

Simple quantum error-correcting codes

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Methods of finding good quantum error-correcting codes are discussed, and many example codes are presented. The recipe $C_2^\perp \subseteq C_1$, where C_1 and C_2 are classical codes, is used to obtain codes for up to 16 information quantum bits (qubits) with correction of small numbers of errors. The results are tabulated. More efficient codes are obtained by allowing C_1 to have reduced distance, and introducing sign changes among the code words in a systematic manner. This systematic approach leads to single-error-correcting codes for 3, 4, and 5 information qubits with block lengths of 8, 10, and 11 qubits, respectively. [S1050-2947(96)07611-1]

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Two recent papers have shown that efficient quantum error-correcting codes exist [1,2]. A quantum error-correcting code is a method of storing or transmitting K bits of quantum information using $n > K$ quantum bits (qubits) [3], in such a way that if an arbitrary subset of the n qubits undergoes arbitrary errors, the transmitted quantum information can nevertheless be recovered exactly. The only condition is that the values of n and K , and the structure of the error-correcting code, place an upper limit on the number of qubits which can undergo errors before quantum information is irretrievably lost. In this context an ‘‘error’’ is any physical process which influences the quantum information but whose effect cannot be ‘‘undone’’ simply by applying, upon reception, a time-reversed version of the ‘‘error’’ process. In practice this will be because the errors are unpredictable (e.g., caused by unknown stray fields) or they entangle the information-bearing system with another system which is not accessible to detailed manipulation (e.g., the environment). The latter case, entanglement with the environment, includes as a subset relaxation processes, such as, spontaneous emission and phase decoherence.

A simple such error-correcting code was presented in [4], and a general method of encoding a single qubit with correction of multiple errors was presented in [5]. More importantly, further work [1,2] derived whole classes of codes, for multiple correction of many qubits, and showed that *efficient* codes exist. The word ‘‘efficient’’ in this context refers to the fact that the ratio K/n , which is called the rate of the code, need not become smaller and smaller as K increases, for a given probability of error per qubit. However, whereas the pioneering works just mentioned established the possibility of efficient quantum error-correction, and presented the simplest codes possible, they did not address the more pragmatic issue of identifying other specific useful codes. This is the subject addressed in this paper.

It is important to identify codes for more than a single qubit, since it is known that codes involving more information can be more powerful than simple repetition of single-qubit codes. In comparing two coding techniques, one may appear more powerful because it can encode a single qubit more efficiently. However, if it cannot also be applied to many qubits in an efficient way, for example, because find-

ing a good code is too difficult, then a simpler technique for which multiple-qubit codes can be found may end up being the better choice.

The earlier simple codes have now been improved upon with the discovery [6,7] of a ‘‘perfect’’ code, that is, one which fills a lower bound (elucidated below) on the number of qubits n required to do the job. This code is more efficient than those constructed by the recipe of [1] and [2], and introduces an important class of codes. However, it is not yet known how to generalize the construction method in order to obtain other efficient codes, so once again the task of identifying specific examples of useful codes is an important one. These more efficient codes are also discussed here. Simple methods to find good codes are described, and three examples presented.

Note that the errors which we wish to correct are completely random, and we have no knowledge of their nature other than that they affect different qubits independently. If we are in possession of further information about an error process, this can be used to construct codes which are more resistant to the errors caused by that process [8,9]. An example is when the dominant error process is spontaneous emission. In this case the ‘‘error’’ process is in fact almost completely known, but causes an unavoidable coupling to the environment. Efficient coding for this situation has recently been considered [10,11]. In the present work we make the standard assumption that the only predictable feature of the errors is their random nature, so we wish the code to correct as many arbitrary errors as possible using as few qubits as possible.

Arbitrary errors of qubits can usefully be divided into ‘‘amplitude errors,’’ that is, changes of the form $|0\rangle \leftrightarrow |1\rangle$, and ‘‘phase errors,’’ that is, changes of the form $|0\rangle + |1\rangle \leftrightarrow |0\rangle - |1\rangle$. This division is not meant to imply that these simple state rotations in Hilbert space are the only form of error considered, but rather a completely general error can be described as a combination of such amplitude and phase errors, with associated entanglement with the environment. Less obviously, but importantly, a method which can correct both amplitude and phase errors is sufficient to correct general errors [1,2,12].

An essential result which relates the problem of error-correction coding of a quantum channel to that of a classical channel is embodied in theorems 3.1 and 3.3 of [2] and theorem 1 of [1]:

Theorem (“quantum correction theorem”): If C_1 and C_2 are both linear $[n, k, d]$ codes with $\{0\} \subseteq C_2^\perp \subseteq C_1 \subseteq F_2^n$, then the quantum code Q_{C_1, C_2} is a t -error-correcting code, where $t = \lfloor (d-1)/2 \rfloor$.

This statement is taken from [1]. Note, however, that we have defined k to be the dimension of C_1 rather than of its dual as in [1], and replaced C_2 by C_2^\perp , to make the reasoning more symmetric. The symbol F_2^n refers to the n -dimensional vector space over a binary field. The phrase “ t -error-correcting code” refers to the fact that this form of encoding allows correction after arbitrary errors of t qubits. This is stated and proved as a separate theorem (Theorem 3.3) in [2]. In [2], the equivalent statement of the quantum correction theorem is the following:

To encode K qubits with minimum distance d_1 in one basis, and minimum distance d_2 in the other, it is sufficient to find a linear code of minimum distance d_1 , whose K th order subcode is the dual of a distance d_2 code.

Since C_2^\perp has dimension $n-k$, and $C_2^\perp \subseteq C_1$, clearly C_2^\perp is a K th order subcode of C_1 , with $K = 2k - n$, which links the first statement to the second. In the second statement a more general form is considered, in which the minimum distances of C_1 and C_2 need not be equal. This has consequences for the type of error which can be corrected, as elucidated in [2]. In particular, if only phase errors are present, then d_1 can be 1, that is, C_1 is simply the set of all words, so the coding problem reduces to that of finding $C_2 = [n, K, d_2]$, which means it is equivalent to the classical coding problem. The same simplification applies when only amplitude errors are present, in which case C_2 is the set of all words and C_1 is the error correcting code.

The second statement above refers to two different bases in which the state of the information-bearing quantum system may be expressed. For a single qubit, basis 1 is $\{|0\rangle, |1\rangle\}$, and basis 2 is $\{(|0\rangle + |1\rangle)/\sqrt{2}, (|0\rangle - |1\rangle)/\sqrt{2}\}$. For multiple qubits, bases 1 and 2 are defined to be the respective product bases. The reason why these bases are introduced is that correction of amplitude errors is essentially a classical error correction operating in basis 1, while that of phase errors is essentially a classical error correction operating in basis 2 [2].

The rest of this paper is concerned with finding quantum error-correcting codes. We begin with codes obtained according to the recipe of the quantum correction theorem. In Sec. I we consider arbitrary d_1 and d_2 , for the case $K=0$. This is not a fruitless exercise since it will be shown that a code with $K>0$ can be obtained from one with $K=0$. In Sec. II the case $d_1 = d_2 \equiv d$ is considered, for arbitrary d and K . This case is important because a completely general error of less than $d/2$ qubits can be corrected by the use of such a code (quantum correction theorem above, and Theorem 6 of [2]). As well as specific codes, simple methods for finding codes, and for deducing one code from another are given.

Note that whereas the method of Secs. I and II will produce a set of useful codes, whose rate does not reduce as K is increased at constant error probability per qubit, it will not produce the most efficient codes possible. However, the codes may be regarded as a starting point from which more efficient codes can be derived by judicious use of sign

changes with code augmentation (adding of code words) or puncturing (deletion of bits). Such methods are discussed in Sec. III, where optimal or almost optimal quantum codes are presented for encoding 3 to 5 qubits with single-error correction.

Many powerful mathematical (group theoretical) techniques have been applied in the pursuit of classical coding theory [8], and since completing the present work I have learned of two studies which apply such techniques with much success to quantum coding [15,16]. In this paper, the approach is to use simple concepts such as Hamming distance, parity check, and generator matrices, and examine methods to convert classical codes into good quantum codes. The simplicity of these concepts has the advantage of being suggestive of useful coding techniques, since they make the structure of the codes simple to appreciate, but they do not always lead to analytical proofs of the properties of the quantum codes, for which one must resort to computer testing. The two approaches of trying simple ideas and applying powerful analytical methods are both useful in the quest to find good codes. The matrix methods used here have the further advantage that the quantum networks for encoding and correction can be derived quite straightforwardly from the parity check and generator matrices, by generalizing the method described in [2].

To distinguish the various types of code, it is helpful to have a concise notation. Classical linear error-correcting codes are identified by the notation $[n, k, d]$, meaning a code by which n classical bits can store k bits of classical information with minimum distance d , hence allowing correction of up to $\lfloor (d-1)/2 \rfloor$ errors. The expression $\lfloor x \rfloor$ denotes the largest integer less than or equal to x . The notation $\{n, K, d_1, d_2\}$ is here introduced to identify a “quantum code,” meaning a code by which n quantum bits can store K bits of quantum information and allow correction of up to $\lfloor (d_1-1)/2 \rfloor$ amplitude errors, and simultaneously up to $\lfloor (d_2-1)/2 \rfloor$ phase errors. For codes with $d_1 = d_2 \equiv d$ the notation will be abbreviated to $\{n, K, d\} \equiv \{n, K, d, d\}$. Such codes allow recovery after arbitrary error of up to $\lfloor (d-1)/2 \rfloor$ of the quantum bits. It may be argued that the Hamming distance d is no longer a useful term in the quantum context, since it is not clear whether $2t+1$ always corresponds to a quantity with the correct properties to be called a “distance” between code vectors of a quantum t -error-correcting code. However, I retain the use of d , both because it implies the distinction between error detection and correction, and because the concept of distance remains useful in searching for quantum codes, as will be shown in Sec. III.

The recipe $C_1 = [n, k, d_1]; C_2 = [n, k, d_2]; C_2^\perp \subseteq C_1$ of the quantum correction theorem leads to a code construction in which each code vector (i.e., encoded version of a given logical symbol) consists of a superposition of words with coefficients equal in sign as well as magnitude in one of the bases 1 or 2 (though not in the other). One may therefore choose the sign of all coefficients in the superposition to be positive, in the chosen basis. A code having this special form (i.e., all those discussed in Secs. I and II) will be indicated by appending a superscript “+” sign to the notation, i.e., $\{n, K, d_1, d_2\}^+$. In general, by allowing more general code vectors, a code having the same correction ability but higher rate can be obtained from one with all-positive signs in the original basis. In symbols,

TABLE I. Lower bound on n permitting a dual pair of codes of distances d and d^\perp . The underlined figures (the column $d=3$ and the row $d^\perp=7$) indicate that the bound is sharp and the code is given below. The asterisks indicate the following possibilities: a reduced Golay code can be used to obtain $d=d^\perp=5$ with $n=18$. The $[31,16,5]$ BCH code has $d^\perp=12$, so $d,d^\perp=5,11$ is possible with $n=30$. There are quadratic residue self-dual codes with parameters $[48,24,12]$ and $[80,40,16]$ so $d=d^\perp=11, 15$ is possible with $n=46, 78$, respectively [8].

n	d						
	3	5	7	9	11	13	15
3	<u>6</u>						
5	<u>11</u>	16*					
7	<u>14</u>	<u>20</u>	<u>22</u>				
9	<u>20</u>	25	30	34			
11	<u>23</u>	28*	33	39	42*		
13	<u>27</u>	33	38	43	46	52	
15	<u>30</u>	37	42	47	51	56	60*

$$\{n, K, d_1, d_2\}^+ \Rightarrow (\{n' < n, K, d_1, d_2\} \text{ and/or } \{n, K' > K, d_1, d_2\}), \quad (1)$$

where the implication sign is used to mean that once the left-hand side code is known, the right-hand side code can be obtained easily.

I. ZERO INFORMATION QUBITS

This section will consider $\{n, 0, d_1, d_2\}^+$ codes. If d_1 is the minimum distance of a classical linear code C , then by the quantum correction theorem (second statement above), d_2 is just the minimum distance of the dual code C^\perp , when $K=0$. In symbols,

$$\{n, k, d\}, d^\perp \equiv \{n, 0, d, d^\perp\}^+. \quad (2)$$

TABLE II. Properties of codes making up Table I. The size of the code is k , that of the dual is $n-k$. Where the code is identified as a sum of two or more, the first code in the sum is extended by the others in a manner explained in the text. Other codes may be possible, having the same $\{n, k, d, d^\perp\}^+$ but a different structure. However, there are no linear codes of smaller n for the same d, d^\perp , with the exception of the final entry: two-error-correcting BCH codes are not necessarily optimal. They are included here because they are close to optimal and easily constructed.

n	k	d	d^\perp	code C
2	1	2	2	Repetition
6	3	3	3	Hamming
11	4	3	5	$[8,4,4]$ extended Hamming + $[3,3,1]$
14	10	3	7	Hamming
20	15	3	9	$[16,11,4]$ extended Hamming + $[4,4,1]$
23	18	3	11	$[16,11,4]$ + $[7,4,3]$ Hamming
27	22	3	13	$[16,11,4]$ + $[8,4,4]$ + $[3,3,1]$
30	25	3	15	Hamming
24	12	8	8	Golay
n	1	n	2	Repetition \leftrightarrow even weight
$2^r - 1$	$n - r$	3	2^{r-1}	Hamming \leftrightarrow Simplex
$2^r - 1$	$n - 2r$	5	$2^{r-1} - 2^{\lfloor r/2 \rfloor}$	BCH \leftrightarrow BCH $^\perp$

Helgert and Stinaff [13] have prepared a table of the minimum distance d of linear codes $[n, k, d]$ for given n and k . Specifically, the interesting quantity is the highest $d = d_{\max}(n, k)$ permitted for the given values of n and k . If $d_{\max}(n, k)$ is not known then Helgert and Stinaff give upper and lower bounds on it. For brevity, Helgert and Stinaff's table will be referred to as HS. It is possible to convert such a table into one providing a lower bound on the smallest number of bits $n = n_{\min}(d, d^\perp)$ necessary in order that a code can have distance d and its dual have distance d^\perp . For a given d , one commences with $n = d$, which gives a code $[d, 1, d], d^\perp = 2$. To allow larger values of d^\perp , n must be increased, and k set to the largest value allowing an $[n, k, d]$ code, as indicated by HS. The values of n and k are increased together in this way until HS indicates that an $[n, n-k, d^\perp]$ code is possible. Clearly, there is no code with n smaller than the value thus obtained, for which both $[n, k, d]$ and $[n, n-k, d^\perp]$ codes are possible. This does not prove, however, that an $[n, k, d]$ code exists whose dual has distance d^\perp . A necessary but not sufficient existence condition is established, or in other words, a lower bound on the value of $n_{\min}(d, d^\perp)$. This lower bound is given in Table I.

To find out whether the lower bound in Table I is sharp, I have attempted to identify codes which satisfy the bound. Success at identifying such a code is indicated by an underlined n value in Table I, and the code identified is described in Table II. An asterisk in Table I indicates that a code with n close to the lower bound exists and is identified in the caption.

In Table II, the identification $[n, k, d] + [n', k', d']$ refers to a code built by combining two others as follows. To the check matrix of the first code ($[n, k, d]$) in the sum, additional columns are added as specified by the generator matrix of the second code ($[n', k', d']$) in the sum. This lengthens the minimum distance of the dual by d' while increasing n by n' and reducing d . For example, the code identified as

“ $[16,11,4]$ extended Hamming $+ [4,4,1]$ ” is the $[20,15,3]$ code with the following check matrix:

$$H_1 = \begin{pmatrix} 111111111111110000 \\ 101010101010101000 \\ 01100110011001100100 \\ 00011110000111100010 \\ 0000000111111100001 \end{pmatrix} \quad (3)$$

Its dual has minimum distance $8+1=9$.

One can “navigate” around Table I to some extent by use of the following two constructions:

$$[n,k,d], d^\perp \Rightarrow \begin{cases} [n-1, k-1, d], & d^\perp - 1 \\ [n-1, k, d-1], & d^\perp. \end{cases} \quad (4)$$

In these two constructions, the code on the right-hand side is derived from the $C=[n,k,d]$ code on the left-hand side by removing a single row from the generator ($k \rightarrow k-1$) or parity check ($d \rightarrow d-1$) matrix. To see how the minimum distance of the dual code is affected, recall that the generator matrix of C is the parity check matrix of C^\perp , therefore deleting a row from the generator matrix of C means deleting a row from the check matrix of C^\perp , and vice versa. A single row deleted from a generator matrix leaves the minimum distance either unaffected (the most likely result) or increased. A single row deleted from a check matrix leaves the minimum distance either reduced by one (the most likely result) or unaffected.

II. K INFORMATION QUBITS

The case $K \neq 0$ will now be addressed. The simplest case to consider is that of a classical code which contains its own dual: $C^\perp \subseteq C = [n,k,d]$. This is only possible when $2k \geq n$. Such codes have been called “weakly self-dual” [8]. Since C^\perp is a subcode of C , clearly the quantum correction theorem can be satisfied with $d_1 = d_2 = d$ and $K = 2k - n$, since C^\perp is itself the subcode required by the theorem. In symbols,

$$C^\perp \subseteq C = [n,k,d] \Rightarrow C = \{n, 2k-n, d\}^+. \quad (5)$$

In such a case, the error corrector is the same in basis 1 and basis 2. An example is the Hamming code discussed in [2,5]. This result transforms the search for quantum $\{n,K,d\}^+$ codes to a large extent to a search for classical weakly self-dual codes. This was recognized in [1], where a proof was given that weakly self-dual codes exist which satisfy the Gilbert-Varshamov bound. However, there exist $\{n,K,d\}^+$ codes which cannot be derived from weakly self-dual codes (examples are given below), and these can be more efficient (higher K/n for given d/n) than the best weakly self-dual codes.

A code contains its dual if and only if all the rows of the parity check matrix satisfy all the parity checks [i.e., $\text{wt}(H_i \cdot H_j)$ is even, for all $i, j = 0, \dots, n-k-1$, where $\text{wt}(z)$ is the number of 1's in z]. This implies that when a single row is deleted from the parity check matrix, the resulting code again contains its dual. Using the second construction given in (4), combined with Eq. (5), one finds

$$C^\perp \subseteq C = \{n, K, d\}^+ \Rightarrow C'^\perp \subseteq C' = \{n-1, K+1, d-1\}^+. \quad (6)$$

This allows one to generate codes encoding more quantum information (having greater K) from ones of smaller K , at the expense of reduced d . Note that d is not *required* to fall by 1, but implication (6) states that d does not fall by more than 1 in this construction.

Next the following question will be addressed: we wish to encode K qubits with given $d = d_1 = d_2$. What is the necessary value of n ? The quantum correction theorem implies that if subcodes of an $[n,k,d]$ code are used, then $K = 2k - n$. In the case of single-error correction, i.e., $d = 3$, Hamming's construction implies

$$k \leq n - \lceil \log_2(n+1) \rceil, \quad (7)$$

therefore, for an $\{n, K, 3\}^+$ code,

$$K \leq n - 2 \lceil \log_2(n+1) \rceil. \quad (8)$$

When $n = 2^r - 1$ we have a perfect Hamming code, and for this case the code contains its dual. Therefore equality holds in (8), and $K = n - 2 \log_2(n+1)$. The smallest n allowing $d = 3$ for values of K in the range 1 to 16 is indicated in Table III.

For $K = 2$ Eq. (8) implies $n \geq 10$. In fact $n = 10$ is possible using the following code:

$$H_1 = \begin{pmatrix} 1011001000 \\ 0101100100 \\ 1010110010 \\ 0110010001 \end{pmatrix}, \quad H_2 = \begin{pmatrix} 1111001000 \\ 0111100100 \\ 1010010010 \\ 1110110001 \end{pmatrix}, \quad (9)$$

$$D = \begin{pmatrix} 0001001100 \\ 0000010011 \end{pmatrix}. \quad (10)$$

Here, H_1 and H_2 give the correctors in bases 1 and 2, respectively, and the generator works as follows. Let C_1 be the classical code of which H_1 is the check matrix. The two rows of D are the fourth and sixth rows of the generator G_1 of C_1 , which is obtained from the well-known relation

$$H_1 = (A | I_{n-k}) \Leftrightarrow G_1 = (I_k | A^T), \quad (11)$$

where I_j is the $j \times j$ identity matrix, and A is the rest of the check matrix. Adding these two extra checks to H_1 , we obtain the check matrix for a subcode C_2^\perp of C_1 . The four states (code vectors) in the quantum $\{10,2,3\}^+$ code are the subcode $|C_2^\perp\rangle$, whose generator is H_2 , and its three cosets $|C_2^\perp \oplus D_0\rangle, |C_2^\perp \oplus D_1\rangle, |C_2^\perp \oplus D_0 \oplus D_1\rangle$, where D_0 and D_1 are the rows of D (the letter D is chosen here for “displacement”). In symbols, one may write this generation procedure as

$$G = \begin{pmatrix} H_2 \\ D \end{pmatrix}. \quad (12)$$

This equation may be regarded as a summary of the quantum network which will encode the two qubits of information. Note that since the rows of D are members of the code C_1 ,

TABLE III. Upper bound on d for $\{n, K, d\}^+$ codes of small K and n . Entries which are identical to the one immediately above them (i.e., with n reduced by 1) are left blank, in order to bring out the pattern in the results. A pair of figures is given when the table of HS [13] indicates a range of distance values rather than a precise upper limit. The underlined values are produced by codes given in Table IV or obtained from them by the methods discussed in the text. For these codes the listed upper bound is thus shown to be obtainable. Some codes are obtained from one another by deleting a row of G_1 (moving upwards and to the left in the table), or by deleting a row of H_1 (moving upwards and to the right in the table). An asterisk (*) indicates a self-dual or weakly self-dual code.

d	K															
	1	2	3	4	5	6	7	8	9	10	11	12	13	14	15	16
1	<u>1</u> *															
2		<u>1</u> *														
3			<u>1</u> *													
4	<u>2</u>	<u>2</u> *		<u>1</u> *												
5		<u>2</u>			<u>1</u> *											
6			<u>2</u>	<u>2</u> *		<u>1</u> *										
7	<u>3</u> *			<u>2</u>			<u>1</u> *									
8*					<u>2</u>	<u>2</u> *		<u>1</u> *								
9	<u>3</u>					<u>2</u>			<u>1</u> *							
10		<u>3</u>					<u>2</u>	<u>2</u> *		<u>1</u> *						
11	4		3					<u>2</u>			<u>1</u> *					
12		4	<u>3</u>	3					<u>2</u>	<u>2</u> *		<u>1</u> *				
13			<u>4</u>		<u>3</u>					<u>2</u>			<u>1</u> *			
14				4		<u>3</u>					<u>2</u>	<u>2</u> *		<u>1</u> *		
15					4		<u>3</u> *					<u>2</u>			<u>1</u> *	
16		<u>4</u>				<u>4</u> *						<u>2</u>	<u>2</u> *		<u>1</u> *	
17	5					<u>4</u>	<u>3</u>						<u>2</u>			
18				<u>4</u>		4		3							<u>2</u>	<u>2</u> *
19	<u>5</u>						4	<u>3</u>	3							<u>2</u>
20		<u>5</u>		<u>4</u> *				<u>4</u>	<u>3</u>	3						
21	<u>6</u>		<u>5</u> *						<u>4</u>	<u>3</u>	3					
22		<u>6</u> *		5						<u>4</u>	<u>3</u>	3				
23	<u>7</u> *		6		5						<u>4</u>	<u>3</u>	3			
24*				6		45						4	<u>3</u>	3		
25					56		45						4	<u>3</u>	3	
26				6		56		45						4	<u>3</u>	3
27	78				6		56		45						4	<u>3</u>
28		67				6		56		45						4
29	<u>6</u> 8		67				6		56		<u>4</u> *5					
30		<u>6</u> *8		67		<u>5</u> *6		6		56		<u>4</u> *				
31	<u>7</u> *8			68			67		6		<u>5</u> *		4			

they satisfy all the checks of H_1 , and so the cosets they generate are all subsets of C_1 . Also, since the rows of D have odd weight, the coset $|C_2^{\perp} \oplus D_i\rangle$ fails the parity check D_i , so the cosets are distinct. In general D need not have rows of odd weight. The nonoverlapping of the cosets is ensured by the fact that C_2 is not a zero-distance code.

The matrix formed by H_1 plus the extra rows given by D is the generator of C_2 , and the corrector in basis 2, H_2 , is obtained from this generator using relation (11). All these relationships may be summarized as follows:

$$H_1 \leftrightarrow G_1 \rightarrow D, \tag{13}$$

$$\left(\frac{H_1}{D}\right) \leftrightarrow H_2, \tag{14}$$

$$H_1 \leftrightarrow \left(\frac{H_2}{D}\right) = G. \tag{15}$$

From this one may see that an equivalent code is obtained by using H_1 and D as the generator in Eq. (12) instead of H_2 and D . A further equivalent quantum code can be obtained by using the first two rows of G_1 for D , instead of the fourth and sixth rows [cf. Eqs. (16) and (25) below].

The above approach can clearly be applied to any classical $[n, k, d]$ code. That is, one produces a subcode by using $2k - n$ words from the code as extra parity checks, with the aim that the check matrix thus obtained is the generator of another (or the same) $[n, k, d]$ code. However, it is not clear whether this method can always succeed in producing a useful quantum code. For example, whereas the cyclic check

TABLE IV. Example $\{n, K, d\}^+$ codes. The final entries give some assorted values of n and K larger than those covered by Table III.

n	K	d	code C^{+K}
4	2	2	
n	$n-2$	2	Even weight, for even n
2^r-1	$n-2 \log_2(n+1)$	3	Hamming, $r>2$
eg. 7	1	3	
15	7	3	
31	21	3	
2^r-1	$n-4 \log_2(n+1)$	5	BCH, $r>4$
eg. 31	11	5	
29	11	4	reduced BCH, $C^\perp \subseteq C$
2^r-1	$n-2t \log_2(n+1)$	$2t+1$	t -error correcting BCH
eg. 31	1	7	BCH
30	6	5	delete three rows from H of previous entry
16	6	4	extended Hamming
17	7	3	see text
19-27	8-16	3	cyclic, see text
23	1	7	Golay
48	0	12	quadratic residue (self-dual)
63	3	11	BCH
63	15	9	BCH
80	0	16	quadratic residue (self-dual)
104	0	20	quadratic residue (self-dual)
127	15	17	BCH

matrix H_1 of Eqs. (9) leads quickly to a quantum code, the Hamming check matrix for the same parameters [10,6,3] does not lead to a generator in the form given by Eq. (11) whose rows can be used to form D . It would be interesting to try to prove or disprove the hypothesis that the existence of a classical $[n, k, d]$ code is sufficient to imply at least the existence of a quantum $\{n, 2k-n, d\}^+$ code. The author's current impression is that this hypothesis is untrue in general. However, it is true for weakly self-dual codes, and probably gives a close estimate of the parameter values possible for other cases.

The single-error-correcting codes indicated in Table III were all obtained by using the above method of using code words as extra parity checks, but note that whereas I have thus found single-error correcting $\{13,5,3\}^+$ and $\{14,6,3\}^+$ codes, filling the lower limit on n set by the Hamming bound, I have not found $\{11,3,3\}^+$ or $\{12,4,3\}^+$ codes even though classical [11,7,3] and [12,8,3] codes exist. These single-error-correcting quantum codes are all obtained from the cyclic code given by the irreducible primitive polynomial $x^4=1+x$. The check matrix H_1 in Eqs. (9) is the check matrix of this cyclic code for the case $n=10$, and for higher n , up to $n=15$, further columns are added to the front of the matrix following the standard procedure. Once we have H_1 , the quantum code is fully defined once the relevant displacement matrix D is given. For $n=12$ to 14 the following matrices D fulfill the requirements for single-error correction:

$$D_{\{12,3,3\}^+} = \begin{pmatrix} 001000001010 \\ 000100000101 \\ 000010001011 \end{pmatrix}, \quad (16)$$

$$D_{\{13,5,3\}^+} = \begin{pmatrix} 1000000001111 \\ 0100000001110 \\ 0010000000111 \\ 0001000001010 \\ 0000000010011 \end{pmatrix}, \quad (17)$$

$$D_{\{14,6,3\}^+} = \begin{pmatrix} 1000000001101 \\ 0100000001111 \\ 00100000001110 \\ 0001000000111 \\ 00001000001010 \\ 0000000010011 \end{pmatrix}. \quad (18)$$

Note that the cases $n=13$ and $n=14$ are similar to one another, and can be obtained by reducing the $n=15$ code. For $n=15$ the code C_1 contains its own dual.

Using the above methods, and once again the table of HS, lists of quantum $\{n, K, d\}^+$ codes can be compiled. The results are summarized in Tables III and IV. The upper bound on d is found from the classical bound $d_{\max}[n, k=(K+n)/2]$ given by HS. As it stands, Table III is incomplete in that for most entries I have not found codes which realize the upper bound, thus proving that it is obtainable. However, classical self-dual codes supply efficient quantum codes of low rate, high distance (low K/n , high d), and Bose-Chaudhuri-Hocquengham (BCH) codes supply efficient quantum codes of high rate, low distance. Therefore

we have identified infinite series of codes, as n increases, at the two ends of the range $K=0$ to $\sim[n - \log_2(n)]$.

The $\{17,7,3\}^+$ code in Table III is specified by the following check matrix:

$$H_1 = \begin{pmatrix} 01100111100110000 \\ 10111100101101000 \\ 11010010111100100 \\ 11101001110000010 \\ 00011111110000001 \end{pmatrix} \quad (19)$$

with the D matrix equal to the last seven rows of G_1 . This code can be obtained by adding a check bit to the $[16,11,4]$ extended Hamming code.

The single-error-correcting codes indicated in Table III as certainly obtainable (i.e., which I have succeeded in finding) for $7 < K < 17$ do not realize the minimum implied by the results of HS, but require one additional qubit, similar to the cases $K=3$ and 4 already remarked. Using the clue mentioned above that a cyclic classical code rather than a Hamming code is a good choice, all the codes from $K=8$ to 16 were obtained from the cyclic classical codes of primitive polynomial $x^5=1+x^3$. In this series of codes, a classical $[n,k,d]$ code gives rise to a quantum $\{n,2k-n-1,d\}^+$ code. For example, the generator for $K=16$ is

$$G_{\{27,16,3\}^+} = \begin{pmatrix} 011001111100011011101010000 \\ 101100111110001101110101000 \\ 001111100011011101010000100 \\ 100111110001101110101000010 \\ 11001111000110111010100001 \\ 0010000000000000000001100 \\ 0000100000000000000000111 \\ 0000010000000000000001111 \\ 00000001000000000000011001 \\ 00000000010000000000001100 \\ 00000000001000000000001100 \\ 0000000000010000000000011 \\ 00000000000010000000010101 \\ 00000000000001000000011110 \\ 00000000000000010000001111 \\ 000000000000000001000010011 \\ 0000000000000000001000011101 \\ 0000000000000000000100011010 \\ 0000000000000000000010001101 \\ 0000000000000000000001010010 \\ 000000000000000000000101001 \end{pmatrix} \quad (20)$$

The generators for the other codes in this series have a similar form and will not be listed.

A $\{20,9,3\}^+$ code can also be obtained from the classical code having the minimum n for $d=3, d^\perp=9$, referred to in Tables I and II. Its check matrix H_1 is given in Eq. (3), and the displacement matrix is formed from rows of G_1 as follows:

$$D_{\{20,9,3\}^+} = \begin{pmatrix} 0100000000000010100 \\ 00000100000000010110 \\ 00000001000000010001 \\ 00000000010000010101 \\ 00000000001000011101 \\ 00000000000100010011 \\ 00000000000010011011 \\ 00000000000001010111 \\ 00000000000000111111 \end{pmatrix} \quad (21)$$

III. MORE EFFICIENT CODES

Implication (1) was used in the introduction to encapsulate the twin facts that in general $\{n,K,d\}^+$ codes are not the most efficient possible, and that they can be used as a starting point to obtain more efficient $\{n,K,d\}$ quantum codes. The simplest example is the perfect $\{5,1,3\}$ quantum code described in [6,7], which can be obtained by deleting any two bits from the $\{7,1,3\}^+$ code described in [5,2], and changing the signs of a subset of the words in each of the two code vectors. The relevant sign changes can be found for this simplest case by an exhaustive computer search. The computer search is a useful tool in the task of finding good codes, which may be likened to a search for the best fruit on a many-branched tree. However, a complete search of all possible allocations of signs rapidly becomes too time consuming, as the parameters $\{n,K,d\}$ are increased. Intelligent search techniques must be used, and barren branches of the tree ruled out as efficiently as possible, while fruitful branches must be identified before the search begins, which is the demanding task of the human researcher. In this section a set of quantum codes will be presented, all of which were found by taking advantage of two simple methods to identify fruitful branches and thus find suitable sign allocations quickly. Before discussing these sign allocations, however, we will consider ways of combining classical codes which go beyond the simple recipe $C_2^\perp \subseteq C_1$.

The generator matrix of a quantum code, Eq. (12), creates one classical code (forming the first quantum code vector) and $2^K - 1$ cosets (which form the remaining quantum code vectors). Thus we may picture the first quantum code vector as a lattice of points in an n -dimensional Hamming space, and the other code vectors as this lattice displaced around the Hamming space by distances of order d . The codes described in previous sections used lattices displaced so that each point in any given lattice was at least a distance d from any point in another lattice. In other words, the set of all the lattices formed a classical code of distance d , and this ensured that error correction was possible in basis 1. However, in forming a quantum code, it is not necessary to displace the lattices as far as this. Bit flips in basis 1, i.e., amplitude errors, will cause a given lattice to move towards some other lattice, i.e.,

the code vectors approach, but if we now allow the signs in basis 1 to be negative as well as positive, then lattices (i.e., cosets) which overlap, in that they contain the same sets of words, may nevertheless correspond to orthogonal quantum states since there is an equal number of positive and negative contributions to the inner product $\langle i, e_k | j, e_l \rangle$, where $|i, e_k\rangle, |j, e_l\rangle$ are code vectors affected by errors e_k, e_l . Thus if we start from a set of code vectors with all-positive signs when written in basis 1, then the introduction of sign changes permits the distance between cosets in basis 1 to be reduced.

Clearly, we must not hope for too much from this ability to allow the cosets to approach. The minimum assumption is that we may permit the distance between cosets in basis 1 to be reduced by one. In other words, the recipe for a $\{n, K, d\}$ quantum code becomes $C_2^\perp \subseteq C_1$ where C_2 is a distance d classical code as before, but now C_1 is a distance $d-1$ code. In addition, we wish C_2^\perp to have as large a minimum distance as possible, in order to allow a lot of ‘‘room’’ to move the lattice around in Hamming space before it overlaps itself. These two conditions, together with a judicious application of sign changes, will be used to find optimal single-error-correcting quantum codes. First, however, we must consider how to apply sign changes to the words in the code vectors.

The first method to allocate sign changes to the words in each code vector is to restrict the possible sign allocations to those given by rows of the Hadamard matrix [8]. That is to say, we use the $w \times w$ Hadamard matrix to supply w different allocations of w signs. A sign allocation is a set of $w+1$'s and -1 's, giving the signs of each of the w words in the superposition forming the code vector to be tested. The Hadamard matrix can be used in this way since in all the codes considered here, each code vector contains a number of words w equal to a power of 2. It is not hard to convince oneself that a row of the Hadamard matrix is an intelligent choice of sign allocation for any quantum code derived from a linear classical code by the methods discussed in previous sections (with possibly one or more bits deleted).

Once we have a code vector, that is, a set of words with a proposed sign allocation, it is tested. The test consists of first testing whether errors in the code vector lead to states orthogonal to the code vector itself and to each other, and then testing whether such erroneous states are also orthogonal to all the other code vectors in the code and their erroneous versions. The possible errors included in the test are all those which the code is supposed to be able to correct. If a code survives such a test, then errors of different syndrome lead to orthogonal states, and the orthogonality of different code vectors is also preserved. Such an ‘‘orthogonal coding’’ implies that error correction is certainly possible. This latter fact is a central part of the argument presented in [1] and [2]. An elegant presentation of it is also provided by Ekert and Macchiavello [12], which enables the latter authors to deduce a quantum version of the Hamming bound, based on counting the number of possible orthogonal directions in Hilbert space, see also [9,7,16]. Their bound is

$$2^K \sum_{i=0}^t 3^i \binom{n}{i} \leq 2^n \quad (22)$$

for the size K of a possible quantum code which can correct

t general errors using n qubits. This bound is more general than that required for $\{n, K, 2t+1\}^+$ codes derived in [2], since it includes the possibility of the more general $\{n, K, 2t+1\}$ codes which we are considering in this section. However, there is an intriguing possibility that it is not strictly necessary for all possible error syndromes of all possible code vectors to be associated with mutually orthogonal states, since some error correction techniques may be able to correct errors of different syndrome without needing to distinguish the syndromes explicitly. This is the subject of active research [7,9,14,15] and will not be addressed in the present work. In other words, we use ‘‘orthogonal coding’’ throughout.

The second of our two methods to identify fruitful trial codes is to consider the sign allocation, that is the row of $w+1$'s and -1 's, as itself a binary vector of length w , and then to use linear combinations of such vectors in a fashion to be explained shortly. To keep the notation concise, we replace $+1$ in the sign allocation vector by 0, and -1 by 1, to get a vector in the usual binary form, but one which is understood to represent a sign allocation among w superposed words. Note that this vector has length w , which is a power of two and usually larger than the length n of the words in the superposition forming a code vector. For example, the sign allocation for the two code vectors of the five-qubit code of [6] can be written

$$00010100, 01110010, \quad (23)$$

where the least significant (i.e., rightmost) bit in the sign vector gives the sign of the first word in the code vector, and we assume the order of the words in the code vectors is that obtained when they are generated using the generator matrix

$$G = \begin{pmatrix} 10101 \\ 10011 \\ \hline 01111 \\ 11111 \end{pmatrix} \quad (24)$$

(cf. Eq. 12 for the notation). Note that the signs in (23) are *not* rows of the Hadamard matrix, showing that the Hadamard method will not pick up all good codes. However, both sign vectors in (23) are offset from rows of the Hadamard matrix by the same code vector 00010100, so (23) is a coset of a sign allocation obtained by the Hadamard method (i.e., a coset of a subset of a first-order Reed-Muller code).

The second of our two methods to allocate signs only applies to codes of more than one encoded qubit, i.e., having more than two code vectors. The method is to let the set of 2^K w -bit sign vectors itself be a classical linear code (or a coset of a linear code if necessary), and to allocate each sign vector thus generated to the corresponding code vector generated by G . For example, once we have found sign vectors s_{00}, s_{01}, s_{10} for the first three code vectors of an $\{n, 2, d\}$ code, we try the sign vector $s_{11} = s_{00} \oplus s_{01} \oplus s_{10}$ for the fourth code vector. By this process, we only need K sign vectors (plus possibly one more to form a coset) to specify all the signs for an $\{n, K, d\}$ code, rather than finding 2^K independent vectors which is a much more demanding task.

The quantum Hamming bound (22) states that for single-error correction ($t=1, d=3$), at least $n=5,7,8,9,10$ qubits are required to encode $K=1,2,3,4,5$ qubits, respectively, and $n=10$ qubits are required to correct $K=1$ qubit with double-error correction ($t=2, d=5$). The $n=5$ case is a perfect code since it fills the bound, and is that discussed in [6,7]. The next most simple case is $n=7$, for which we search for an encoding of two qubits with single-error correction. I have not found such a $\{7,2,3\}$ code, despite a wide but not complete search (this search was not restricted to the two methods just discussed). The best codes I have found are ones which encode two qubits using 7 but for which the third and fourth code vectors are not quite compatible with the first and second. That is to say, there are ten cases in which a single-qubit error in one code vector leads to the same quantum state as a different single-qubit error in another code vector, causing an ambiguity for any error corrector. These ten cases are taken out of the $88^2/2=3872$ possible comparisons between one code vector with its erroneous versions and another code vector with its erroneous versions, so the code comes close to single-error correction, while not realizing it completely.

An encoding of three qubits permitting complete single-error correction can be obtained with $n=8$, which is optimal in that this is the lower limit given by (22). To find the code, we begin with a classical code C_2 having a minimum distance of at least 3, to allow correction of errors in basis 2 (phase errors) and having a dual C_2^\perp of minimum distance as large as possible, since this dual code defines the lattice in basis 1 whose various displaced versions constitute the code vectors in basis 1. The $[8,4,3]$ Hamming code is not a good choice since its dual has a minimum distance of only 1.

Instead we adopt the extended Hamming code or Reed-Muller code $C_2=[8,4,4]$, which is self-dual so C_2^\perp has a minimum distance of 4. Since we want $d=3$, we allow C_1 to have distance $d-1=2$. This suggests the even-weight $[8,7,2]$ code, which has the correct number of code vectors to allow $K=3$. Thus we obtain the following generator:

$$G_{\{8,2,3\}} = \begin{pmatrix} 01010101 \\ 00110011 \\ 00001111 \\ \underline{11111111} \\ 11000000 \\ 10100000 \\ 10001000 \end{pmatrix}. \quad (25)$$

The sign vectors are found by computer search using the two shortcuts described above, which leads to eight sign vectors generated by

$$S_{\{8,3,3\}} = \begin{pmatrix} 0011001100110011 \\ 0000111100001111 \\ 0110011001100110 \end{pmatrix} \equiv \begin{pmatrix} 3333 \\ 0F0F \\ 6666 \end{pmatrix}, \quad (26)$$

where the second version is the first written in hexadecimal to bring out the structure.

Equations (25) and (26) are quite concise and combine several notations introduced in this paper. To make sure the notation is correctly understood, the $\{8,2,3\}$ code defined by these equations is now written out in full:

$$\begin{aligned} |v^{000}\rangle = & +|00000000\rangle + |01010101\rangle + |00110011\rangle + |01100110\rangle \\ & + |00001111\rangle + |01011010\rangle + |00111100\rangle + |01101001\rangle \\ & + |11111111\rangle + |10101010\rangle + |11001100\rangle + |10011001\rangle \\ & + |11110000\rangle + |10100101\rangle + |11000011\rangle + |10010110\rangle \end{aligned} \quad (27)$$

$$\begin{aligned} |v^{001}\rangle = & -|11000000\rangle - |10010101\rangle + |11110011\rangle + |10100110\rangle \\ & -|11001111\rangle - |10011010\rangle + |11111100\rangle + |10101001\rangle \\ & -|00111111\rangle - |01101010\rangle + |00001100\rangle + |01011001\rangle \\ & -|00110000\rangle - |01100101\rangle + |00000011\rangle + |01010110\rangle \end{aligned} \quad (28)$$

$$\begin{aligned} |v^{010}\rangle = & -|10100000\rangle - |11110101\rangle - |10010011\rangle - |11000110\rangle \\ & + |10101111\rangle + |11111010\rangle + |10011100\rangle + |11001001\rangle \\ & -|01011111\rangle - |00001010\rangle - |01101100\rangle - |00111001\rangle \\ & + |01010000\rangle + |00000101\rangle + |01100011\rangle + |00110110\rangle \end{aligned} \quad (29)$$

$$\begin{aligned} |v^{011}\rangle = & +|01100000\rangle + |00110101\rangle - |01010011\rangle - |00000110\rangle \\ & -|01101111\rangle - |00111010\rangle + |01011100\rangle + |00001001\rangle \\ & +|10011111\rangle + |11001010\rangle - |10101100\rangle - |11111001\rangle \\ & -|10010000\rangle - |11000101\rangle + |10100011\rangle + |11110110\rangle \end{aligned} \quad (30)$$

$$|v100\rangle = \begin{aligned} & +|10001000\rangle -|11011101\rangle -|10111011\rangle +|11101110\rangle \\ & +|10000111\rangle -|11010010\rangle -|10110100\rangle +|11100001\rangle \\ & +|01110111\rangle -|00100010\rangle -|01000100\rangle +|00010001\rangle, \\ & +|01111000\rangle -|00101101\rangle -|01001011\rangle +|00011110\rangle \end{aligned} \tag{31}$$

$$|v101\rangle = \begin{aligned} & -|01001000\rangle +|00011101\rangle -|01111011\rangle +|00101110\rangle \\ & -|01000111\rangle +|00010010\rangle -|01110100\rangle +|00100001\rangle \\ & -|10110111\rangle +|11100010\rangle -|10000100\rangle +|11010001\rangle, \\ & -|10111000\rangle +|11101101\rangle -|10001011\rangle +|11011110\rangle \end{aligned} \tag{32}$$

$$|v110\rangle = \begin{aligned} & -|00101000\rangle +|01111101\rangle +|00011011\rangle -|01001110\rangle \\ & +|00100111\rangle -|01110010\rangle -|00010100\rangle +|01000001\rangle \\ & -|11010111\rangle +|10000010\rangle +|11100100\rangle -|10110001\rangle, \\ & +|11011000\rangle -|10001101\rangle -|11101011\rangle +|10111110\rangle \end{aligned} \tag{33}$$

$$|v111\rangle = \begin{aligned} & +|11101000\rangle -|10111101\rangle +|11011011\rangle -|10001110\rangle \\ & -|11100111\rangle +|10110010\rangle -|11010100\rangle +|10000001\rangle \\ & +|00010111\rangle -|01000010\rangle +|00100100\rangle -|01110001\rangle, \\ & -|00011000\rangle +|01001101\rangle -|00101011\rangle +|01111110\rangle \end{aligned} \tag{34}$$

This code has also recently been derived by Gottesman [16]. He presents a general construction for $\{2^r, 2^r - r - 2, 3\}$ codes. These parameters are consistent with the supposition that such codes are obtained from the above method applied to the classical pair $C_2^\perp =$ first order Reed-Muller $[2^r, r+1, 2^{r-1}]$ code, $C_1 =$ even weight $[2^r, 2^r - 1, 2]$ code.

Proceeding to the encoding of four qubits, the bound (22) implies that single error correction is possible with $n=9$. However, $n=9$ is not large enough to allow a significant improvement on the properties of the $[8,4,4]$ classical code, so it seems unlikely that $\{9,4,3\}$ is possible, and I have not been able to find such a code. With $n=10$, on the other hand, we can adopt the 11-bit code indicated in Tables I and II which allows $d=3, d^\perp=5$, reducing it by the first construction in (4) to obtain $C_2 = [10,6,3]$, $C_2^\perp = [10,4,4]$. This leads to the following quantum code:

$$G_{\{10,4,3\}} = \begin{pmatrix} 0101010110 \\ 0011001101 \\ 0000111100 \\ 1111111100 \\ 1100000000 \\ 1010000000 \\ 1000100000 \\ 0000000011 \end{pmatrix}, \tag{35}$$

$$S_{\{10,4,3\}} = \begin{pmatrix} 0011001100110011 \\ 0000111100001111 \\ 0101010110101010 \\ 0000111111110000 \end{pmatrix} \equiv \begin{pmatrix} 3333 \\ 0F0F \\ 55AA \\ 0FF0 \end{pmatrix}. \tag{36}$$

Calderbank *et al.* [15] have also obtained a code of these parameters. The quantum Hamming bound (22) does not rule out the possibility of a further information qubit without increasing n , i.e., $\{10,5,3\}$, but I have been unsuccessful in finding such a code.

To encode five qubits with single-error correction, the classical 11-bit code with $d=3, d^\perp=5$ just mentioned can be used to obtain the following quantum code:

$$G_{\{11,5,3\}} = \begin{pmatrix} 01010101100 \\ 00110011010 \\ 00001111001 \\ 11111111000 \\ 11000000000 \\ 10100000000 \\ 10001000000 \\ 00000000110 \\ 00000000101 \end{pmatrix}, \quad S_{\{11,5,3\}} = \begin{pmatrix} 3333 \\ 0F0F \\ 55AA \\ 0FF0 \\ 3333 \end{pmatrix}. \tag{37}$$

This is optimal for five information qubits and single-error correction if $\{10,5,3\}$ does not exist.

IV. CONCLUSION

Much research is currently directed to finding the most efficient quantum error-correction techniques. Commonly in these efforts only the simplest example code, encoding a single qubit of information, is actually identified. However, to convey many bits of information, it is known from classical theory that more advanced codes, involving many infor-

mation bits, are more powerful than a repetition of single-information-bit codes. This implies that a more efficient coding technique is not useful unless the task of applying it to many information bits is mathematically tractable. A simpler coding technique, which is less efficient than other methods for one qubit, may become more efficient than the competing methods when many qubits are involved, simply because powerful many-qubit codes can be identified for the simpler method but not for its competitors.

In this paper many examples have been given of quantum error-correcting codes of reasonably high efficiency. In the process, several simple techniques for manipulating codes and guessing additional ones have been described. Quantum networks to encode and correct each code have not been given, since they can be deduced directly from the relevant generator and parity check matrices [2] [Theorem 3.1].

Starting with the simplest general method of quantum error correction, based on dual pairs of classical linear codes, and specified here by the notation $\{n, K, d\}^+$, we have tabulated codes which can be lifted almost directly from classical coding theory because they are self-dual or weakly self-dual. In addition, classical dual code pairs with maximal d, d^\perp have been tabulated, since they form a useful starting point for finding quantum codes. We have then examined in more detail the case of single-error correction, obtaining many good quantum codes from classical codes which are not weakly self-dual, and whose conversion to the quantum case therefore requires more ingenuity. A method using a subset of the rows of the generator matrix as extra parity checks has been described, and used to find quantum codes of param-

eters $\{10,2,3\}^+, \{12,3,3\}^+, \{13,5,3\}^+, \{14,6,3\}^+, \{17,7,3\}^+, \{19 \cdots 27, 8 \cdots 16, 3\}^+$.

Next, we have improved on the $\{n, K, d\}^+$ codes by allowing one of the classical codes (C_1) used to generate the quantum code to have its minimum distance reduced, and compensating for this by allowing the signs of words in the code vectors to be either positive or negative in all bases. To find out how to allocate the signs in this case, it is necessary to use insight rather than trial and error. By making the sign allocations themselves form a classical linear code, and by using the Hadamard matrix to supply useful sets of signs, we have introduced further structure into the quantum code. As well as making the design of generator and corrector networks easier, this allows the set of possible sign allocations to be vastly restricted, which greatly aids the search for good codes. These methods have enabled us to identify single-error correcting $\{n, K, 3\}$ quantum codes of $n=8, 10, 11$, $K=3, 4, 5$ qubits, respectively. The first is optimal, and it is possible that the others are also, though $\{10, 5, 3\}$ and $\{11, 6, 3\}$ are not ruled out by the quantum Hamming bound.

Developing general methods for producing good error-correcting codes is notoriously difficult. The task of finding good codes can be framed as a computational problem. It may be an example of a practically important computation whose solution on an ideal quantum computer is more efficient (has lower computational complexity) than any algorithm for a classical computer.

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- [1] A. R. Calderbank and P. W. Shor, *Phys. Rev. A* **54**, 1098 (1996).
 - [2] A. M. Steane, *Proc. R. Soc. London A* (to be published).
 - [3] B. Schumacher, *Phys. Rev. A* **51**, 2738 (1995).
 - [4] P. W. Shor, *Phys. Rev. A* **52**, R2493 (1995).
 - [5] A. M. Steane, *Phys. Rev. Lett.* **77**, 793 (1996).
 - [6] R. Laflamme, C. Miquel, J. P. Paz, and W. H. Zurek, *Phys. Rev. Lett.* **77**, 198 (1996).
 - [7] C. H. Bennett, D. P. DiVincenzo, J. A. Smolin, and W. K. Wootters (unpublished).
 - [8] F. J. MacWilliams and N. J. A. Sloane, *The Theory of Error-Correcting Codes* (North-Holland, Amsterdam, 1977).
 - [9] E. Knill and R. Laflamme, *Phys. Rev. A* (to be published).
 - [10] I. L. Chuang and R. Laflamme (unpublished).
 - [11] M. B. Plenio, V. Vedral, and P. L. Knight (unpublished).
 - [12] A. Ekert and C. Macchiavello, *Phys. Rev. Lett.* **77**, 2585 (1996).
 - [13] H. J. Helgert and R. D. Stinaff, *IEEE Trans. Inf. Theory* **19**, 344 (1973).
 - [14] P. Shor and J. A. Smolin (unpublished).
 - [15] A. R. Calderbank, E. M. Rains, N. J. A. Sloane, and P. W. Shor (unpublished).
 - [16] D. Gottesman, *Phys. Rev. A* **54**, 1862 (1996).