequence $\{x_n\}$ will converge to a zero of $f(x)$ if x_1 and x_2 are close enough to a zero and if the zero is a simple zero

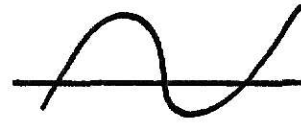


Not a simple zero

(slope = 0)



Simple zero



Simple zero

Derivation:

$$y - f(x_1) = \frac{f(x_2) - f(x_1)}{x_2 - x_1} (x - x_1)$$

Next x is where $y = 0$ rearranging,

$$x = x_1 - \frac{f(x_1) [x_1 - x_2]}{f(x_2) - f(x_1)}$$

APPENDIX 3.3

This appendix contains the program to generate the parameters X_i and Y_i , which completely describe the optimum gaussian quantizer for a finite number of output levels, N .

THE OPTIMUM QUANTIZER

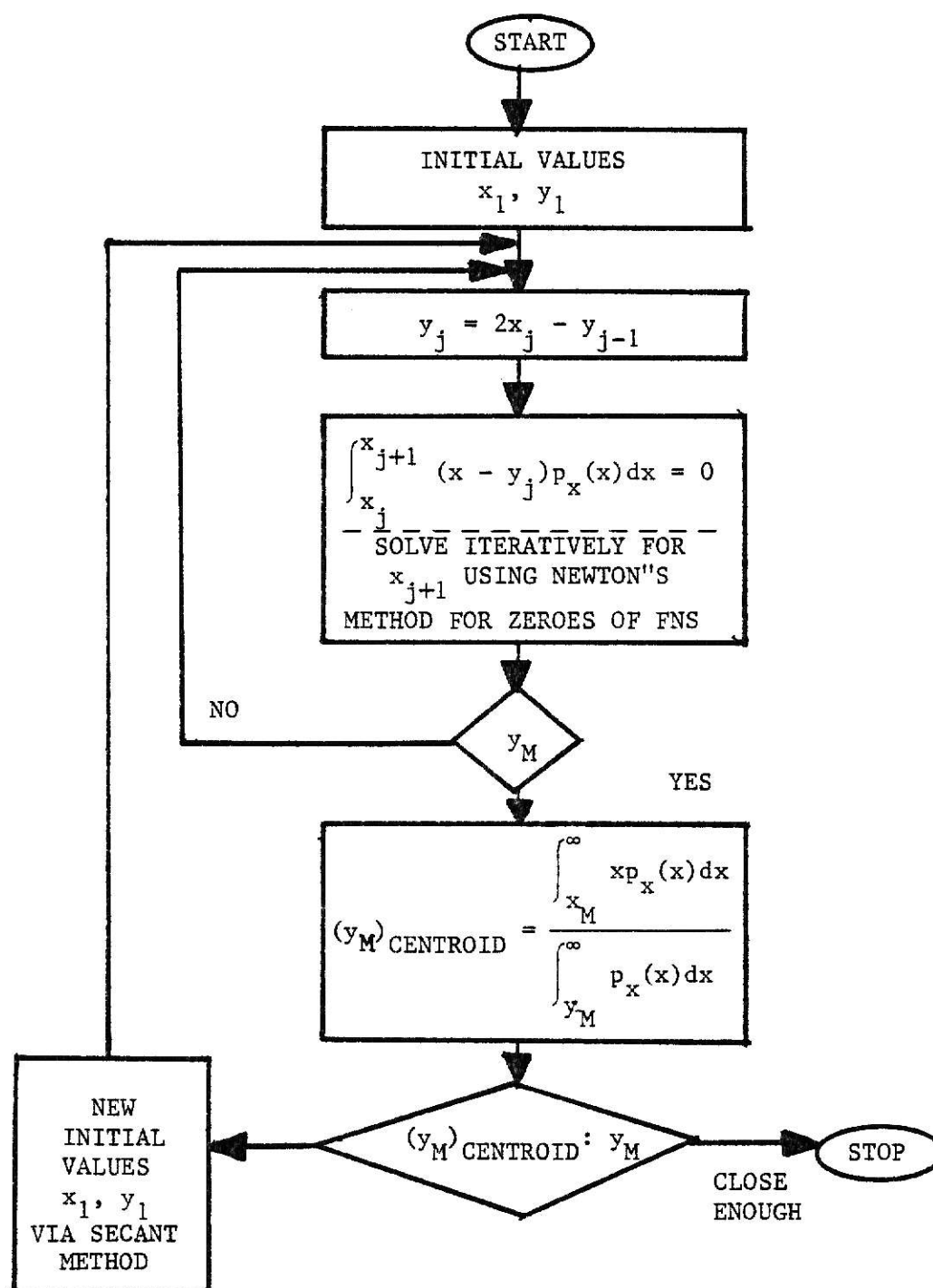


FIGURE 3.14

TABLGENS.FR

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Page 1

```

C      THIS ROUTINE COMPUTES THE CUT-POINTS AND REP-VALUES FOR
C      A GAUSSIAN PROBABILITY DENSITY AND MSE DISTORTION MEASURE.
C
C      COMPILER FREE
C      DOUBLE PRECISION X,Y,MSE,ENTROPY,AREA1,AREA2
C      DOUBLE PRECISION X0,X1,X2,P0,P2,XN,YN
C      DOUBLE PRECISION P1,Y0,GAP,Z0,Z1,W0,W1,AREA,TAIL,A0,A1,W2,DABS,Q1,Q0
C      REAL MSE1,ENTR1
C      DIMENSION TABLG(128,2)
C      CALL OPENW(4,'MEAN SQUARED ERROR OUTPUT FILENAME ? ',0,SIZE)
C      CALL OPENW(5,'ENTROPY OUTPUT FILENAME ? ',0,SIZE)
C      ITTO=10
C      ITTI=11
C      ITTIV=12
C
C      GET THE NUMBER OF LEVELS
C
C      WRITE(ITTO,5)
C      FORMAT(' ')
C      5000 ACCEPT 'NUMBER OF OUTPUT LEVELS ? ', NLEV
C      N=NLEV/2
C
C      ESTIMATE THE VALUE OF THE FIRST CUT-POINT OR REP-VALUE
C
C      P1=2.1613D0/DBLE(FLCAT(NLEV))-1.653D0/DELE(FLOAT(NLEV)**2.0)
C
C      PERFORM TEN ITERATIONS
C
C      WRITE(ITTV,6000) NLEV
C      6000 FORMAT('I',12X,I3,' OUTPUT LEVELS')
C      DO 1073 II=1,10
C      WRITE(ITTV,6500) NLEV
C      6500 FORMAT('J',50X,'N=',I3)
C      WRITE(ITTV,2040) II
C      2040 FORMAT(' ',12X,'ITERATION NUMBER',I3,Z)
C      WRITE(ITTV,2045)
C      2045 FORMAT('I',4X,'J',7X,'X(J)',12X,'Y(J)')
C
C      INITIALIZE THE FIRST CUT-POINT AND REP-VALUE
C
C      IF(N*2.EQ.NLEV) GO TO 1100
C
C      ODD NUMBER OF LEVELS
C
C      X0=P1
C      Y0=2.0D0*P1
C      GO TO 1110
C      1100 CONTINUE
C
C      EVEN NUMBER OF LEVELS
C
C      X0=0.0D0
C      Y0=P1
C      1110 CONTINUE
C      I=1
C
C      THE FIRST CUT-POINT VALUE

```

TABLGEN5.FR 7/15/1982 01:132 DIR DFO Page 2

```

C
2000 WRITE(ITTYY,2000) I,X0,Y0
      FORMAT(' ',I35,F13.6,F16.6)
      TABLEG(I,1)=X0
      TABLEG(I,2)=Y0
      IF(NLEV.LT.4) GO TO 1000

C
C      DETERMINE THE REST OF THE CUT-POINTS AND REP-VALUES
C
      DO 1000 I=2,N
      CALL CUTPT(X0,Y0,X1)

C
      X0=X1
      Y0=2.0D0*X0-Y0
      WRITE(ITTYY,2000) I,X0,Y0
      TABLEG(I,1)=X0
      TABLEG(I,2)=Y0
1000  CONTINUE

C
C      USING AN ALTERNATE METHOD,
C      DETERMINE THE LAST REP-VALUE, Y(N), BY COMPUTING
C      THE CENTROID OF THE AREA FROM X(N) TO INFINITY
C
      XN=X0
      CALL ERF(XN,AREA)
      YN=DEXP(-XN**2/2.0D0)/DSQRT(2.0D0*3.141592D0)/(0.5D0-AREA)
      WRITE(ITTYY,3700) YN
3700  FORMAT(' ',17X,'SEDANT METHOD.....',6X,'Y(N) =',1X,F16.6)

C
C      COMPUTE THE DIFFERENCE BETWEEN THE TWO VALUES AND UPDATE
C      THE STARTING VALUE, P1
C
      Q1=Y0-YN
      IF(DABS(Q1/YN).LT.1.0D-6) GO TO 3000
      IF(II.EQ.1) P2=0.99104P1
      IF(II.NE.1) P2=(P1*Q0-P0*Q1)/(Q0-Q1)
      P0=P1
      P1=P2
      Q0=Q1
1073  CONTINUE

C
C      NOW THAT THE OPTIMUM PARAMETERS HAVE BEEN
C      DETERMINED, CALCULATE THE MEAN SQUARED
C      ERROR FOR THE CURRENT NUMBER OF
C      OUTPUT LEVELS
C
3000  MSE=0.0D0
      IF (NLEV.EQ.2) GO TO 6500
      IF (N*2.EQ.NLEV) GO TO 7000

C
C      ODD NUMBER OF LEVELS
C
      X=TABLEG(1,1)
      Y=0.0D0
      CALL ERF(X,AREA)
      MSE=2.0D0*(((X+2.0D0*Y)*DEXP(-X**2/2.0D0)
      & -2.0D0*Y)/DSQRT(2.0D0*3.141592D0)

```

```

TABLENS,FR          7/15/1982  01:132  DIR DPO          Page 3

      1  +(1.000+Y**2)*AREA)
      IF (NLEV.EQ.3) GO TO 8500
C
C      ODD OR EVEN NUMBER OF LEVELS
C
7000  DO 8000 I=1,N-1
      X1=TABLEG(I,1)
      X2=TABLEG(I+1,1)
      Y=TABLEG(I,2)
      CALL ERF(X1,AREA1)
      CALL ERF(X2,AREA2)
      AREA=AREA2-AREA1
      MSE=2.000*((-X2+2.000*Y)*DEXP(-X2**2/2.000)
      1  -(-X1+2.000*Y)*DEXP(-X1**2/2.000))
      1  /DSQRT(2.000*3.14159200)
      1  +(1.000+Y**2)*AREA)+MSE
8000  CONTINUE
C
C      INTEGRATE FROM X(N) TO INFINITY
C
8500  X=TABLEG(N,1)
      Y=TABLEG(N,2)
      CALL ERF(X,AREA)
      MSE=2.000*(-(-X+2.000*Y)*DEXP(-X**2/2.000)
      1  /DSQRT(2.000*3.14159200)
      1  +(1.000+Y**2)*(0.500-AREA))+MSE
      MSE1=SNGL(MSE)
      WRITE (ITTY,1) NLEV,MSE
1      FORMAT(' ',12X,'MEAN SQUARED ERROR FOR ',I3,' OUTPUT LEVELS = ',F9.6)
      WRITE BINARY(4) MSE1
C
C      CALCULATE THE ENTROPY
C
      ENTROPY=0.000
      IF(NLEV.EQ.2) GO TO 9300
      IF (N*2.EQ.NLEV) GO TO 9000
C
C      ODD NUMBER OF LEVELS
C
      X=TABLEG(1,1)
      CALL ERF(X,AREA)
      ENTROPY=(2.000*AREA*DLOG(2.000*AREA))/DLOG(2.000)
      IF (NLEV.EQ.3) GO TO 9300
C
C      ODD OR EVEN NUMBER OF LEVELS
C
9000  DO 9500 I=1,N-1
      X1=TABLEG(I,1)
      X2=TABLEG(I+1,1)
      CALL ERF(X1,AREA1)
      CALL ERF(X2,AREA2)
      AREA=AREA2-AREA1
      ENTROPY=2.000*(AREA*DLOG(AREA))/DLOG(2.000)+ENTROPY
9500  CONTINUE
C
C      INTEGRATE FROM X(N) TO INFINITY
C
9300  X=TABLEG(N,1)
      CALL ERF(X,AREA)
      AREA=0.500-AREA
      ENTROPY=-((2.000*(AREA*DLOG(AREA))/DLOG(2.000)+ENTROPY)
      1  ENTR1=SNGL(ENTROPY)
      WRITE (ITTY,2) NLEV,ENTROPY
2      FORMAT(' ',12X,'ENTROPY FOR ',I3X,I3,' OUTPUT LEVELS = ',F9.6)
      WRITE BINARY(5) ENTR1
      IF (NLEV.GT.0) GO TO 5000
      STOP
      END

```

CUTPT2.FR 7/15/1982 0: 0:22 DIR DPO Page 1

```

C      THIS SUBROUTINE COMPUTES THE NEXT CUT-POINT, X1, GIVEN THE
C      PREVIOUS CUT-POINT AND REP-VALUE
C
      SUBROUTINE CUTPT(X0,Y0,X1)
      DOUBLE PRECISION X0,Y0,X1,EPS,AREA,AREA0,AREA1,DELX1,PROB,FUNC,DABS
      DOUBLE PRECISION TAIL
      EPS=1.0D-9
C
C      INTEGRATE THE FUNCTION FROM X0 TO Y0
C
      CALL FFX(X0,Y0,AREA0)
C
      X1=2.0D0*Y0-X0
C
C      ITERATE UP TO 100 TIMES TO DETERMINE X1
C
      DO 1000 I=1,100
      CALL FFX(X1,Y0,AREA1)
      AREA=AREA1-AREA0
      DELX1=-(AREA)/(PROB(X1)*FUNC(X1,Y0))
      IF(DABS(DELX1/X1).LT.EPS) RETURN
      X1=X1+DELX1
1000  CONTINUE
      RETURN
      END

```

ERF2.FR 7/15/1982 0: 0:21 DIR DPO Page 1

```

C      THIS SUBROUTINE EVALUATES THE INTEGRAL OF THE
C      GAUSSIAN PROBABILITY DISTRIBUTION VIA AN
C      INFINITE SERIES REPRESENTATION
C
      SUBROUTINE ERF(X,AREA)
      DOUBLE PRECISION X,W,AREA,AREA1,DEXP,DSQRT,DBLE,DABS
      W=1.0D0
      AREA1=(DEXP(-X**2/2.0D0))/DSQRT(2.0D0*3.141592D0)
      DO 100 J=1,100
      K=2*J+1
      W=W*DBLE(FLOAT(K))
      AREA=(X/DSQRT(2.0D0))*K**2.0D0*W/DEXP(-X**2/2.0D0)/DSQRT(3.141592D0)+AREA1
      IF (DABS(AREA1-AREA).LT.1.0D-9) RETURN
      AREA1=AREA
100  CONTINUE
      TYPE 'NEED MORE TERMS OF THE INFINITE SERIES'
      RETURN
      END

```

FPX.FR 7/15/1982 0: 0:25 DIR DPO Page 1

```

      SUBROUTINE FPX(X,Y,AREA)
      DOUBLE PRECISION X,Y,AREA,DEXP,DSQRT,A
      CALL ERF(X,A)
      AREA=(1-DEXP(-X**2/2.0D0))/DSQRT(2.0D0*3.141592D0)-Y**A
      RETURN
      END

```

PROB.FR 7/15/1982 01:01:24 DIR DFO Page 1

```
DOUBLE PRECISION FUNCTION PROB(X)
DOUBLE PRECISION X,DEXP,DSQRT
PROB=DEXP(-X**2/2.0D0)/DSQRT(2.0D0*3.141592D0)
RETURN
END
```

FUNC.FR 7/15/1982 01:01:24 DIR DFO Page 1

```
DOUBLE PRECISION FUNCTION FUNC(X,Y0)
DOUBLE PRECISION X,Y0
FUNC=X-Y0
RETURN
END
```

CHAPTER IV

EXPERIMENTAL RESULTS

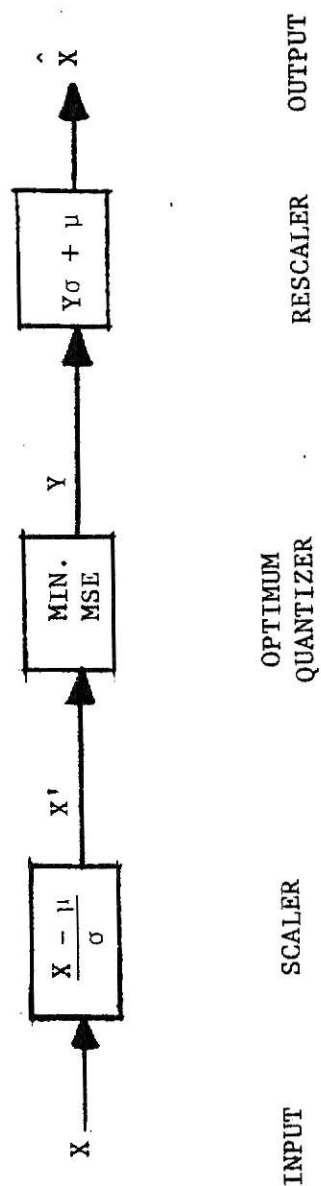
The experimental results reported in this chapter were carried out at the Signal Processing Laboratory of the Electrical Engineering Department at KSU.

A random gaussian input sequence of specified mean and variance was generated using the program NORMGEN, listed in Appendix 4.1. The mean and variance of this input sequence were then calculated via the program, XBAR. These actual values are necessary since the input sequence is scaled to zero mean and unit variance, as depicted in Figure 4.1. The scaled sequence is then processed through the quantizer. The quantizer's output is also gaussianly distributed with zero mean and unit variance. Hence it must be rescaled to the actual mean and variance of the input sequence. The mean square error was then calculated to access whether the rescaled output sequence, for the given number of output levels N, is a faithful reproduction of the input sequence. The mean square error can be calculated at two places in Figure 4.1; at X' , Y and at X , $\hat{X}$. Both results are useful. The first, at X' , Y, has been discussed earlier in (2-1). It provides an experimental measure for comparison against the theoretical values tabulated in Table 3.3. The second, at X , $\hat{X}$, establishes how faithful a reproduction the rescaled quantizer output sequence is of the input sequence. Theses two mean square errors are related. Using the notation of Figure 4.1 we have

$$x = x'\sigma + \mu$$

$$\hat{x} = y\sigma + \mu$$

$$\frac{(x - \hat{x})^2}{NPTS} = \frac{((x'\sigma + \mu) - (y\sigma + \mu))^2}{NPTS} \quad (4-1)$$



X : Input random variable
 μ : Mean of X
 σ : Standard deviation of X
 $\hat{X}$: Reconstruction of X ; i.e., quantizer output

THE OPTIMUM QUANTIZATION PROCESS

FIGURE 4.1

where NPTS is the total number of data points. Rewriting (4-1) produces the desired relationship

$$(x - \hat{x})^2 = \sigma^2(x' - y)^2 \quad (4-2)$$

The entire optimum quantization process of Figure 4.1 is implemented by the program QUANTIZE1, listed in Appendix 4.2.

Two examples of the entire quantization process are now presented. The first example has gaussian input sequence with mean 0.044654 and variance 0.93953. The second example has a gaussian input sequence with mean 250.11 and variance 5.6372. Both examples follow the same format. The input sequence and the rescaled quantizer output sequences are shown for 2,4,8,16,32,64, and 128 output levels, which correspond to 1,2,3,4,5,6, and 7 output bits respectively. Perhaps more informative of what is actually happening are the histograms of the rescaled quantizer outputs, each of which consist of 2048 data points. The histograms were formed by dividing the range between the minimum and maximum values of each sequence into 450 bins, and then determining how many of the 2048 data points fall in each of these 450 bins. The program HISTPLOT was used to compute these histograms. We observe that the histograms resemble gaussian distributions as the number of output levels is increased. Finally the mean square error is calculated for each of the output sequences relative to the input sequence. These values are compared to the theoretical values published by Max [1], except N=64 and N=128 which Max did not calculate. These two exceptions were earlier tabulated in Table 3.3. The mean square error is then rescaled, due to its dependence on the variance of the input sequence, via the relationship obtained in (4-2).

EXAMPLE 1

A random gaussian input sequence with a mean and variance equal to 0.044654 and 0.93953, respectively. Note that the input sequence is in floating point arithmetic, while the output sequences are in fixed point arithmetic.

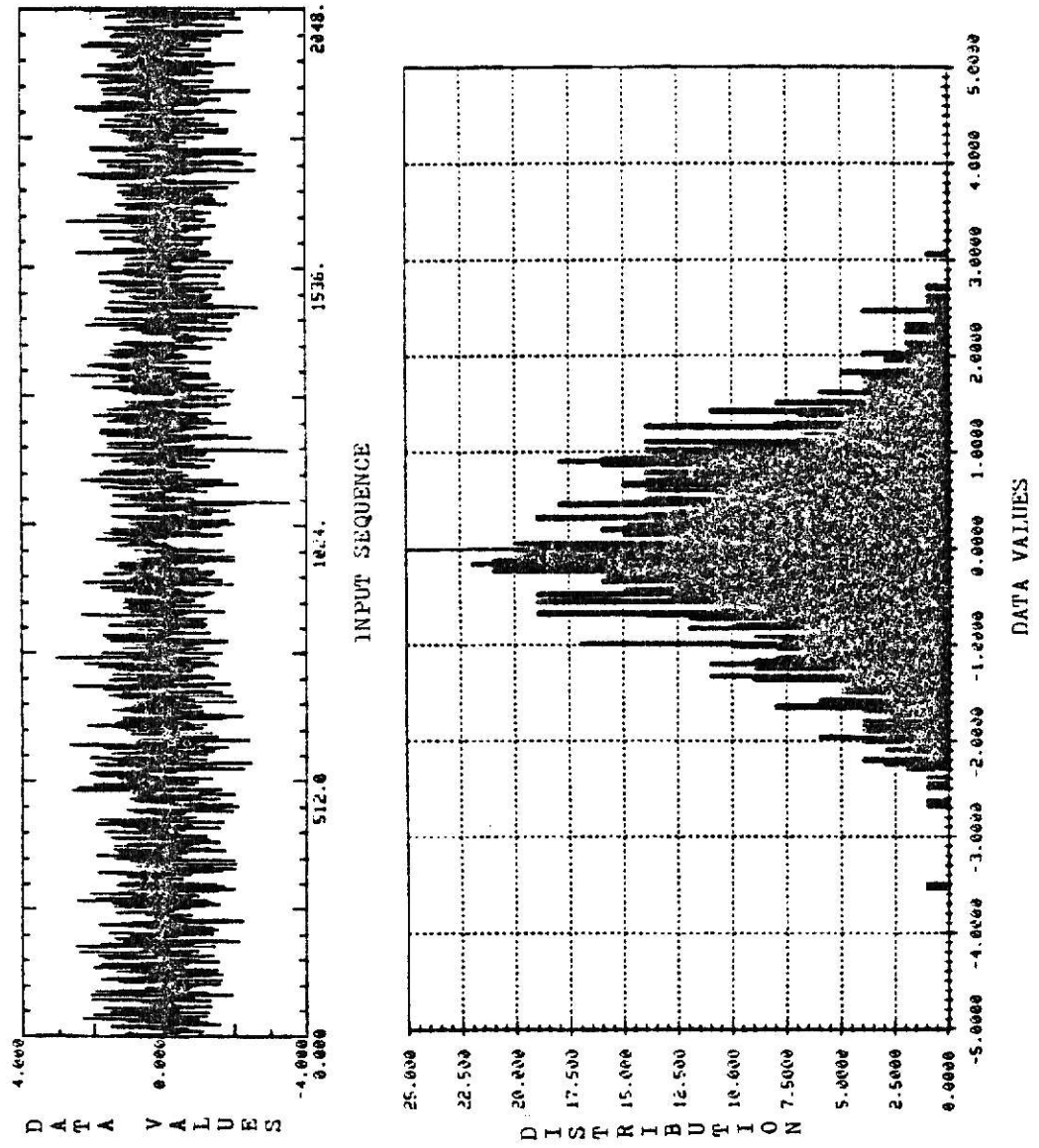
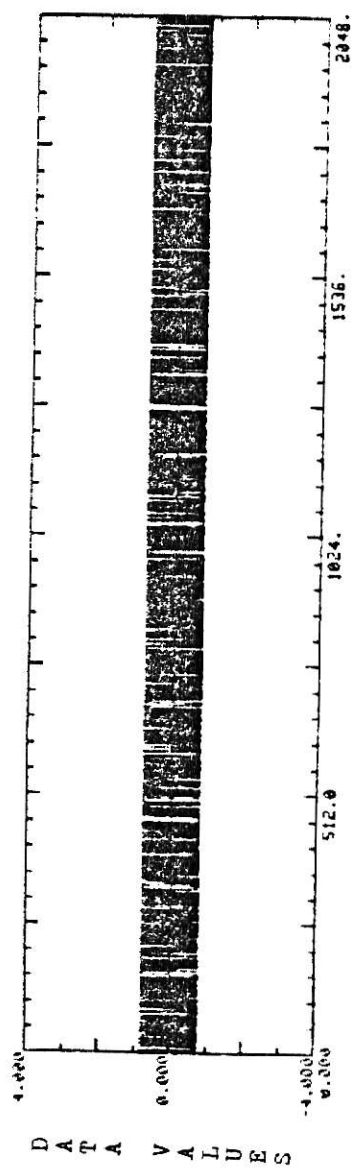


FIGURE 4.2



2 OUTPUT LEVELS

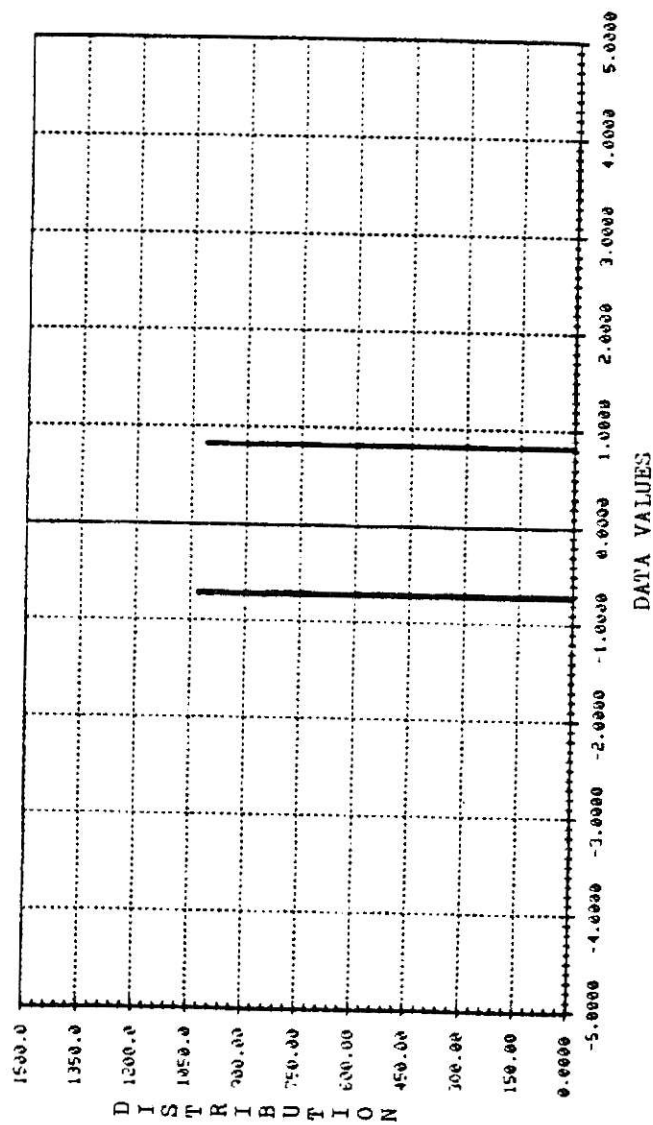
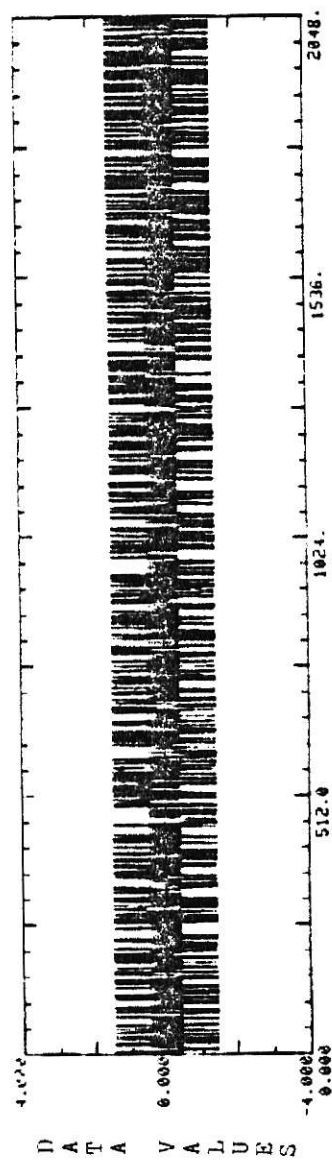


FIGURE 4.3



4 OUTPUT LEVELS

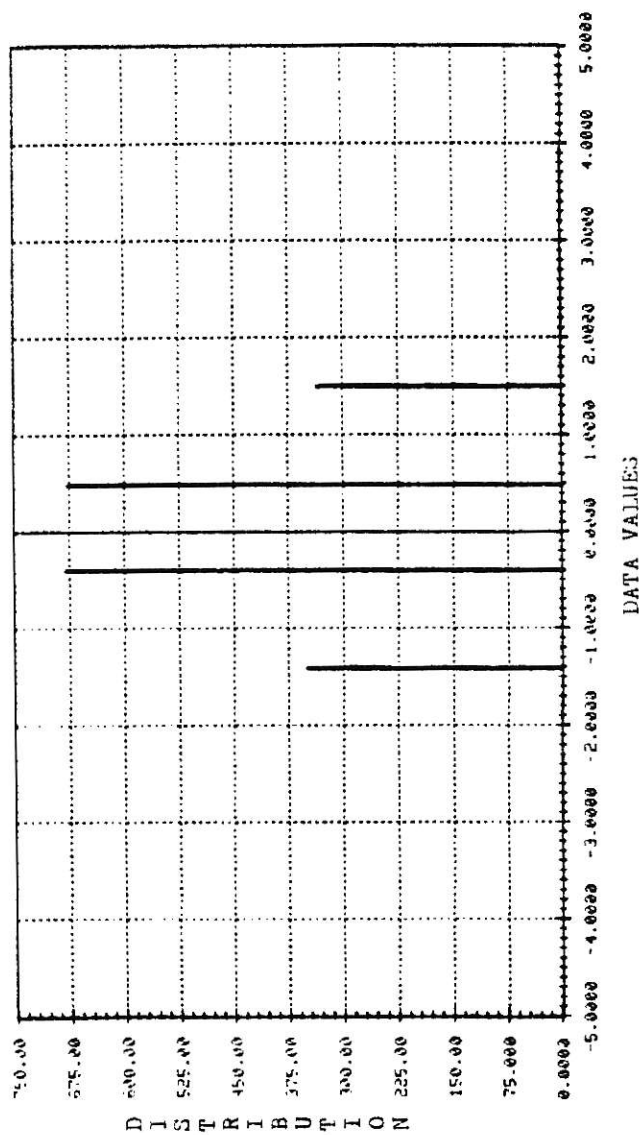
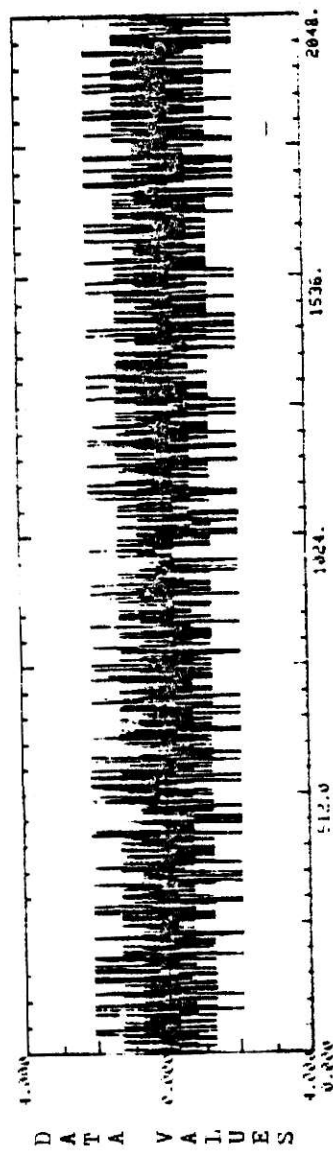


FIGURE 4.4



8 OUTPUT LEVELS

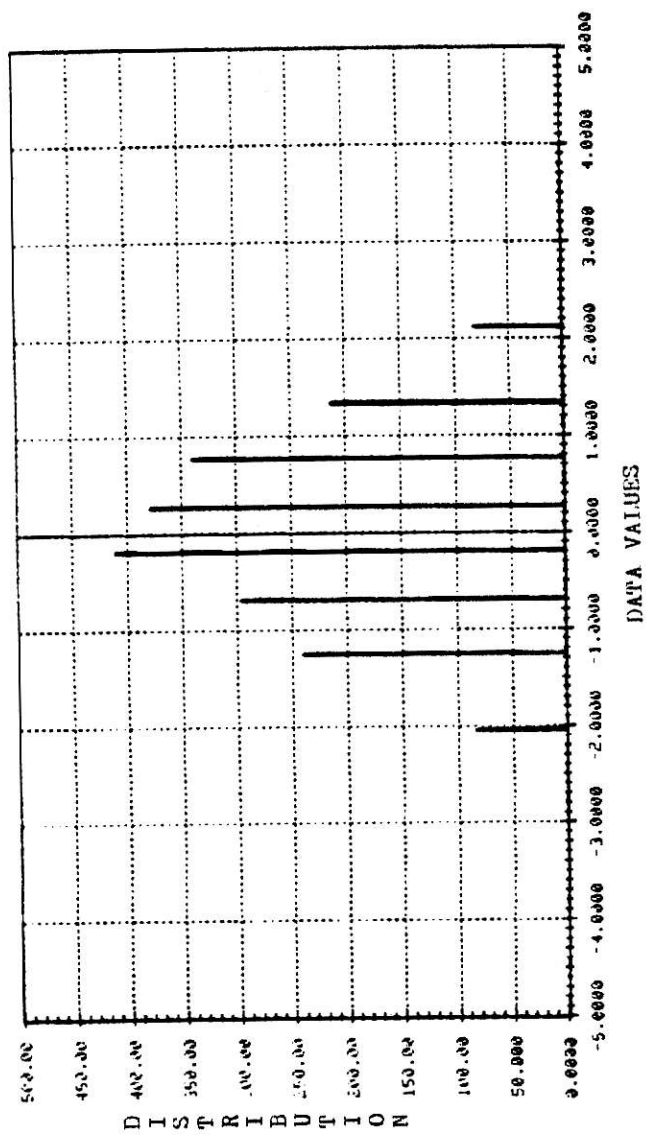
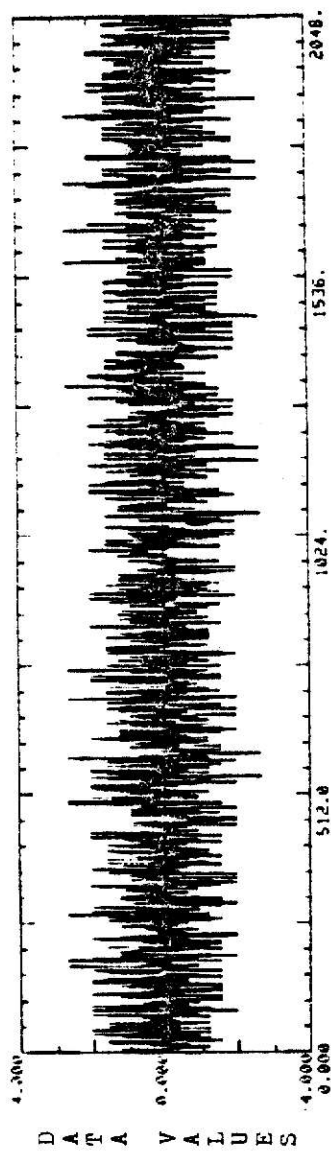


FIGURE 4.5



16 OUTPUT LEVELS

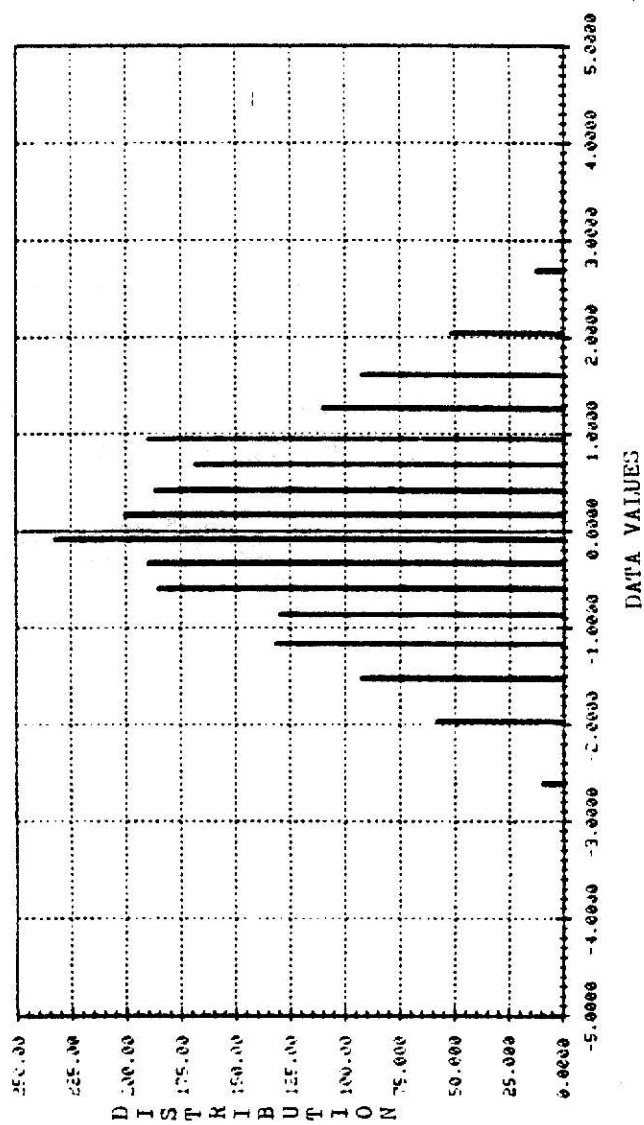


FIGURE 4.6

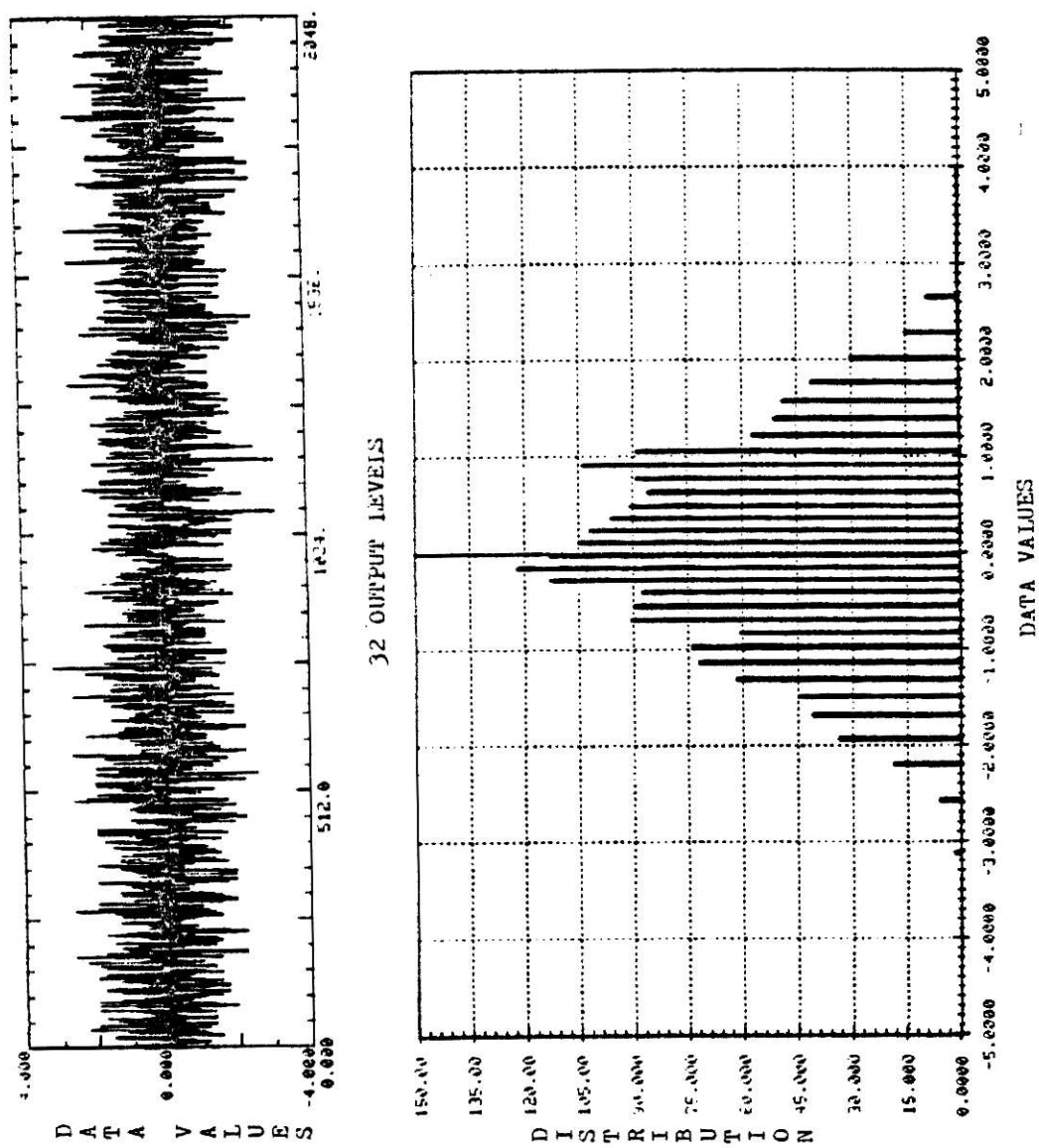
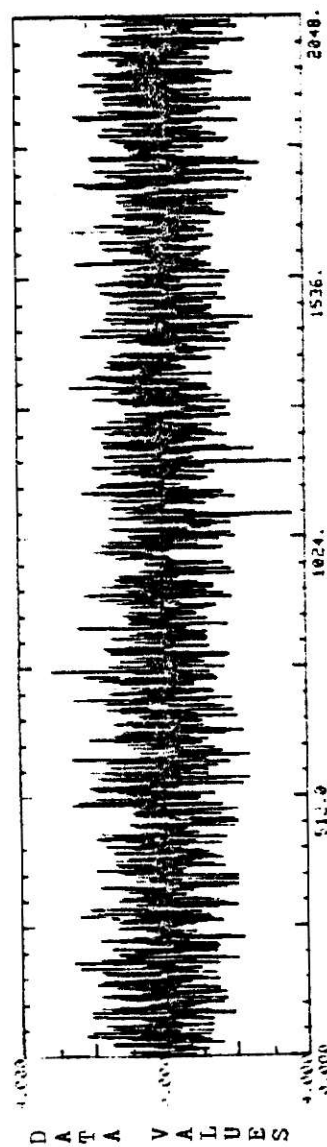


FIGURE 4.7



64 OUTPUT LEVELS

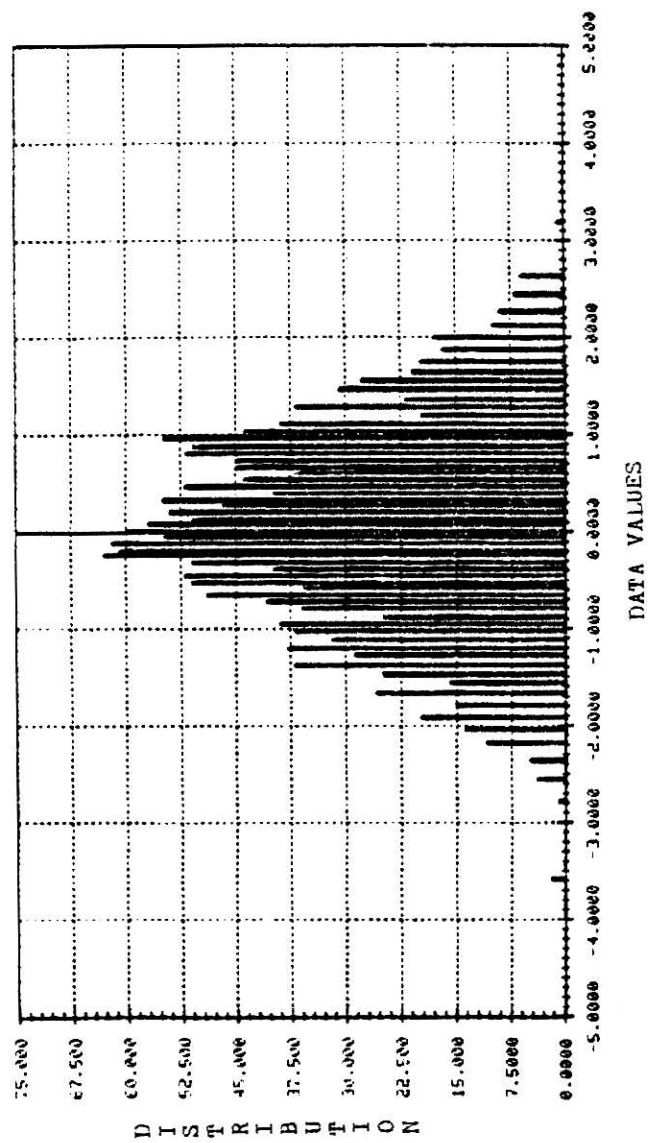


FIGURE 4.8

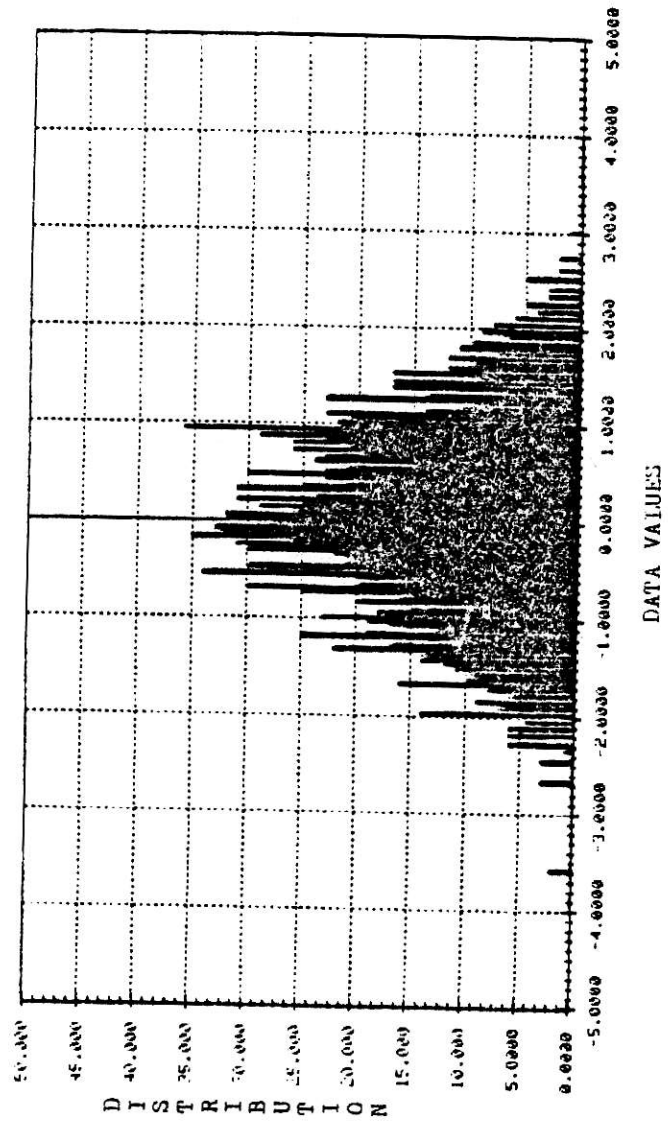
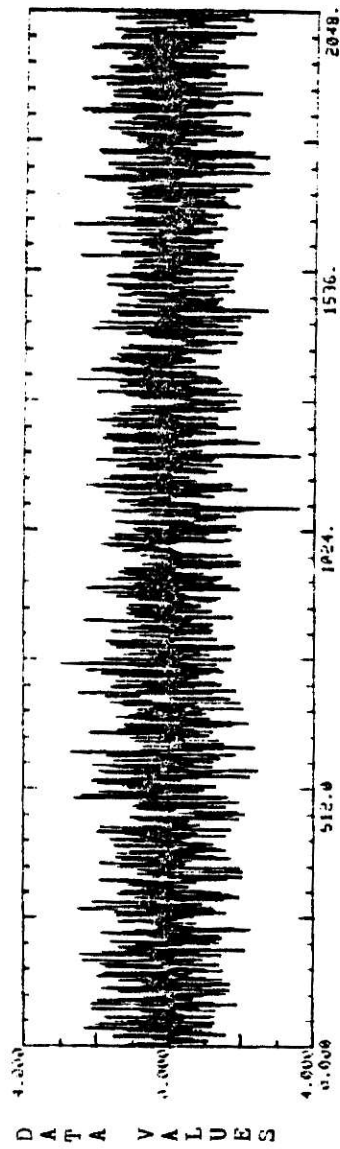


FIGURE 4.9

TABLE 4.1

EXPERIMENTAL VS. THEORETICAL
MEAN SQUARE ERROR
(EXAMPLE 1)

NUMBER OF OUTPUT LEVELS	THEORETICAL MSE [1]	EXPERIMENTAL MSE	RESCALED EXPERIMENTAL MSE
2	0.3634	0.382545	0.337679
4	0.1175	0.120184	0.106088
8	0.03454	0.034559	0.030506
16	0.009497	0.010177	0.008983
32	0.002499	0.002453	0.002165
64	0.000642	0.000666	0.000588
128	0.000162	0.000159	0.000140

EXAMPLE 2

A random gaussian input sequence with mean and variance equal to 250.11 and 5.6372 respectively. Note that the input sequence is in floating point arithmetic, while the output sequences are in fixed point arithmetic.

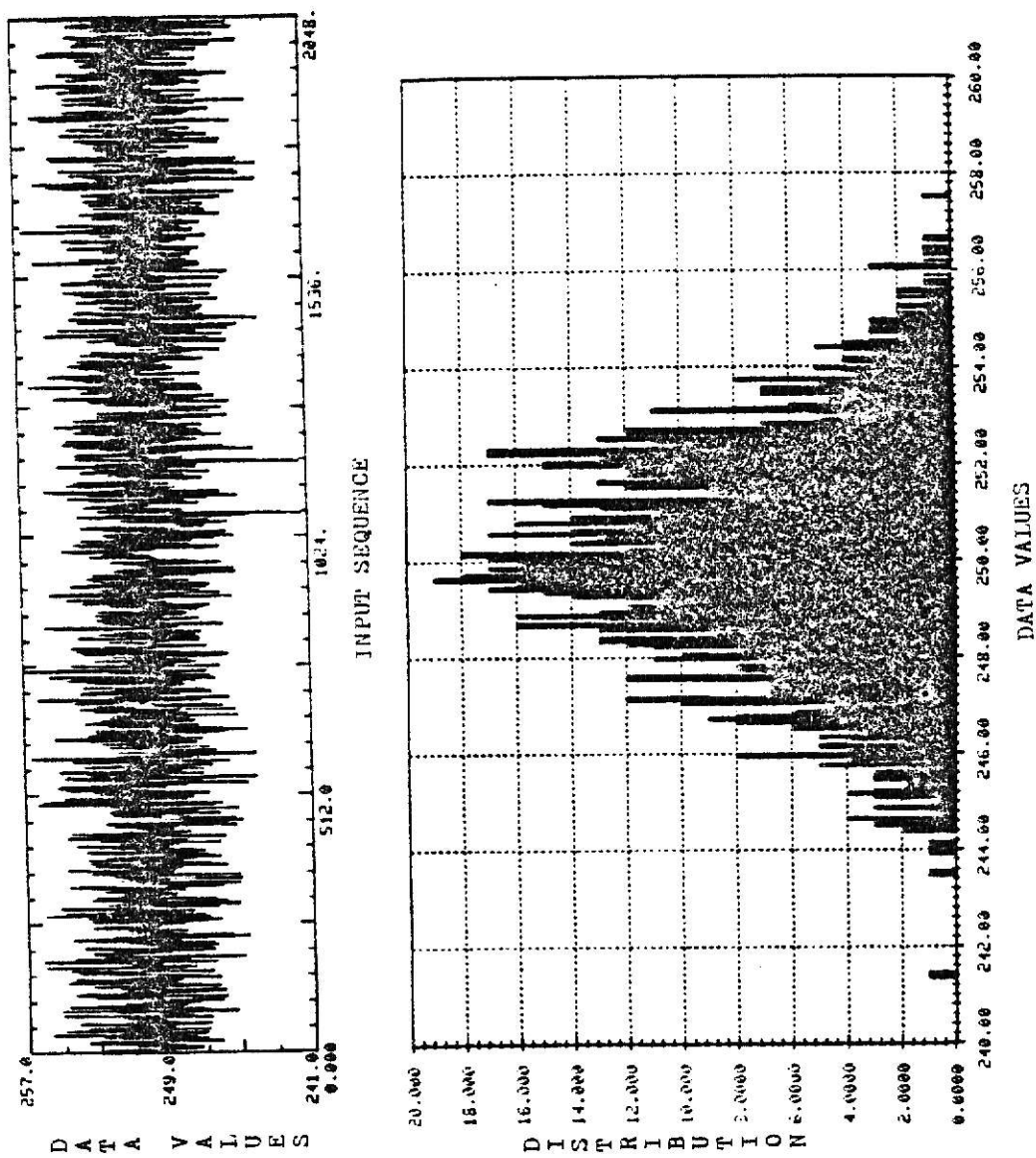
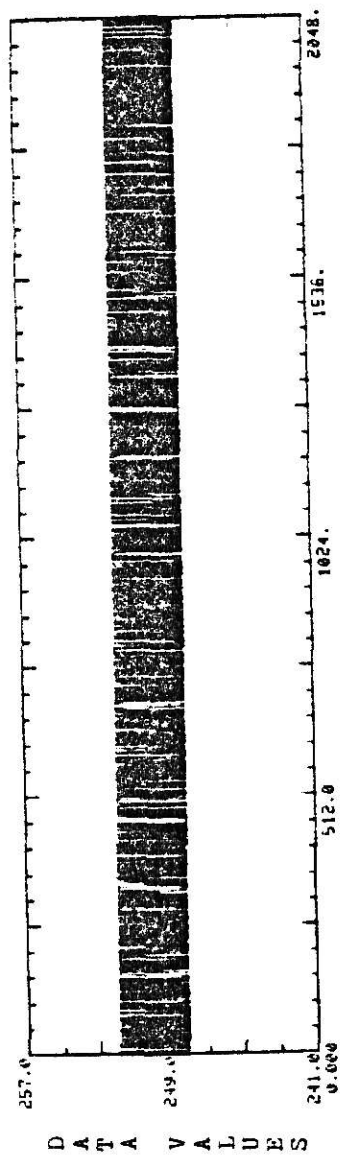


FIGURE 4.10



2 OUTPUT LEVELS

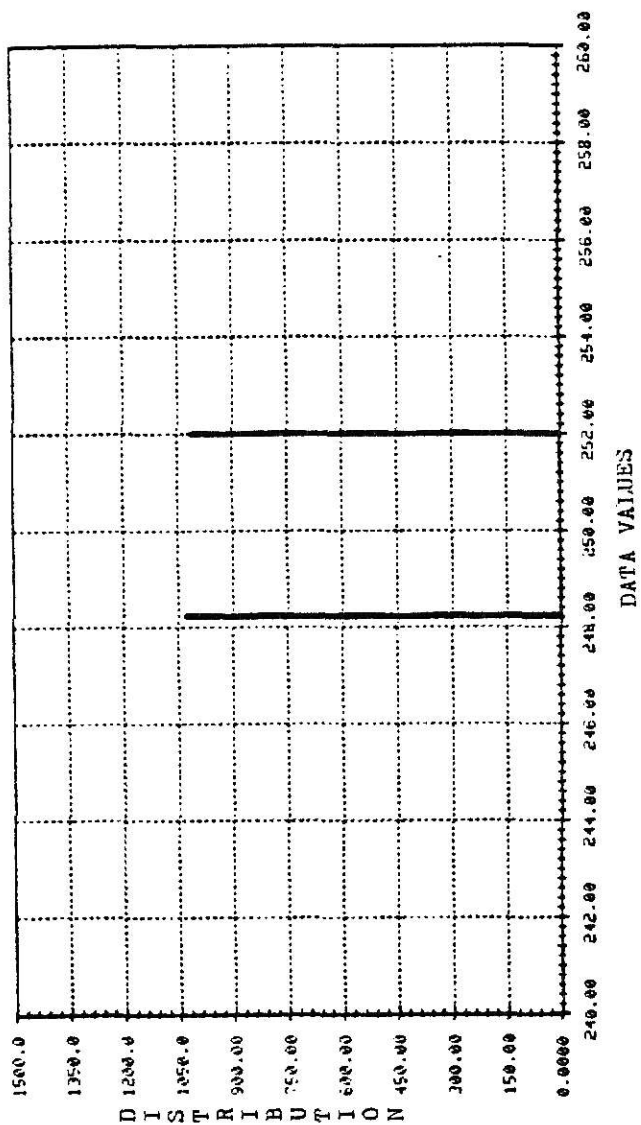
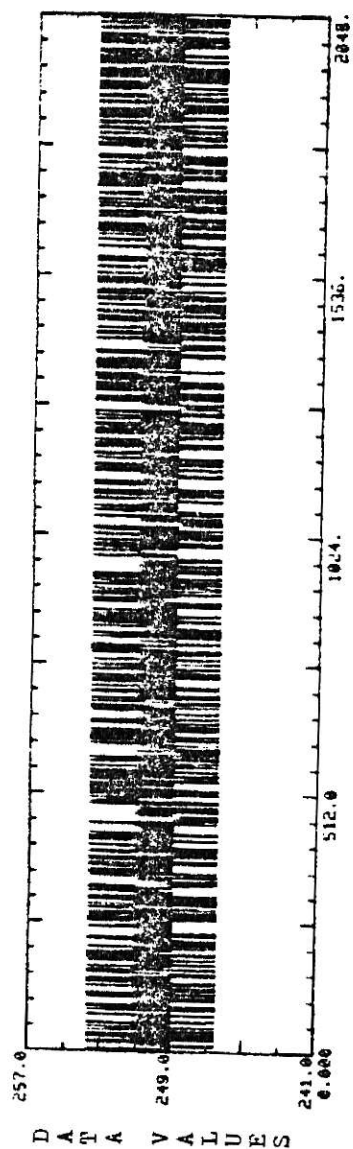


FIGURE 4.11



4 OUTPUT LEVELS

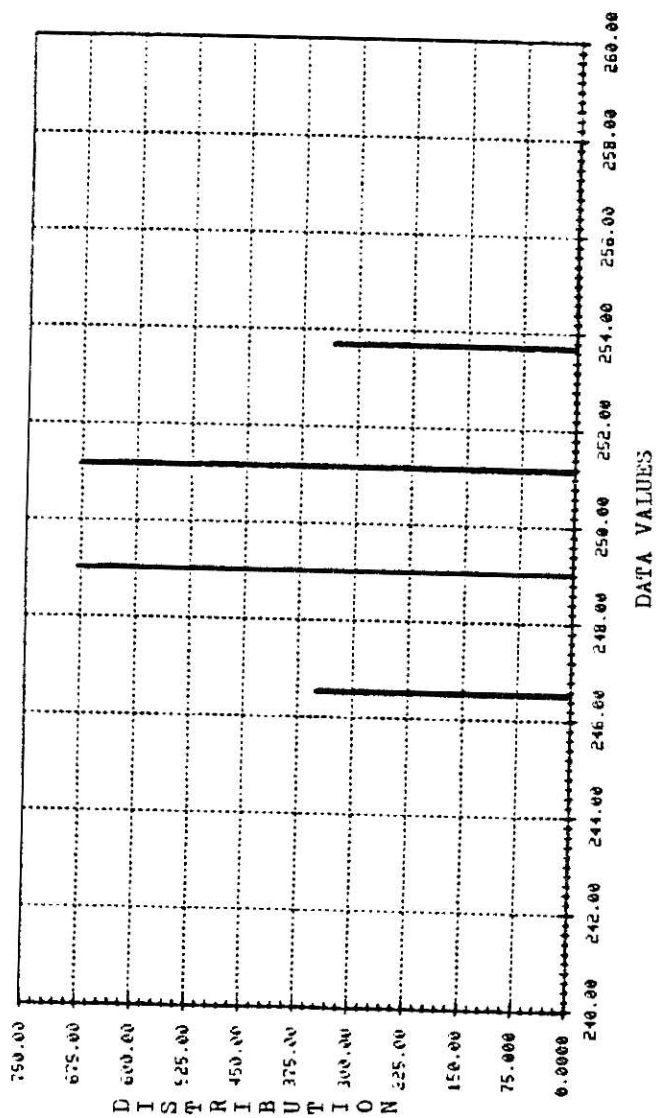
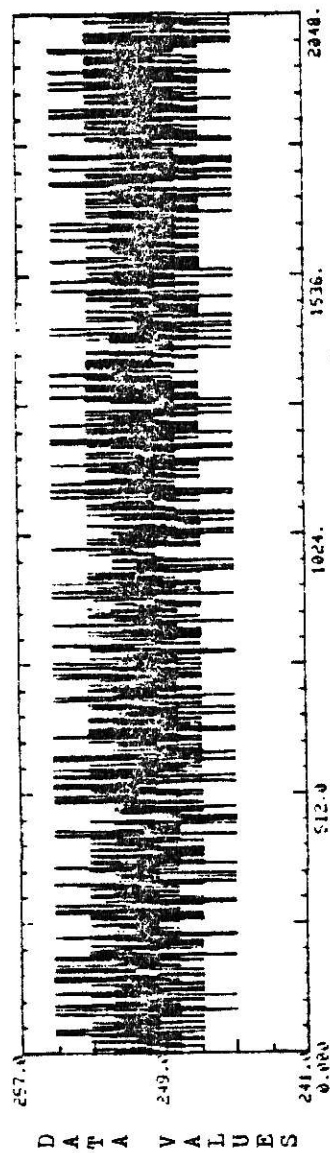


FIGURE 4.12



8 OUTPUT LEVELS

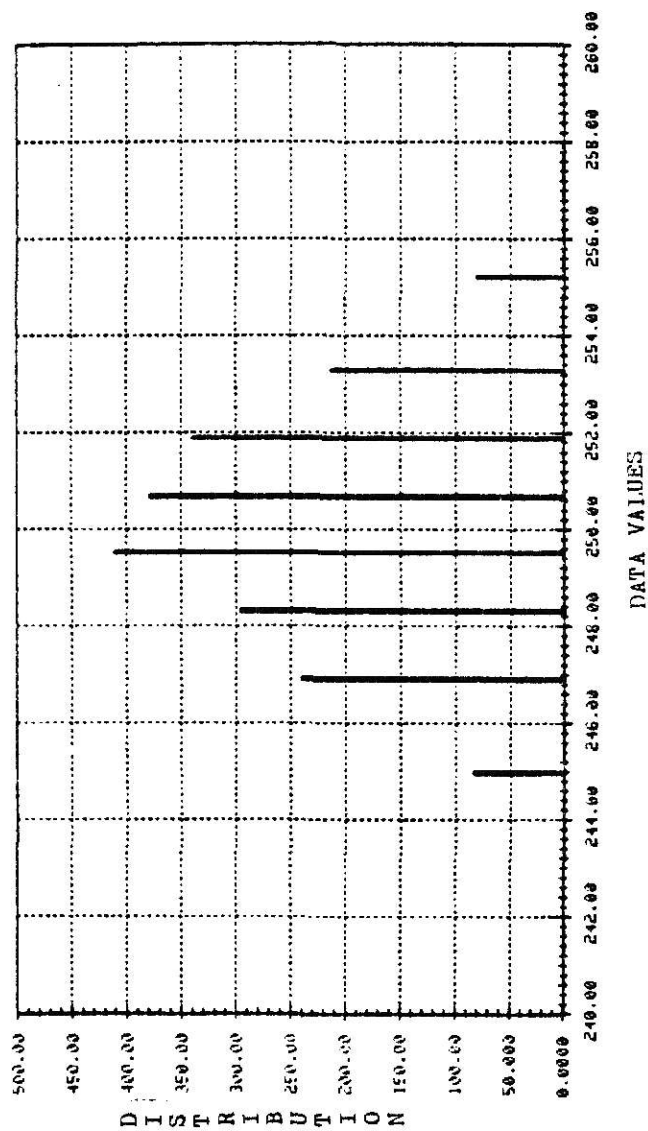


FIGURE 4.13

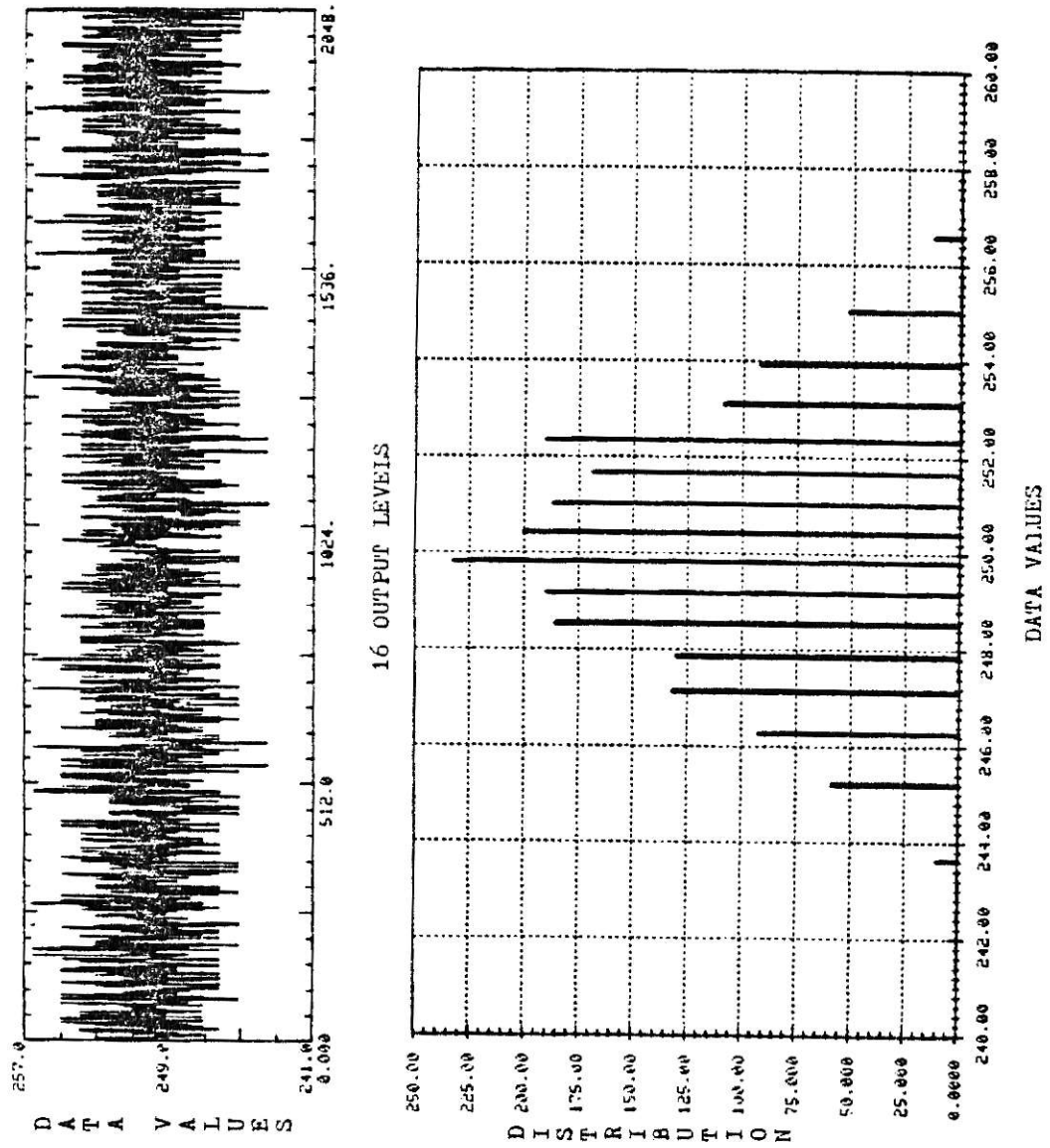


FIGURE 4.14

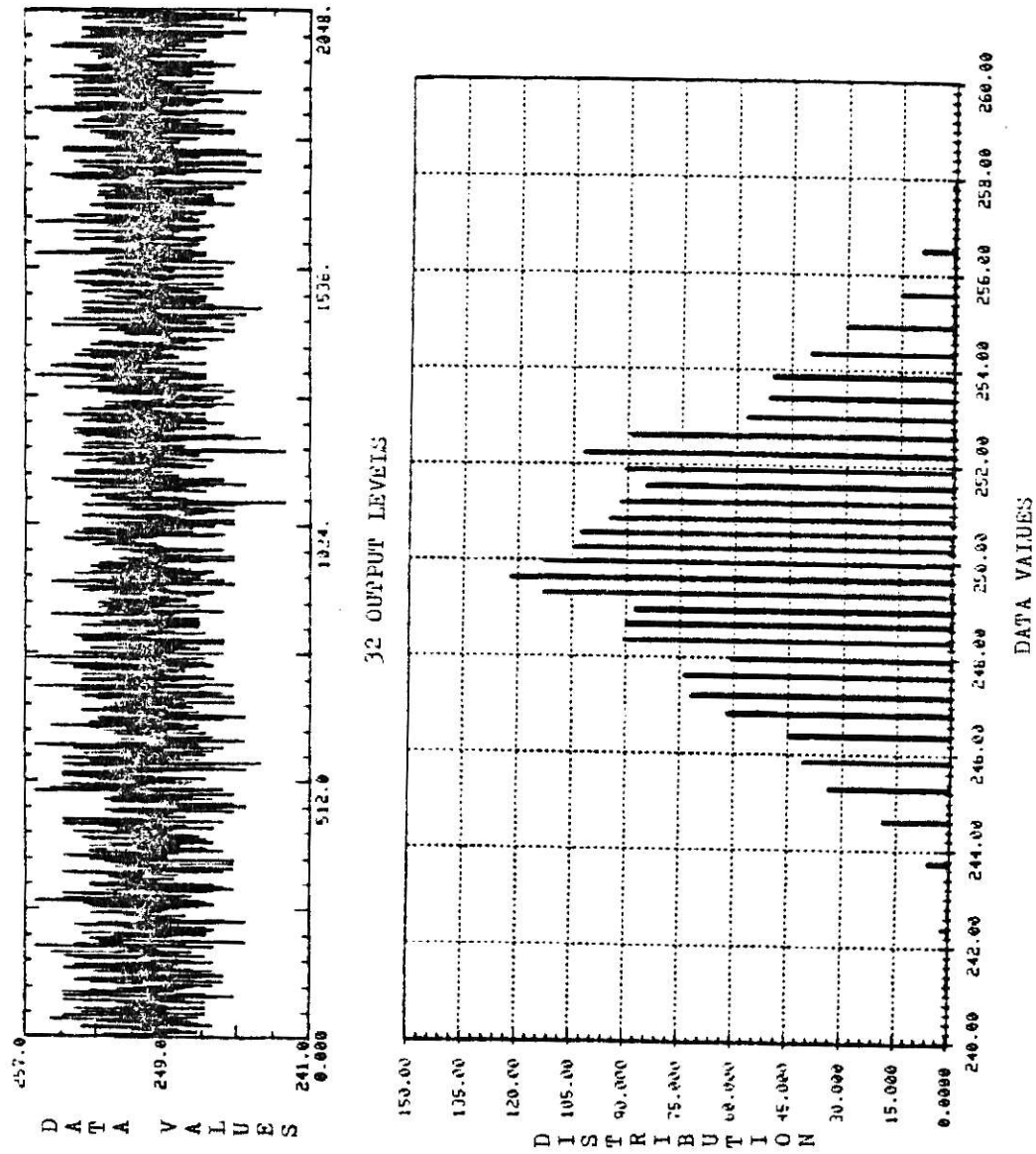


FIGURE 4.15

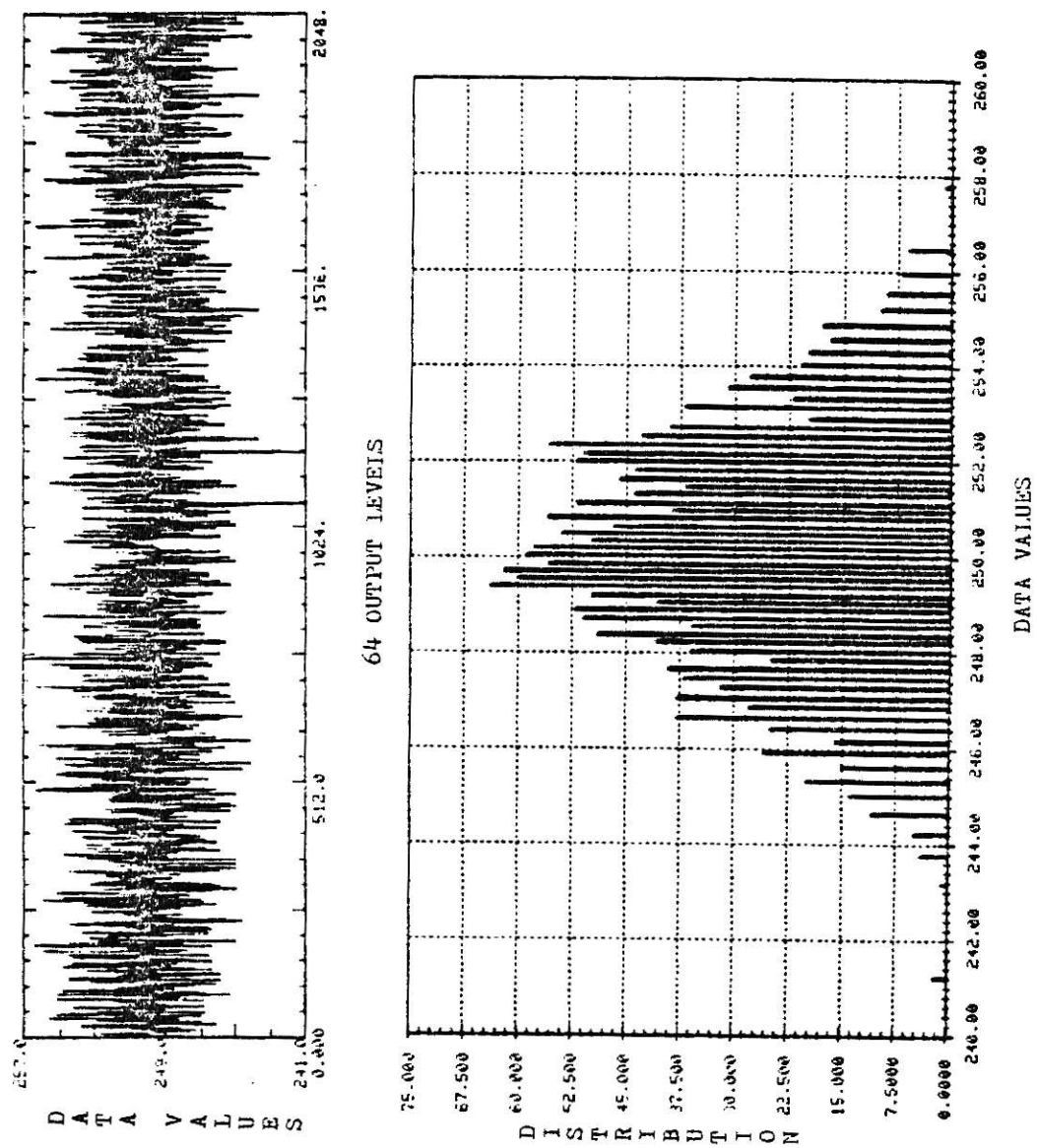


FIGURE 4.16

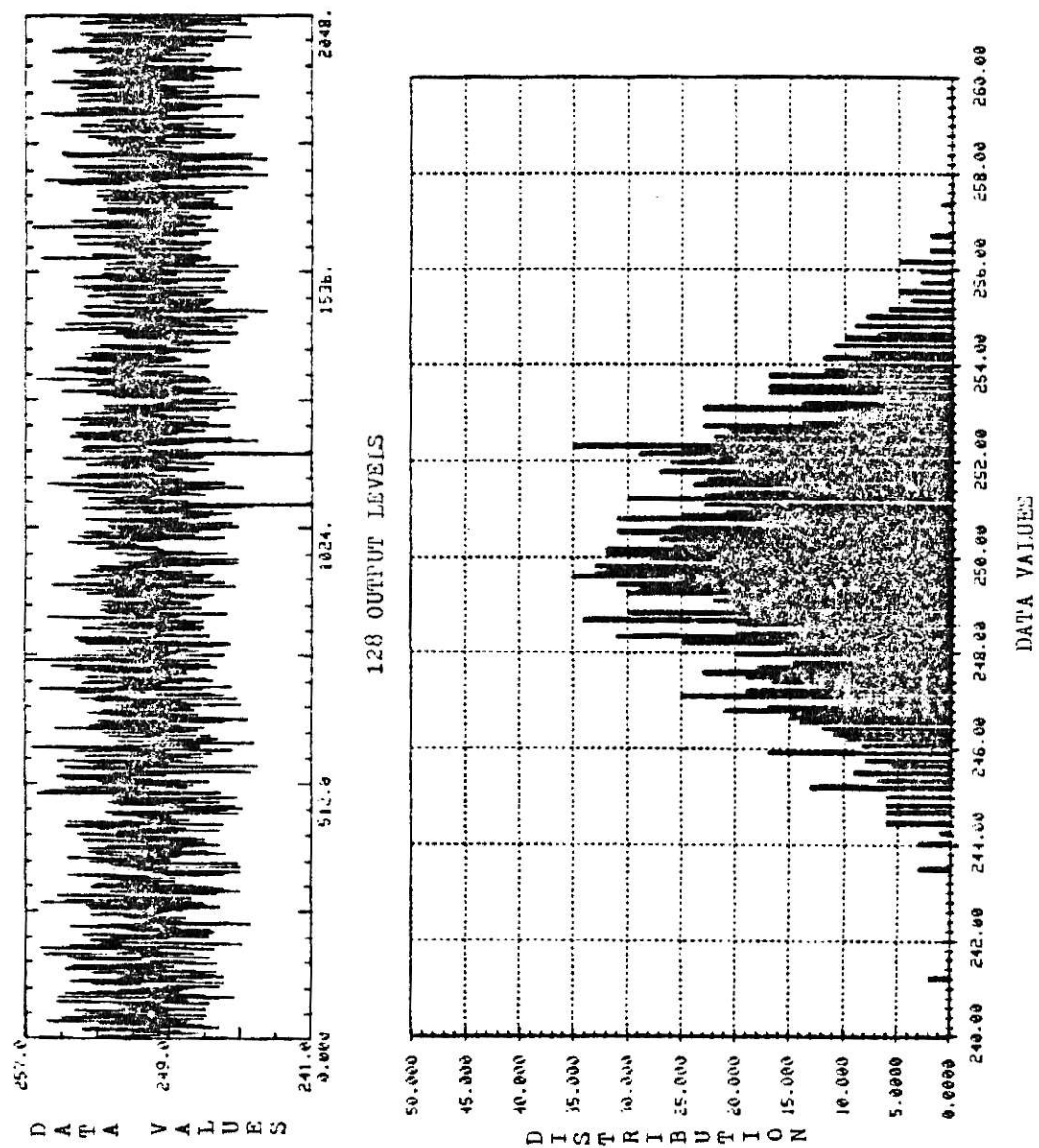


FIGURE 4.17

TABLE 4.2

EXPERIMENTAL VS. THEORETICAL
MEAN SQUARE ERROR
(EXAMPLE 2)

<u>NUMBER OF OUTPUT LEVELS</u>	<u>THEORETICAL MSE [7]</u>	<u>EXPERIMENTAL MSE</u>	<u>RESCALED EXPERIMENTAL MSE</u>
2	0.3634	0.063750	2.025848
4	0.1175	0.020033	0.636594
8	0.03454	0.005759	0.182994
16	0.009497	0.001697	0.053914
32	0.002499	0.000409	0.012989
64	0.000642	0.000111	0.003526
128	0.000162	0.000027	0.000843

APPENDIX 4.1

This appendix contains the program to generate a random gaussian data sequence of specified mean and variance.

NORMGEN.FR

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```

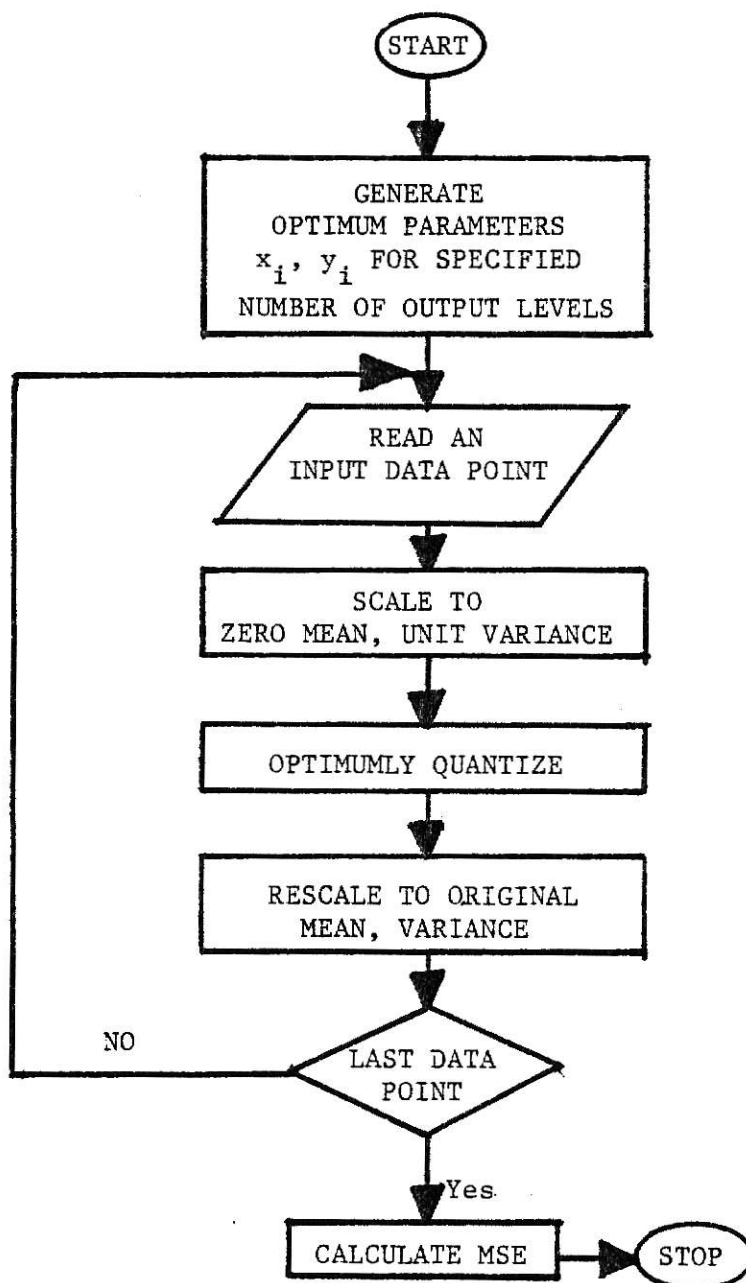
C      THIS PROGRAM GENERATES A RANDOM GAUSSIAN
C      DATA FILE OF SPECIFIED MEAN AND VARIANCE.
C
      REAL MEAN
      CALL OPENW(0,'OUTPUT FILENAME ? ',4,SIZE)
      ACCEPT 'NUMBER OF DATA POINTS IN THE OUTPUT FILE ? ',NPTS
      ACCEPT 'DESIRED MEAN AND VARIANCE OF THE OUTPUT FILE: MEAN,VAR ? ',MEAN,VAR
C
      DO 100 I=1,NPTS
      X=GAUSS(VAR,MEAN)
      CALL WRITR(0,I,X,1,IERR)
      CALL CHECK(IERR)
100   CONTINUE
      STOP
      END

```

APPENDIX 4.2

This appendix contains the program to optimumly quantize a gaussian input data sequence. This program also calculates the mean square error for the entire quantization process; i.e., the mean square error with respect to the input and output shown in Figure 4.1 for the entire sequence of data points.

FIGURE 4.18
THE OPTIMUM QUANTIZATION PROCESS



QUANTIZE1.FR 7/15/1982 01:11:2 DIR DPO Page 1

```

C THIS SUBROUTINE QUANTIZES A NORMALLY DISTRIBUTED
C INPUT DATA FILE INTO AN OUTPUT FILE
C CONTAINING A SPECIFIED NUMBER OF ASSIGNED
C OUTPUT VALUES (IE, LEVELS).
C
C ASSIGNMENT OF VARIABLES USED IN THIS SUBROUTINE
C
C NLEV...THE NUMBER OF OUTPUT LEVELS
C .....NOTE: 2**B=NLEV WHERE B IS THE NUMBER OF
C OUTPUT BITS
C
C MEAN...THE MEAN OF THE INPUT DATA FILE
C
C VAR...THE VARIANCE OF THE INPUT DATA FILE
C
C TABLES...QUANTIZATION TABLE (JOEL MAX, M.I.T. 1961)
C .....NOTE: TABLE(ROW,COLUMN) WHERE COLUMN 1 ARE
C THE CUT POINTS OR XI'S, AND COLUMN 2
C ARE THE REP-VALUES OR YI'S
C
C X(I)...THE INPUT DATA VALUES
C
C Y(I)...THE QUANTIZED OUTPUT VALUES
C
C
C COMPILER FREE
C REAL MEAN, MSE
C DIMENSION TABLE(128,2)
C DIMENSION X(2048), Y(2048)
C CALL OPENR(2,'INPUT FILENAME ? ',0,SIZE)
C CALL OPENW(3,'OUTPUT FILENAME ? ',0,SIZE)
C 2 FORMAT(' NUMBER OF OUTPUT LEVELS ? ',Z)
C 3 FORMAT(' VARIANCE OF THE INPUT FILE ? ',Z)
C 4 FORMAT(' MEAN OF THE INPUT FILE ? ',Z)
C
C ITTO=10
C ITTI=11
C ITTY=12
C WRITE(ITTO,2)
C READ(ITTI) NLEV
C WRITE(ITTO,3)
C READ(ITTI) VAR
C WRITE(ITTO,4)
C READ(ITTI) MEAN
C MSE=0.0
C
C GENERATE TABLE
C
C CALL TABLE(NLEV,TABLE)
C
C READ INPUT FILE INTO TABLE
C
C DO 100 I=1,2048
C READ BINARY(2,END=990) X(I)
C NPTS=I
C CALL LOOKUP(NLEV,MEAN,VAR,TABLE,I,X,Y)
C 100 CONTINUE
C TYPE 'INPUT FILE LENGTH MAY EXCEED 2048 POINTS'
C 990 CONTINUE
C
C CALCULATE MEAN SQUARED ERROR
C
C DO 995 I=1,NPTS
C MSE=((Y(I)-X(I))**2)/NPTS+MSE
C WRITE BINARY(3) Y(I)
C 995 CONTINUE
C WRITE(ITTO,1) MSE
C 1 FORMAT(' MEAN SQUARED ERROR = ',F16.6)
C
C CALL CLOSE(2,IERR)
C CALL CLOSE(3,IERR)
C STOP
C END

```

TABLGENB.FR

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```

C      THIS SUBROUTINE COMPUTES THE CUT-POINTS AND REP-VALUES FOR
C      A GAUSSIAN PROBABILITY DENSITY AND MSE DISTORTION MEASURE.
C
      SUBROUTINE TABLE(NLEV,TABLEG)
      COMPILER FREE
      DOUBLE PRECISION X,Y,MSE,ENTROPY,AREA1,AREA2
      DOUBLE PRECISION X0,X1,X2,P0,P2,XN,YN
      DOUBLE PRECISION P1,Y0,GAP,Z0,Z1,W0,W1,AREA,TAIL,A0,A1,W2,DABS,Q1,Q0
      DIMENSION TABLEG(128,2)
      ITT0=10
      ITT1=11
      ITT2=12
C
C      GET THE NUMBER OF LEVELS
C
      N=NLEV/2
C
C      ESTIMATE THE VALUE OF THE FIRST CUT-POINT OR REP-VALUE
C
      P1=2.1613D0/DBLE(FLOAT(NLEV))-1.653D0/DBLE(FLOAT(NLEV)**2.0)
C
C      PERFORM TEN ITERATIONS
C
      DO 1073 II=1,10
C
C      INITIALIZE THE FIRST CUT-POINT AND REP-VALUE
C
      IF(N*2.EQ.NLEV) GO TO 1100
C
C      ODD NUMBER OF LEVELS
C
      X0=P1
      Y0=2.0D0*P1
      GO TO 1110
1100  CONTINUE
C
C      EVEN NUMBER OF LEVELS
C
      X0=0.0D0
      Y0=P1
1110  CONTINUE
      I=1
C
C      THE FIRST CUT-POINT VALUE
C
      TABLEG(I,1)=X0
      TABLEG(I,2)=Y0
      IF(NLEV.LT.4) GO TO 1000
C
C      DETERMINE THE REST OF THE CUT-POINTS AND REP-VALUES
C
      DO 1000 I=2,N
      CALL CUTPT(X0,Y0,X1)
C
      X0=X1
      Y0=2.0D0*X0-Y0
      TABLEG(I,1)=X0

```

T4BLGENS.FR 7/15/1982 01:14:47 DIR DPO Page 2

```

      TABLE(I,2)=Y0
1000  CONTINUE
C
C      USING AN ALTERNATE METHOD,
C      DETERMINE THE LAST REP-VALUE, Y(N), BY COMPUTING
C      THE CENTROID OF THE AREA FROM X(N) TO INFINITY
C
      XN=X0
      CALL ERF(XN,AREA)
      YN=DEXP(-XN**2/2.0D0)/DSORT(2.0D0*3.141592D0)/(0.5D0-AREA)
C
C      COMPUTE THE DIFFERENCE BETWEEN THE TWO VALUES AND UPDATE
C      THE STARTING VALUE, P1
C
      Q1=Y0-YN
      IF(0ABS(Q1/YN).LT.1.0D-6) RETURN
      IF(11.EQ.1) P2=0.99D0*P1
      IF(11.NE.1) P2=(P1*Q0-P0*Q1)/(Q0-Q1)
      P0=P1
      P1=P2
      Q0=Q1
1073  CONTINUE
      STOP
      END

```

CUTPT2,FR 7/15/1982 01:11:8 DIR DPO Page 1

```

C      THIS SUBROUTINE COMPUTES THE NEXT CUT-POINT, X1, GIVEN THE
C      PREVIOUS CUT-POINT AND REP-VALUE
C
      SUBROUTINE CUTPT(X0,Y0,X1)
      DOUBLE PRECISION X0,Y0,X1,EPS,AREA,AREA0,AREA1,DELX1,PROB,FUNC,DABS
      DOUBLE PRECISION TAIL
      EPS=1.0D-9
C
C      INTEGRATE THE FUNCTION FROM X0 TO Y0
C
      CALL FPX(X0,Y0,AREA0)
C
      X1=2.0D0*Y0-X0
C
C      ITERATE UP TO 100 TIMES TO DETERMINE X1
C
      DO 1000 I=1,100
      CALL FPX(X1,Y0,AREA1)
      AREA=AREA1-AREA0
      DELX1=-(AREA)/(PROB(X1)*FUNC(X1,Y0))
      IF (DABS(DELX1/X1).LT.EPS) RETURN
      X1=X1+DELX1
1000  CONTINUE
      RETURN
      END

```

ERF2,FR 7/15/1982 01:11:6 DIR DPO Page 1

```

C      THIS SUBROUTINE EVALUATES THE INTEGRAL OF THE
C      GAUSSIAN PROBABILITY DISTRIBUTION VIA AN
C      INFINITE SERIES REPRESENTATION
C
      SUBROUTINE ERF(X,AREA)
      DOUBLE PRECISION X,U,AREA,AREA1,DEXP,DSQRT,DBLE,DABS
      U=1.0D0
      AREA1=X*DEXP(-X**2/2.0D0)/DSQRT(2.0D0*3.141592D0)
      DO 100 J=1,100
      K=2*J+1
      U=U*DBLE(FLOAT(K))
      AREA=(X/DSQRT(2.0D0))**K*2.0D0**J/U*DEXP(-X**2/2.0D0)/DSQRT(3.141592D0)+AREA1
      IF (DABS(AREA1-AREA).LT.1.0D-9) RETURN
      AREA1=AREA
100  CONTINUE
      TYPE 'NEED MORE TERMS OF THE INFINITE SERIES'
      RETURN
      END

```

FPX,FR 7/15/1982 01:11:11 DIR DPO Page 1

```

      SUBROUTINE FPX(X,Y,AREA)
      DOUBLE PRECISION X,Y,AREA,DEXP,DSQRT,A
      CALL ERF(X,A)
      AREA=(1-DEXP(-X**2/2.0D0))/DSQRT(2.0D0*3.141592D0)-Y**A
      RETURN
      END

```

FUNC.FR 7/15/1982 01:11:10 DIR BFO Page 1

```
DOUBLE PRECISION FUNCTION FUNC(X,YO)
DOUBLE PRECISION X,YO
FUNC=X-YO
RETURN
END
```

PROB.FR 7/15/1982 01:11:09 DIR BFO Page 1

```
DOUBLE PRECISION FUNCTION PROB(X)
DOUBLE PRECISION X,DEXP,DSQRT
PROB=DEXP(-X**2/2.0D0)/DSQRT(2.0D0*3.141592D0)
RETURN
END
```

LOOKUP.FR 7/15/1982 01:15 DIR DPO Page 1

```

SUBROUTINE LOOKUP(NLEV,MEAN,VAR,TABLEG,I,X,Y)
COMPILER FREE
REAL MEAN
DIMENSION TABLEG(128,2)
DIMENSION X(2048), Y(2048)
C
C   ITT0=10
C   ITT1=11
C   ITT2=12
C
C   MAXJ=NLEV/2
C   LAST=MAXJ-1
C   SCALE INPUT DATA TO ZERO MEAN AND UNIT VARIANCE
C
C   XSCALE=(X(I)-MEAN)/SQRT(VAR)
C
C   LOOKUP TABLE VALUES ARE POSITIVE
C
C   XSCAL=ABS(XSCALE)
C
C   ODD NUMBER OF LEVELS
C
C   IF(XSCAL.GE.TABLEG(1,1)) GO TO 50
C   Y(I)=0.0
C   RETURN
C
C   EVEN OR ODD NUMBER OF LEVELS
C
C   50 DO 100 J=1,LAST
C   IF(XSCAL.GE.TABLEG(J+1,1)) GO TO 100
C   J2=J
C   GO TO 200
C   100 CONTINUE
C
C   J2=MAXJ
C
C   200 YTABLE=TABLEG(J2,2)*XSCALE/XSCAL
C
C   NOTE XSCALE/XSCAL EQUALS +1 OR -1
C
C   RESCALE TO ORIGINAL MEAN AND VARIANCE
C
C   Y(I)=(YTABLE*SQRT(VAR))+MEAN
C
C   RETURN
C   END

```

CHAPTER V

PROGRAM USAGE

Two programs are used. The first, TABLGEN5, generates the optimum gaussian quantizer parameters, X_i and Y_i , in tabular form. The second, QUANTIZE1, takes a random gaussian input sequence, scales it to a normalized gaussian sequence and then passes the entire sequence through the optimum quantizer for a specified number of output levels. The quantizer output is then rescaled to the original mean and variance of the input sequence.

The first program, TABLGEN5, generates the optimum parameters. There are three inputs for which the program will prompt the user. The first two inputs are the names of output files. One of these stores the value of the mean square error for the specified number of output levels. These values can be retrieved at a later time and plotted as in Figure 3.12. Similarly the second output filename is used to store the value of entropy being calculated. The last input that is entered is the number of output levels desired for the optimum quantizer. All values are entered from the CRT terminal. A typical input sequence is illustrated in Figure 5.1. The output consists of the values x_i and y_i , ending with the values for the mean square error and entropy for the specific number of output levels being calculated, as exemplified in Figure 5.2. The output is printed on the line printer.

The second program, QUANTIZE1, requires the prior use of a library program, XBAR, to determine the mean and variance of the gaussian input sequence. QUANTIZE1 requires five inputs to the CRT terminal, all of which are prompted for: the output sequence filename, the input filename of the random gaussian sequence, the mean and variance of the input sequence obtained from XBAR, and finally the desired number of output

FIGURE 5.1

TABLGEN5
Input Sequence

TABLGEN5

MEAN SQUARE ERROR OUTPUT FILENAME ? MSE.FP

ENTROPY OUTPUT FILENAME ? ENTROPY.FP

NUMBER OF OUTPUT LEVELS ? 4

NUMBER OF OUTPUT LEVELS ? 5

INT

R

FIGURE 5.2

TABLGEN5
Output Sequence

4 OUTPUT LEVELS

N= 4

ITERATION NUMBER	1	J	X(J)	Y(J)
	1		.000000	.437013
	2		.941286	1.445560
SECANT METHOD.....			Y(N) =	1.473315

N= 4

ITERATION NUMBER	2	J	X(J)	Y(J)
	1		.000000	.432642
	2		.930265	1.427887
SECANT METHOD.....			Y(N) =	1.469573

N= 4

ITERATION NUMBER	3	J	X(J)	Y(J)
	1		.000000	.432642
	2		.992275	1.511508
SECANT METHOD.....			Y(N) =	1.510958

N= 4

ITERATION NUMBER	4	J	X(J)	Y(J)
	1		.000000	.432776
	2		.981588	1.510400
SECANT METHOD.....			Y(N) =	1.510409

N= 4

ITERATION NUMBER	5	J	X(J)	Y(J)
	1		.000000	.432730
	2		.981500	1.510419
SECANT METHOD.....			Y(N) =	1.510419

MEAN SQUARED ERROR FOR 4 OUTPUT LEVELS = .117481

ENTROPY FOR 4 OUTPUT LEVELS = 1.911098

FIGURE 5.2 (CONT.)

TABLGEN5
Output Sequence

5 OUTPUT LEVELS

N= 5

ITERATION NUMBER	1	J	X(J)	Y(J)
	1		.366140	.732280
	2		1.181830	1.631381
SECANT METHOD.....			Y(N) =	1.672620

N= 5

ITERATION NUMBER	2	J	X(J)	Y(J)
	1		.360479	.724957
	2		1.167906	1.610854
SECANT METHOD.....			Y(N) =	1.661201

N= 5

ITERATION NUMBER	3	J	X(J)	Y(J)
	1		.332718	.766434
	2		1.246066	1.726697
SECANT METHOD.....			Y(N) =	1.725562

N= 5

ITERATION NUMBER	4	J	X(J)	Y(J)
	1		.332272	.764544
	2		1.244311	1.721079
SECANT METHOD.....			Y(N) =	1.724110

N= 5

ITERATION NUMBER	5	J	X(J)	Y(J)
	1		.362284	.764568
	2		1.244358	1.724148
SECANT METHOD.....			Y(N) =	1.724149

MEAN SQUARED ERROR FOR 5 OUTPUT LEVELS = .079941

ENTROPY FOR 5 OUTPUT LEVELS = 2.202916

levels from the optimum quantizer. The output file is written to permanent storage as previously mentioned. The only other output is the value of the mean square error, as mentioned in Appendix 4.2, to the CRT terminal. A typical input/output sequence is illustrated in Figure 5.3.

FIGURE 5.3

QUANTIZE1
Input/Output Sequence

```
QUANTIZE1
INPUT FILENAME ? G2048.01
OUTPUT FILENAME ? G64.01
NUMBER OF OUTPUT LEVELS ? 64
VARIANCE OF THE INPUT FILE ? 93953
MEAN OF THE INPUT FILE ? .044654
INPUT FILE LENGTH MAY EXCEED 2048 POINTS
MEAN SQUARED ERROR = .000530

STOP
R
```

CHAPTER VI

CONCLUSIONS AND RECOMMENDATIONS

The programs developed in this report illustrate the accuracy problems encountered when using fixed-point arithmetic as exemplified in Figures 6.1 and 6.2 (Note only the input sequence is in floating point arithmetic). However, since special purpose fixed-point hardware in the field can be implemented faster and more economically, these optimal programs are of value.

The computer programs presented in this report can be utilized as a subroutine in applications such as speech and image data compression. In both instances the quantization occurs after the raw data has been transformed (FFT, DCT, etc.). Figure 6.3 [7] depicts the image processing scheme incorporating the optimum quantization process of Figure 4.1. It should be noted that not all pixels are quantized with the same number of output levels; for instance zonal coding, Figure 6.4 may be the scheme employed for compression [7]. Thus a subroutine that can quantize, encode, various output levels is of great value in such an application. Finally it should be pointed out that in processing the smaller $M \times M$ pixel blocks of Figure 6.3, all the elements are assumed to have a gaussian distribution, $p_X(x)$, except the very first element (1,1) which is the uniformly distributed dc term. Therefore our gaussian quantizer will apply to all but one term. For this singular term, we could go back and calculate optimum parameters for the uniform distribution, $p_X(x)$. However, the internal hardware of the computer itself assigns numbers based on the uniform distribution. Thus, with our gaussian parameters and the inherent uniform distribution of the computer, we have all the ingredients to implement the optimum quantizer for the image data compression scheme of Figure 6.3.

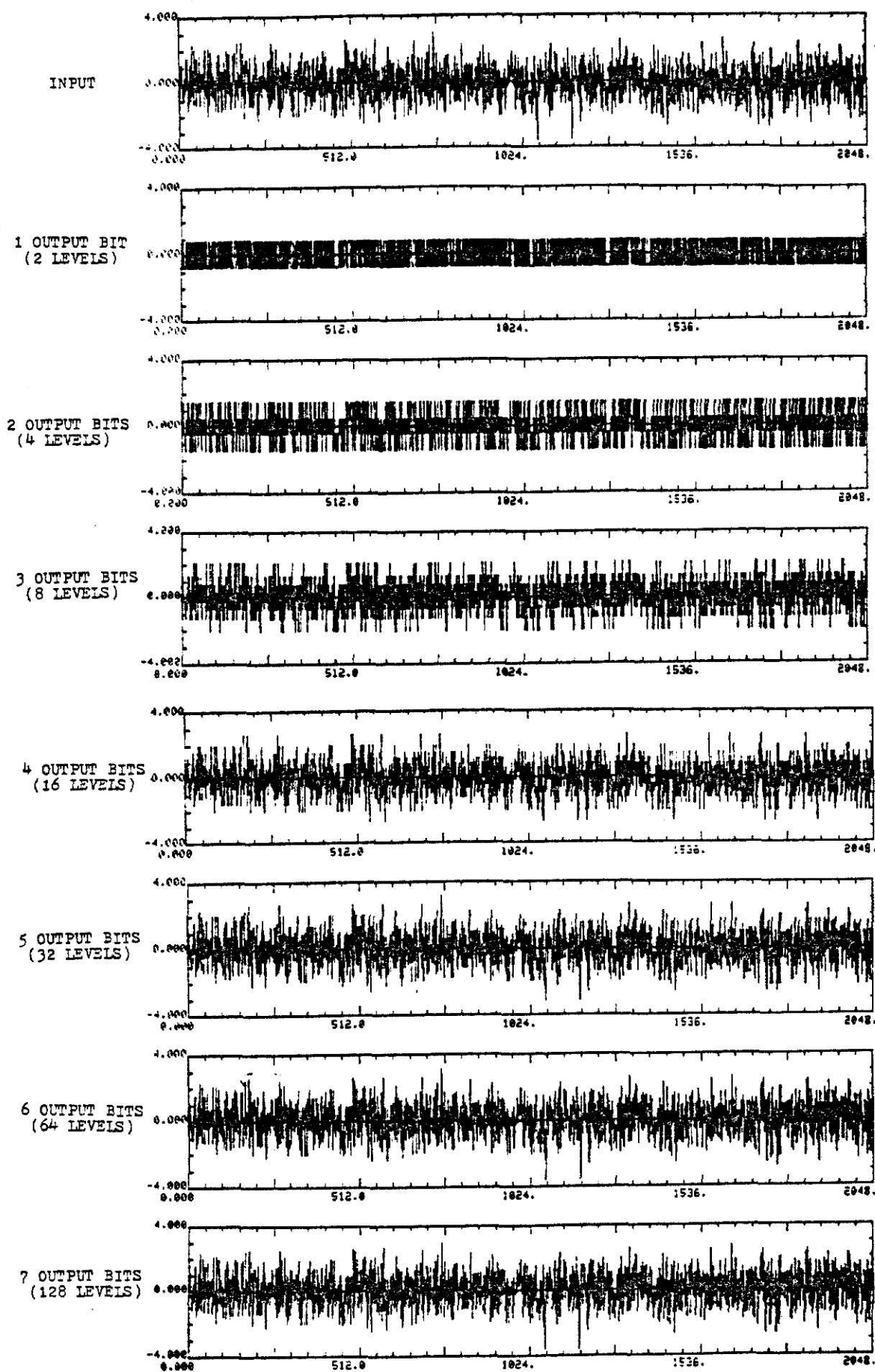
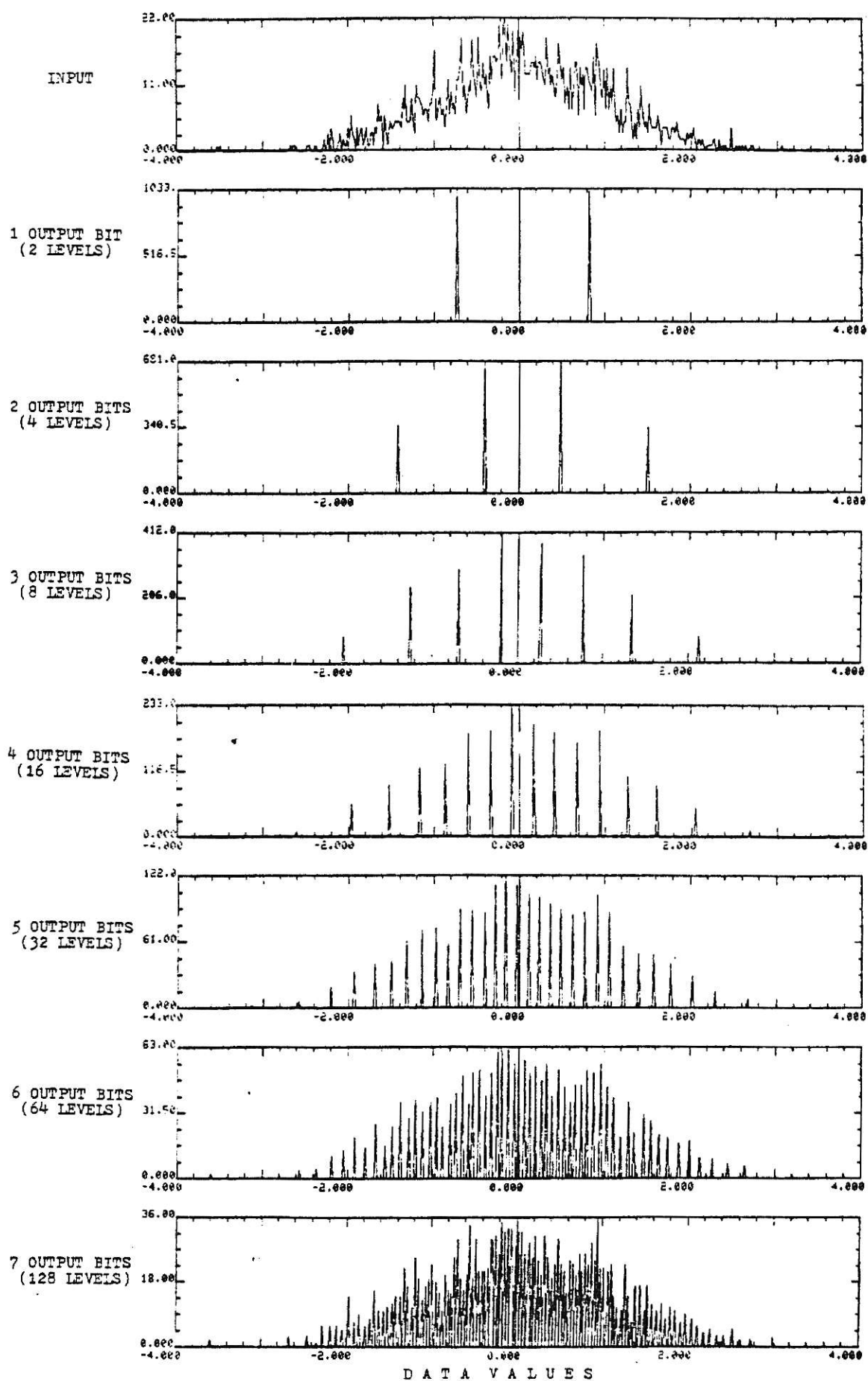


FIGURE 6.1



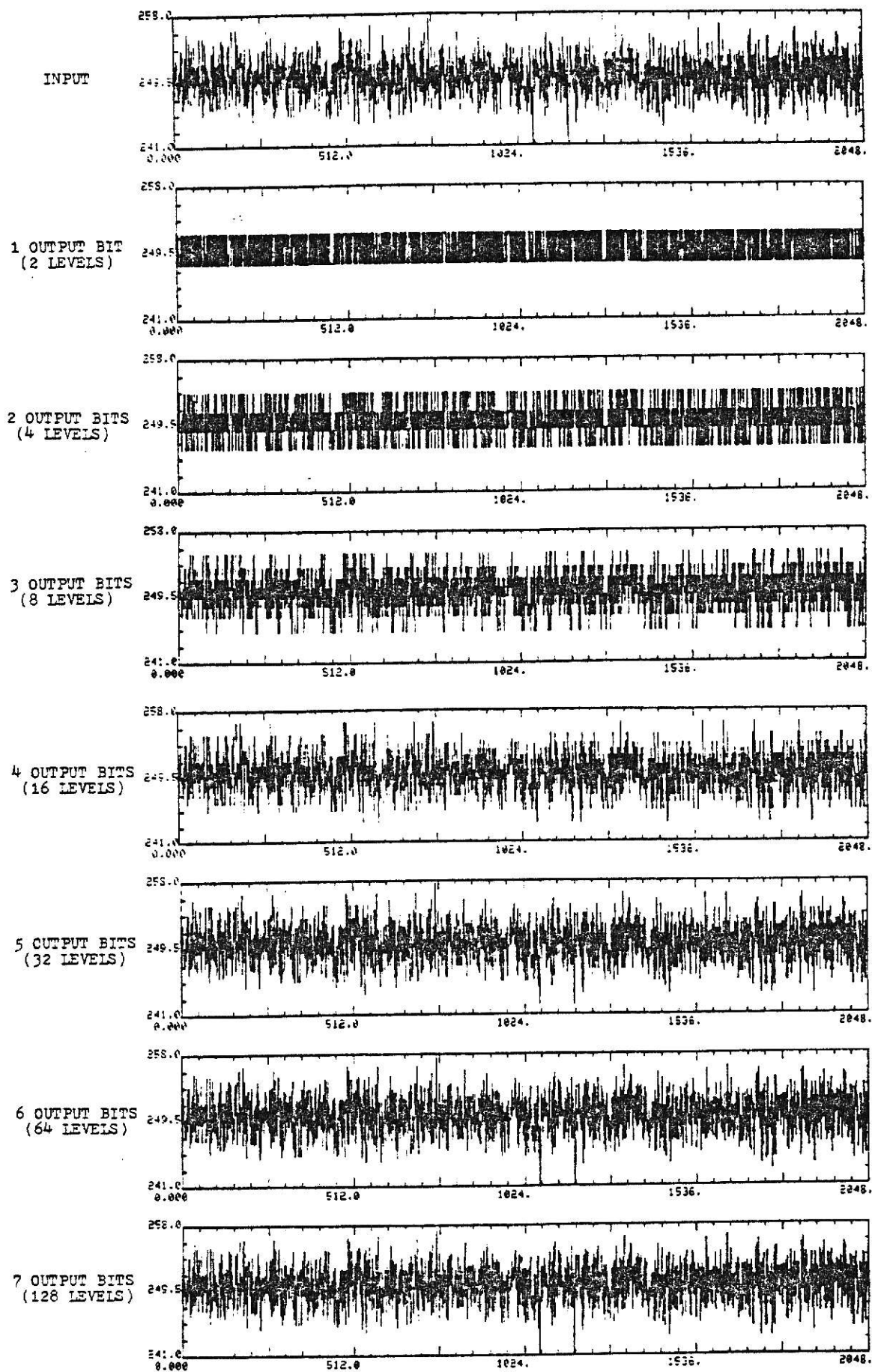


FIGURE 6.2

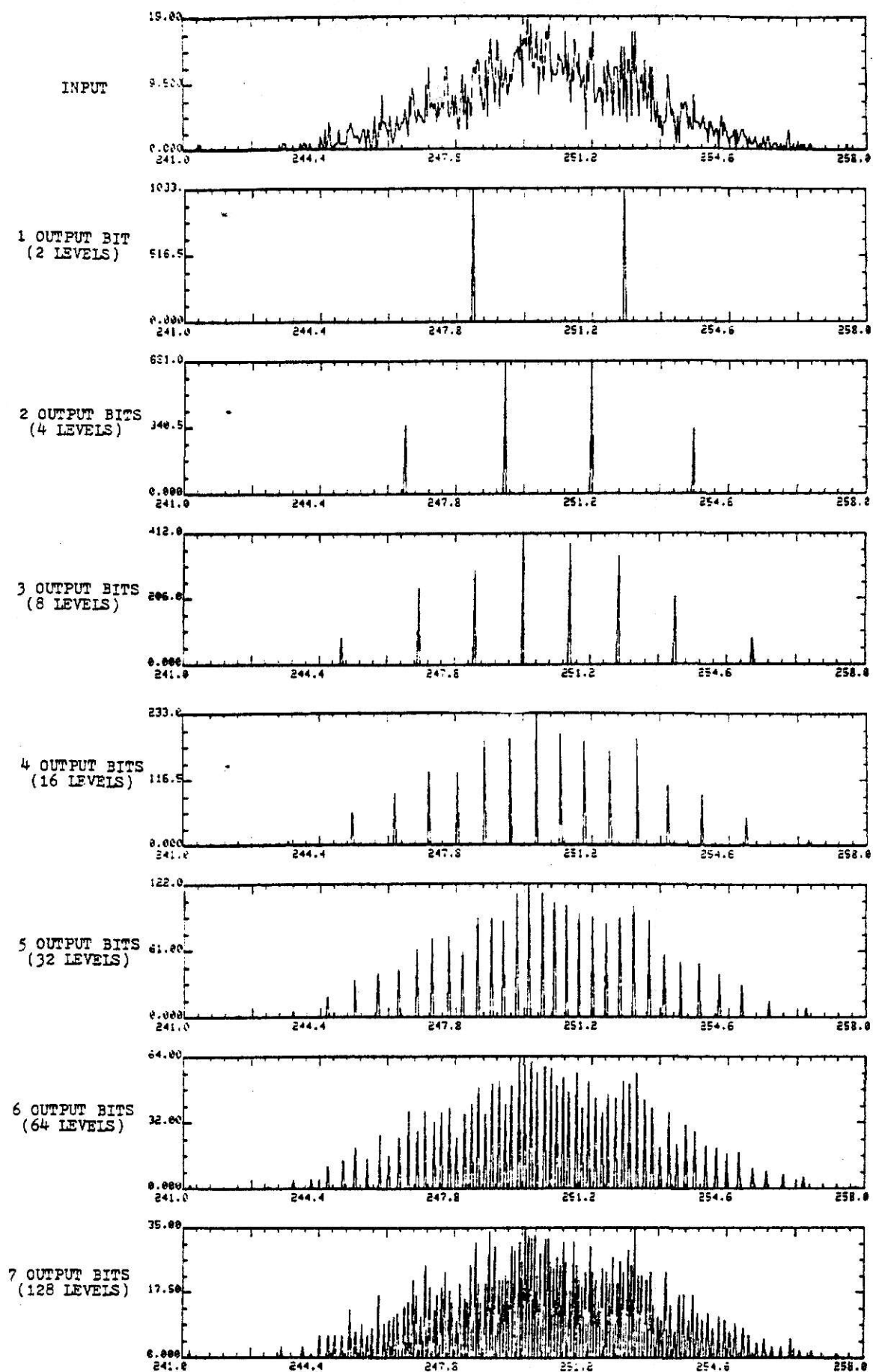


FIGURE 6.2 (CONT.)

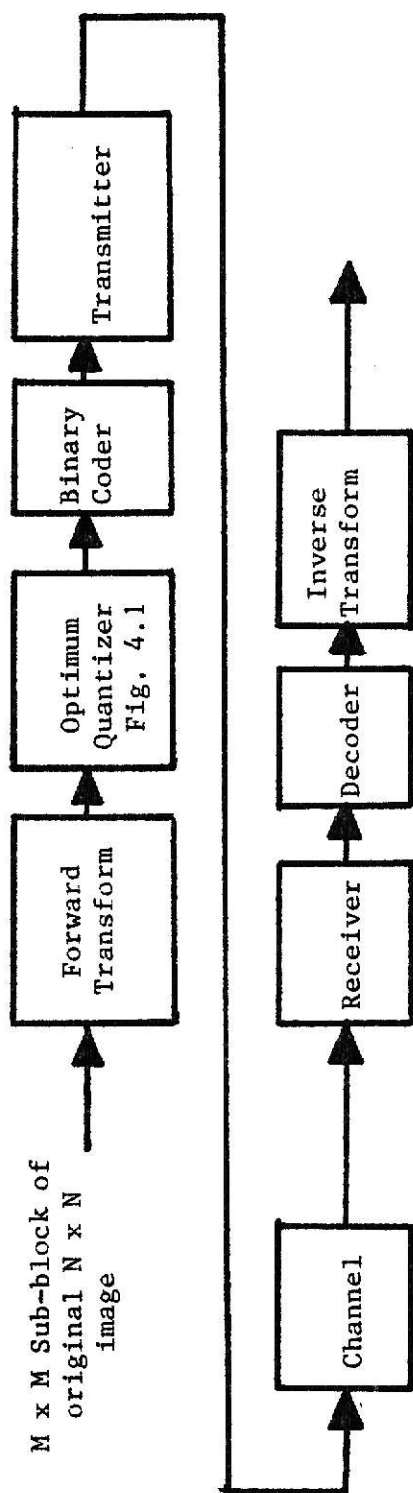


IMAGE PROCESSING SCHEME

FIGURE 6.3

```

8 8 8 7 7 7 7 5 5 4 4 4 4 4 4 4
8 8 7 6 5 5 3 3 3 3 2 2 2 2 2 2
8 7 6 4 4 4 3 3 2 2 2 2 2 2 2 2
7 6 4 3 2 2 2 1 1 1 1 0 0 0 0
7 5 4 2 2 2 1 1 1 0 0 0 0 0 0
7 5 4 2 2 2 1 1 0 0 0 0 0 0 0
5 3 3 2 2 1 1 0 0 0 0 0 0 0 0
5 3 3 2 1 1 0 0 0 0 0 0 0 0 0
4 3 2 1 1 0 0 0 0 0 0 0 0 0 0
4 3 2 1 1 0 0 0 0 0 0 0 0 0 0
4 3 2 1 0 0 0 0 0 0 0 0 0 0 0
4 3 2 1 0 0 0 0 0 0 0 0 0 0 0
4 2 2 0 0 0 0 0 0 0 0 0 0 0 0
4 2 2 0 0 0 0 0 0 0 0 0 0 0 0
4 2 2 0 0 0 0 0 0 0 0 0 0 0 0
4 2 2 0 0 0 0 0 0 0 0 0 0 0 0

```

ZONAL CODING

Typical numbers of output bits assigned for zonal transform coding.
 A 16 x 16 block is illustrated, each element is a pixel.

FIGURE 6.4

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THE OPTIMUM QUANTIZATION PROCESS

by

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B.S., Kansas State University, 1979

AN ABSTRACT OF A MASTER'S REPORT

submitted in partial fulfillment of the

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KANSAS STATE UNIVERSITY

Manhattan, Kansas

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ABSTRACT

The objective of this work is to mathematically describe the optimum quantizer. Once the results have been obtained for the general optimum quantizer, a specific case is considered. A digital computer implementation then establishes experimental values based on the mathematical derivation for the optimum quantizer. The experimental and theoretical values are then compared. The specific case considered is the optimum quantization process for a random gaussian sequence.