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Signal space

The notion of signal space is fundamental in communication. It allows a simple geometric approach to complex communication problems.

1.1 Review on vector space

Vector spaces form an algebraic structure based on Fields. A vector space is composed of a set V of elements called vectors. This set possesses a composition law called "addition of vector". The set of vectors along with the addition of vectors forms an abelian group. In addition, there exists a multiplication operation between elements (called "scalars") from a field F and vectors from giving vectors.

In this set of notes, we are going to use bold letters to designate vectors and normal letters for scalars. To recapitulate, we must have:

Given $V = \{\mathbf{x}_1, \mathbf{x}_2, \dots\}$, the elements must satisfy:

$$\forall \mathbf{x}, \mathbf{y}, \mathbf{z} \in V \quad (\mathbf{x} + \mathbf{y}) + \mathbf{z} = \mathbf{x} + (\mathbf{y} + \mathbf{z}) \quad \{\text{associativity}\}$$

$$\exists \text{ unique } \mathbf{0} \in V \mid \mathbf{x} + \mathbf{0} = \mathbf{0} + \mathbf{x} = \mathbf{x} \quad \{\text{additive identity}\}$$

$$\forall \mathbf{x} \in V ; \exists -\mathbf{x} \in V \mid \mathbf{x} + (-\mathbf{x}) = (-\mathbf{x}) + \mathbf{x} = \mathbf{0} \quad \{\text{additive inverse}\}$$

$$\forall \mathbf{x}, \mathbf{y} \in V \quad \mathbf{x} + \mathbf{y} = \mathbf{y} + \mathbf{x} \quad \{\text{commutativity}\}$$

and

$$\forall \alpha \in F \text{ and } \forall \mathbf{x} \in V \quad \alpha \mathbf{x} \in F \text{ such that}$$

$$\forall \alpha, \beta \in F \text{ and } \forall \mathbf{x} \in V \quad (\alpha + \beta)\mathbf{x} = \alpha\mathbf{x} + \beta\mathbf{x}$$

$$\forall \alpha \in F \text{ and } \forall \mathbf{x}, \mathbf{y} \in V \quad \alpha(\mathbf{x} + \mathbf{y}) = \alpha\mathbf{x} + \alpha\mathbf{y}$$

We say that we have a vector space V over the field F .

We can define many different types of vector spaces.

Example:

◆ The set of n -tuples from \mathbb{R} is a vector space: $\mathbf{x} = (x_1, x_2, \dots, x_n) \in \mathbb{R}^n$. The field is the set of real numbers \mathbb{R} .

◆ A set that is very important is the set of time limited functions that have a finite energy. The field is the set of complex numbers \mathbb{C} . We denote this set $L_2(T)$. The vector space is the set $L_2(T)$ over the set of complex numbers \mathbb{C} .

$x(t) = 0 \quad t \notin [0, T]$ such that $\int_0^T |x(t)|^2 dt < \infty$. When we consider it as a time function, we denote it $x(t)$ and when we consider it as a vector we designate it \mathbf{x} .

1.2 Linear independence

When we work with vectors, it is convenient to be able to express vectors as linear combination of other vectors. A vector \mathbf{y} is a linear combination of the following n vectors $\mathbf{x}_1, \mathbf{x}_2, \dots, \mathbf{x}_n$ if we can express it as:

$$\mathbf{y} = \sum_{k=1}^n \alpha_k \mathbf{x}_k \text{ where the scalars } \alpha_1, \alpha_2, \dots, \alpha_n \text{ are not all zero.}$$

Now, a set $\{\mathbf{x}_1, \mathbf{x}_2, \dots, \mathbf{x}_n\}$ of vectors from V forms a linearly independent set if no one of the vectors can be expressed as a linear combination of the others. This means that $\sum_{k=1}^n \alpha_k \mathbf{x}_k = 0$ if and only if $\alpha_1 = \alpha_2 = \dots = \alpha_n = 0$.

Dimension of the space

In a vector space, the maximum number of linearly independent vectors is called the dimension of the space. Some spaces are finite dimensional. The vectors that represent fields in electromagnetic theory are three dimensional vectors. However, function spaces are usually infinite dimensional spaces.

Example:

Consider the set of complex functions of the real variable t defined as follows:

$$\varphi_k(t) = \begin{cases} \exp(jkt) & 0 \leq t \leq 2\pi \\ 0 & \text{elsewhere} \end{cases}$$

We can show that $\sum_{k=-\infty}^{+\infty} \alpha_k \varphi_k(t) = 0$ if and only if $\alpha_k = 0$ for all $k \in \mathbb{Z}$.

Basis (plur. Bases)

When we possess a set $B = \{\mathbf{u}_1, \mathbf{u}_2, \dots, \mathbf{u}_n\}$ of vectors from V where $n = \dim V$, we can express any vector from V as a linear combination of the vectors composing B . We say that the set of vectors B forms a "basis". So, for $\forall \mathbf{x} \in V$, we can write $\mathbf{x} = \sum_{k=1}^n \alpha_k \mathbf{u}_k$. We say that the set B spans the space V and

we can write also $V = \text{span}\{\mathbf{u}_1, \mathbf{u}_2, \dots, \mathbf{u}_n\}$. At that time, the vectors can be represented by the sequence of coefficients $\mathbf{x} = (\alpha_1, \alpha_2, \dots, \alpha_n)$.

Example:

Consider the infinite set of functions defined previously. Any function that is nonzero only in the interval $[0, 2\pi]$ and that satisfies the Dirichlet conditions (see EE311 course) can be represented as the following linear combination:

$$x(t) = \sum_{k=-\infty}^{+\infty} \alpha_k \exp(jkt) \quad (1.1)$$

Equation (1.1) is nothing but the Fourier series representation of the signal $x(t)$. This signal is represented using the above basis and its coordinates are the Fourier coefficients α_k .

$$\alpha_k = \frac{1}{2\pi} \int_0^{2\pi} x(t) \exp(-jkt) dt \quad (1.2)$$

The vector \mathbf{x} can then be represented by the infinite sequence:

$$\mathbf{x} = (\dots, \alpha_{-1}, \alpha_0, \alpha_1, \dots, \alpha_k, \dots)$$

Subspace

If we are given a vector space V over the field F , we can define a subspace as follows:

W is a subspace of V if it is a subset of V closed under the addition of vectors and multiplication by scalars from F . In finite spaces, we always have $\dim W < \dim V$.

1.3 Metric space and norm

Distance

If we want to compare elements from a given space, we have to define the notion of metric or distance.

The distance between two elements a and b of a set S is the positive real number $d(a, b)$ satisfying:

1. $\forall a, b \in S \quad d(a, b) \geq 0$, with $d(a, b) = 0$ if and only if $a = b$
2. $\forall a, b, c \in S \quad d(a, c) \leq d(a, b) + d(b, c)$ (triangular inequality)

The set S is called a metric space.

Norm

If now the set S possesses the structure of a vector space over \mathbb{C} , we can define the length of the vectors using the concept of norm. The norm of a vector \mathbf{x} is the positive real number $\|\mathbf{x}\|$ satisfying:

1. $\forall \mathbf{x} \in V \quad \|\mathbf{x}\| \geq 0$, with $\|\mathbf{x}\| = 0$ if and if $\mathbf{x} = 0$
2. $\forall \mathbf{x}, \mathbf{y} \in V \quad \|\mathbf{x} + \mathbf{y}\| \leq \|\mathbf{x}\| + \|\mathbf{y}\|$
3. $\forall \mathbf{x} \in V \text{ and } \forall \alpha \in \mathbb{C} \quad \|\alpha \mathbf{x}\| = |\alpha| \|\mathbf{x}\|$

When a norm is defined in a vector space, we can deduce from it a distance between points in the space. If we consider that a vector is constituted by 2 points: Its origin (at the origin of the basis) and its tip, we can use it to

represent points in the space spanned by the basis. So, the vector \mathbf{x} represents also the point in the above space having as coordinates the coordinates of \mathbf{x} . So, given two points represented by \mathbf{x} and \mathbf{y} , the distance between \mathbf{x} and \mathbf{y} can be expressed by:

$$d(\mathbf{x}, \mathbf{y}) = \|\mathbf{x} - \mathbf{y}\| \quad (1.3)$$

Now if a metric space has the structure of a vector space over \mathbb{C} , a norm can be defined using the distance. Since a vector \mathbf{x} is represented by two points (the origin and its tip) and the origin is represented by zero, then we can write:

$$\|\mathbf{x}\| = d(0, \mathbf{x}) \quad (1.4)$$

1.4 Inner product

The concept of inner product can be defined only for vectors spaces defined over the field of complex (real) numbers \mathbb{C} (\mathbb{R}). Given a vectors space V defined over \mathbb{C} , the inner product of two vectors is defined as follows:

Let $\mathbf{x}, \mathbf{y} \in V$, the inner product of \mathbf{x} and \mathbf{y} is the complex number (\mathbf{x}, \mathbf{y}) satisfying:

1. $\forall \mathbf{x} \in V \quad (\mathbf{x}, \mathbf{x}) \geq 0$, with $(\mathbf{x}, \mathbf{x}) = 0$ if and only if $\mathbf{x} = 0$
2. $\forall \mathbf{x}, \mathbf{y} \in V \quad (\mathbf{x}, \mathbf{y}) = (\mathbf{y}, \mathbf{x})^*$
3. $\forall \mathbf{x}, \mathbf{y}, \mathbf{z} \in V \quad \text{and} \quad \alpha, \beta \in \mathbb{C} \quad (\alpha\mathbf{x} + \beta\mathbf{y}, \mathbf{z}) = \alpha(\mathbf{x}, \mathbf{z}) + \beta(\mathbf{y}, \mathbf{z})$

Example:

Consider the two dimensional real vectors represented by their coordinates: $\vec{\mathbf{x}} = (x_1, x_2)$ where $x_1, x_2 \in \mathbb{R}$. The inner product of the two vectors $\vec{\mathbf{x}}$ and $\vec{\mathbf{y}}$ is usually written as $\vec{\mathbf{x}} \cdot \vec{\mathbf{y}} = x_1 y_1 + x_2 y_2 = |\vec{\mathbf{x}}| |\vec{\mathbf{y}}| \cos(\vec{\mathbf{x}}, \vec{\mathbf{y}})$. It is also called "dot product".

If we consider the space $L_2(T)$ defined previously, the inner product is defined as:

$$(\mathbf{x}, \mathbf{y}) = \int_0^T x(t) y^*(t) dt \quad (1.5)$$

Euclidian Norm

The properties of the inner product can be used to define a norm in the vector space: the Euclidian norm

$$\|\mathbf{x}\| = \sqrt{(\mathbf{x}, \mathbf{x})} \quad (1.6)$$

In the two dimensional space shown above, the length of a vector $\vec{\mathbf{x}}$ having as coordinates x_1 and x_2 is $\|\vec{\mathbf{x}}\| = \sqrt{x_1^2 + x_2^2}$

In the space $L_2(T)$, the norm of a function is:

$$\|\mathbf{x}\|^2 = \int_0^T |x(t)|^2 dt \text{ and this is the energy of the signal } x(t).$$

Properties of the inner product

Cauchy-Schwartz inequality

The Cauchy-Schwartz inequality is an important property of the inner product. It is commonly used to solve some optimization problems that occur in communication theory. It is also used to define the concept of an angle between two vectors.

Given two non zero vectors \mathbf{x} and \mathbf{y} from the vector space V defined over \mathbb{C} , the following inequality applies:

$$|(\mathbf{x}, \mathbf{y})| \leq \|\mathbf{x}\| \|\mathbf{y}\| \text{ with equality if } \mathbf{x} = \alpha \mathbf{y}, \alpha \text{ being a non zero scalar.}$$

Proof:

Let \mathbf{x} and \mathbf{y} be two nonzero vectors from the vector space V defined over \mathbb{C} and the arbitrary nonzero scalar λ , consider the positive number:

$$D(\lambda) = \|\mathbf{x} - \lambda \mathbf{y}\|^2 = (\mathbf{x} - \lambda \mathbf{y}, \mathbf{x} - \lambda \mathbf{y})$$

Developing the product, we obtain:

$$\begin{aligned} D(\lambda) &= \|\mathbf{x}\|^2 - \lambda(\mathbf{y}, \mathbf{x}) - \lambda^*(\mathbf{x}, \mathbf{y}) + |\lambda|^2 \|\mathbf{y}\|^2 \\ &= \|\mathbf{x}\|^2 - \lambda(\mathbf{x}, \mathbf{y})^* - \lambda^*(\mathbf{x}, \mathbf{y}) + |\lambda|^2 \|\mathbf{y}\|^2 \end{aligned}$$

λ being arbitrary, let $\lambda = \frac{(\mathbf{x}, \mathbf{y})}{\|\mathbf{y}\|^2}$. Using this value, $D(\lambda)$ becomes:

$$D(\lambda) = \|\mathbf{x}\|^2 - \frac{|(\mathbf{x}, \mathbf{y})|^2}{\|\mathbf{y}\|^2}. \text{ Since } D(\lambda) \geq 0, \text{ the inequality is proved.}$$

We will have equality if $D(\lambda) = 0$. In this case, $\mathbf{x} = \lambda \mathbf{y}$.

◆

Using the Cauchy-Schwartz inequality, we can define the angle between the vector \mathbf{x} and \mathbf{y} as:

$$\cos \theta = \frac{(\mathbf{x}, \mathbf{y})}{\|\mathbf{x}\| \|\mathbf{y}\|} \quad (1.7)$$

If $\mathbf{x} = \lambda \mathbf{y}$, the angle between the two vectors will be 0 or 180° ($\cos \theta = \pm 1$). This means that the two vectors are collinear. If the two vectors are nonzero,

they will be orthogonal if $(\mathbf{x}, \mathbf{y}) = 0$ ($\theta = \pm \frac{\pi}{2}$). We use the notation $\mathbf{x} \perp \mathbf{y}$ to say \mathbf{x} orthogonal with \mathbf{y} .

In the $L_2(T)$ space, the Cauchy-Schwartz inequality can be expressed as:

$$\left| \int_0^T x(t)y^*(t)dt \right|^2 \leq \int_0^T |x(t)|^2 dt \int_0^T |y(t)|^2 dt$$

Pythagorean Theorem

If two vectors \mathbf{x} and \mathbf{y} are orthogonal, they satisfy the Pythagorean theorem:

$$\|\mathbf{x} + \mathbf{y}\|^2 = \|\mathbf{x}\|^2 + \|\mathbf{y}\|^2$$

Proof:

$\|\mathbf{x} + \mathbf{y}\|^2 = (\mathbf{x} + \mathbf{y}, \mathbf{x} + \mathbf{y}) = \|\mathbf{x}\|^2 + (\mathbf{x}, \mathbf{y}) + (\mathbf{y}, \mathbf{x}) + \|\mathbf{y}\|^2$. However, $\mathbf{x} \perp \mathbf{y}$, so $(\mathbf{x}, \mathbf{y}) = (\mathbf{y}, \mathbf{x}) = 0$.

◆

1.5 Orthonormal basis

In signal spaces (and even in general vector spaces) it is convenient to represent vectors using an "orthonormal basis". Consider a space V such that $\dim V = n$. (n can be infinite). The set of vectors $\{\boldsymbol{\varphi}_1, \boldsymbol{\varphi}_2, \dots, \boldsymbol{\varphi}_n\}$ forms an orthonormal basis if the vectors satisfy:

$$(\boldsymbol{\varphi}_k, \boldsymbol{\varphi}_i) = \begin{cases} 0 & k \neq i \\ 1 & k = i \end{cases}$$

These vectors are linearly independent and their number is equal to the dimension of the space. So, we can write $V = \text{span}\{\boldsymbol{\varphi}_1, \boldsymbol{\varphi}_2, \dots, \boldsymbol{\varphi}_n\}$. This means that any vector in V can be expressed as:

$$\mathbf{x} = \sum_{k=1}^n \alpha_k \boldsymbol{\varphi}_k \tag{1.8}$$

One advantage of an orthonormal basis is the simple expression of the coefficients. If we compute the inner product of \mathbf{x} with one basis vector, we obtain:

$$(\mathbf{x}, \boldsymbol{\varphi}_m) = \left(\sum_{k=1}^n \alpha_k \boldsymbol{\varphi}_k, \boldsymbol{\varphi}_m \right) = \sum_{k=1}^n \alpha_k (\boldsymbol{\varphi}_k, \boldsymbol{\varphi}_m)$$

So:

$$\alpha_m = (\mathbf{x}, \boldsymbol{\varphi}_m) \quad (1.9)$$

In the case of the $L_2(T)$ space, the expression of the coefficients is then

$$\alpha_m = \int_0^T x(t) \boldsymbol{\varphi}_m(t) dt \quad (1.10)$$

In the signal space $L_2(T)$, the expression (1.8) is called Fourier decomposition of the signal $x(t)$ and is written as:

$$x(t) = \sum_{k=1}^n \alpha_k \boldsymbol{\varphi}_k(t) \quad (1.11)$$

The coefficients α_k are called the Fourier coefficients.

Orthogonal spaces

Consider a vector space V of dimension n . It can be decomposed into the "sum" of two "orthogonal" subspaces V_1 and V_2 such that:

$$\begin{aligned} \forall \mathbf{x}_1 \in V_1 \quad \text{and} \quad \forall \mathbf{x}_2 \in V_2 \quad \mathbf{x}_1 \perp \mathbf{x}_2 \\ \forall \mathbf{x} \in V \quad \exists \mathbf{x}_1 \in V_1 \quad \exists \mathbf{x}_2 \in V_2 \quad | \quad \mathbf{x} = \mathbf{x}_1 + \mathbf{x}_2 \end{aligned}$$

Using orthonormal bases, we can easily decompose any vector into the two above components. Assume that we are given an orthonormal basis such that $V = \text{span}\{\boldsymbol{\varphi}_1, \boldsymbol{\varphi}_2, \dots, \boldsymbol{\varphi}_n\}$ and let us define two subspaces $V_1 = \text{span}\{\boldsymbol{\varphi}_1, \dots, \boldsymbol{\varphi}_m\}$ and $V_2 = \text{span}\{\boldsymbol{\varphi}_{m+1}, \dots, \boldsymbol{\varphi}_n\}$. It is clear that V_1 and V_2 are orthogonal. If we consider any vector of V , its first m coordinates define an element of V_1 while the last $n - m$ coordinates define an element of V_2 . So:

$\forall \mathbf{x} \in V \quad \mathbf{x}_1 = \sum_{k=1}^m \alpha_k \boldsymbol{\varphi}_k \in V_1 \quad \text{and} \quad \mathbf{x}_2 = \sum_{k=m+1}^n \alpha_k \boldsymbol{\varphi}_k \in V_2$. The coefficients α_k are computed using equation (1.9): $\alpha_k = (\mathbf{x}, \boldsymbol{\varphi}_k)$. The vector \mathbf{x}_1 is the orthogonal projection of \mathbf{x} on V_1 while the vector \mathbf{x}_2 is the orthogonal projection of \mathbf{x} on V_2 .

The Gram-Schmidt procedure

The Gram-Schmidt procedure is a procedure for deriving an orthonormal basis from a set of M vectors spanning a space of dimension $K \leq M$. So, we are given the set $\{\mathbf{x}_1, \mathbf{x}_2, \dots, \mathbf{x}_M\}$ of vectors, not necessarily linearly independent.

The procedure follows the subsequent steps:

1. $\boldsymbol{\varphi}_1 = \frac{\mathbf{x}_1}{\|\mathbf{x}_1\|}$; this is an initialization step.

2. For $k = 2$ to M let:

$$\mathbf{g}_k = \mathbf{x}_k - \sum_{i=1}^{k-1} (\mathbf{x}_k, \boldsymbol{\varphi}_i) \boldsymbol{\varphi}_i; \text{ We subtract from } \mathbf{x}_k \text{ its projection on the}$$

$k - 1$ dimensional space spanned by the already found basis vectors. This implies that \mathbf{g}_k belongs to a space that is orthogonal to the space spanned by the $k - 1$ basis vectors $\boldsymbol{\varphi}_i$. If \mathbf{g}_k is zero, the vector \mathbf{x}_k must be discarded. It belongs to the previous space and it is not linearly independent. The basis vector is then: $\boldsymbol{\varphi}_k = \frac{\mathbf{g}_k}{\|\mathbf{g}_k\|}$.

At the end of step 2, we will have a set of K orthonormal basis vectors. If the set $\{\mathbf{x}_1, \mathbf{x}_2, \dots, \mathbf{x}_M\}$ is composed of linearly independent vectors, we will have $K = M$. Otherwise, the dimension of the space will be smaller than the number of given vectors.

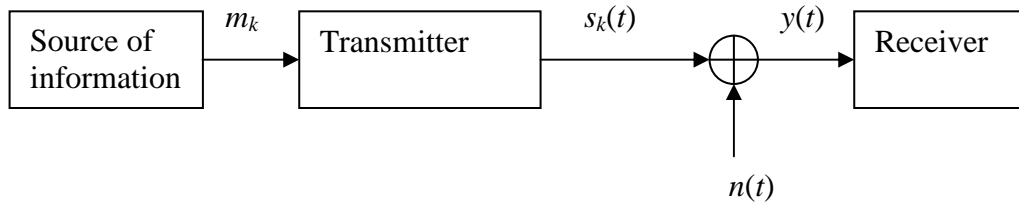
Expression of the inner product in an orthonormal space

Let us consider two vectors \mathbf{x} and \mathbf{y} with coordinates (x_1, x_2, \dots, x_K) and (y_1, y_2, \dots, y_K) in an orthonormal basis. It is easy to show that:

$$(\mathbf{x}, \mathbf{y}) = \left(\sum_{k=1}^K x_k \boldsymbol{\varphi}_k, \sum_{j=1}^K y_j \boldsymbol{\varphi}_j \right) = \sum_{k=1}^K x_k y_k^* \quad (1.12)$$

1.6 Waveform communication system

Let us consider the following communication system



We are given a source of information that can produce M different symbols. The transmitter assigns a signal $s_k(t)$ to the message m_k . The signals belong to the space $L_2(T)$. This means that the duration of a symbol is T seconds and the rate of symbols (it is called the "baud rate") is $r = \frac{1}{T}$. The total number of signals produced by the transmitter is M . These signals belong to a finite

dimensional space. In fact, we can use these M signals to derive an orthonormal basis using the Gram-Schmidt procedure. The derived basis is composed of $K \leq M$ orthonormal signals $\varphi_k(t)$. Let us call this space S_K , $S_K = \text{span}\{\varphi_1, \dots, \varphi_K\}$. The signals generated by the transmitter can be expressed in this space as:

$$s_i(t) = \sum_{k=1}^K \alpha_{i,k} \varphi_k(t) \quad ; \quad i = 1, \dots, M$$

So, we have a correspondence between the signal $s_i(t)$ and the vector $\mathbf{s}_i = (\alpha_{i,1}, \alpha_{i,2}, \dots, \alpha_{i,K})$.

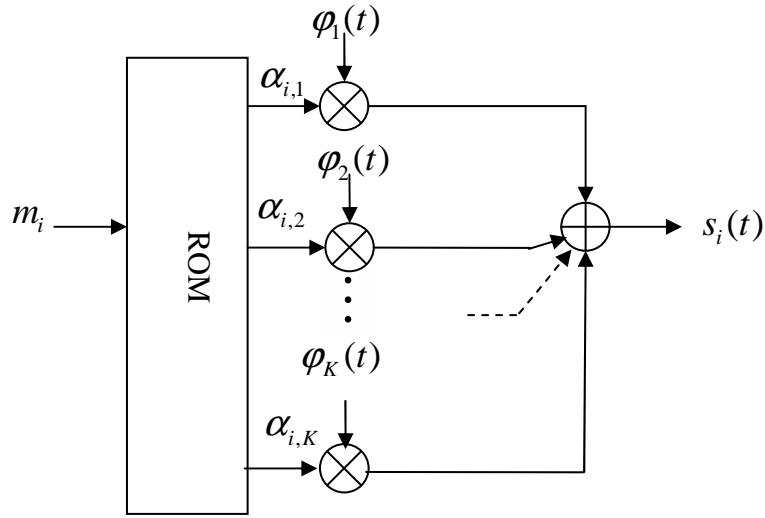


Fig.3- 1 Possible transmitter structure

Fig.3- 1 shows a possible implementation of a transmitter. The symbol m_i selects a set of K coefficients. The waveform produced by the transmitter is a sequence of finite time signals, each one being a weighted sum of basis functions.

The noise signal $n(t)$ is assumed to be a white Gaussian noise with a power spectrum $S_n(f) = \frac{N_0}{2}$. The received signal $y(t)$ is the sum of the transmitted signal and this noise signal. Despite the fact that the signal at the output of the transmitter belongs to a finite dimensional space, the signal at the input of the receiver is a signal which belongs to an infinite dimensional space. We can see this by developing the noise process.

Consider the following orthonormal basis: $\{\varphi_1, \varphi_2, \dots, \varphi_K, \varphi_{K+1}, \dots\}$ where the first K basis vectors are the vector representation of the basis functions used to represent the signals generated by the transmitter. We complete this basis by

an arbitrary set of orthonormal functions to span an infinite dimensional space. In this basis, we can show that the noise process is expressed as:

$$n(t) = \sum_{k=1}^{\infty} N_k \varphi_k(t) = \sum_{k=1}^K N_k \varphi_k(t) + \sum_{k=K+1}^{\infty} N_k \varphi_k(t) \quad (1.13)$$

Equation (1.13) shows that the noise process is the sum of two orthogonal processes. The first one belongs to the same space as the signals generated by the transmitter (S_K). The second one is orthogonal to all these signals. The coordinates N_k are random variables and they are the projections of the noise vector on the basis functions.

$$N_k = (\mathbf{n}, \boldsymbol{\varphi}_k) = \int_0^T n(t) \varphi_k(t) dt \quad (1.14)$$

Since the noise process is Gaussian, these variables are Gaussian random variables. We can derive their statistics as follows:

$$E[N_k] = E\left[\int_0^T n(t) \varphi_k(t) dt\right] = \int_0^T E[n(t)] \varphi_k(t) dt = 0$$

So, they are zero mean. Their covariance is then:

$$E[N_k N_j] = E\left[\int_0^T n(t) \varphi_k(t) dt \int_0^T n(u) \varphi_j(u) du\right] = \int_0^T \int_0^T E[n(t)n(u)] \varphi_k(t) \varphi_j(u) dt du$$

The noise process is white. Its autocorrelation function is:

$$R_n(t-u) = E[n(t)n(u)] = \frac{N_0}{2} \delta(t-u)$$

The covariance is then:

$$E[N_k N_j] = \int_0^T \frac{N_0}{2} \varphi_k(t) \varphi_j(t) dt$$

Since the basis functions are orthonormal, the above covariance is:

$$E[N_k N_j] = \begin{cases} \frac{N_0}{2} & k = j \\ 0 & k \neq j \end{cases}$$

So, the components of the noise process are independent Gaussian random variables with zero mean and variances equal to $\frac{N_0}{2}$. These components will be the same in any orthonormal basis. A white Gaussian noise is always represented by an infinite dimensional vector with coordinates that are independent zero mean Gaussian vector with variance equal to $\frac{N_0}{2}$.

The signal observed by the receiver is the signal $y(t)$. It consists of the sum of a finite dimensional vector ($s_i(t)$) and the infinite dimensional vector

representing the white noise process. However, we can express this received signal as a sum of two orthogonal signals.

$$y(t) = s_i(t) + n_1(t) + n_2(t) = z(t) + n_2(t)$$

where $n_1(t) = \sum_{k=1}^K N_k \phi_k(t) \in S_K$ and $n_2(t) = \sum_{k=K+1}^{\infty} N_k \phi_k(t) \perp S_K$. The signal $z(t)$ is $z(t) = s_i(t) + n_1(t)$. It is clear that $z(t) \in S_K$ and that $z(t) \perp n_2(t)$. We can show that the observation of the vector representing the signal $n_2(t)$ is not needed if we want to detect which message has been transmitted. So, at the receiver, we are going to observe only the signal $z(t)$ and not the signal $y(t)$. The signal $z(t)$ is the orthogonal projection of $y(t)$ on the space S_K and it can be represented by a finite dimensional vector $\mathbf{z} = (z_1, z_2, \dots, z_K)$. The signal $y(t)$ on the other hand must be represented by an infinite dimensional vector $\mathbf{y} = (y_1, y_2, \dots, y_K, y_{K+1}, \dots)$ and the first K coordinates correspond to the vector \mathbf{z} .

We can now write the expression of the conditional pdf of the vector \mathbf{z} given that the symbol m_i is transmitted. The coordinates of \mathbf{z} are:

$z_k = \alpha_{i,k} + N_k$ for $k = 1, \dots, K$. The numbers $\alpha_{i,k}$ are the coordinates of the signal $s_i(t)$. So, the coordinates z_k are independent Gaussian random variables with mean $\alpha_{i,k}$ and variance $\frac{N_0}{2}$. The conditional pdf of \mathbf{z} is:

$$f_{\mathbf{z}|m_i}(\mathbf{z} | m_i) = \prod_{k=1}^K \frac{1}{\sqrt{2\pi} \sqrt{\frac{N_0}{2}}} \exp \left[-\frac{(z_k - \alpha_{i,k})^2}{N_0} \right]$$

or:

$$f_{\mathbf{z}|m_i}(z | m_i) = \frac{1}{(\pi N_0)^{\frac{K}{2}}} \exp \left[-\frac{1}{N_0} \sum_{k=1}^K (z_k - \alpha_{i,k})^2 \right] \quad (1.15)$$

We can use equation (1.15) to determine the structure of an MAP receiver. We have seen that the MAP rule is:

"Decide m_i if $P(m_i) f_{\mathbf{z}|m_i}(\mathbf{z} | m_i) > P(m_j) f_{\mathbf{z}|m_j}(\mathbf{z} | m_j)$ for all $j \neq i ; j = 1, \dots, M$ "

Replacing and taking logarithms, we obtain:

"Decide m_i if $\left\{ \sum_{k=1}^K (z_k - \alpha_{i,k})^2 - N_0 \ln P(m_i) \right\}$ min " (1.16)

The first term in the above expression is just the square of the distance between the vector \mathbf{z} and the vector \mathbf{s}_i . So we can re-express the rule (1.16) as:

$$\left\{ \|\mathbf{z} - \mathbf{s}_i\|^2 - N_0 \ln P(m_i) \right\} \quad \min \quad (1.17)$$

If the symbols are equiprobable (ML receiver), the second term of (1.17) becomes irrelevant and the decision rule simplifies to:

$$\text{Decide } m_i \text{ such that } \|\mathbf{z} - \mathbf{s}_i\| \min \quad (1.18)$$

The receiver is called a "minimum distance classifier". In order to build the receiver, we must determine the coordinates of the vector \mathbf{z} . We have seen that they are the first K coordinates of the vector \mathbf{y} . From (1.9), the coordinates of \mathbf{y} are given by:

$$y_k = z_k = (\mathbf{y}, \boldsymbol{\varphi}_k) = \int_0^T y(t) \varphi_k(t) dt \text{ for } k = 1, \dots, K.$$

So, the receiver is composed of three cascaded sections. The first one computes the coordinates of the vector \mathbf{z} . The second one is a distance computer and the third one applies a bias to take into account the unequal a priori probabilities.

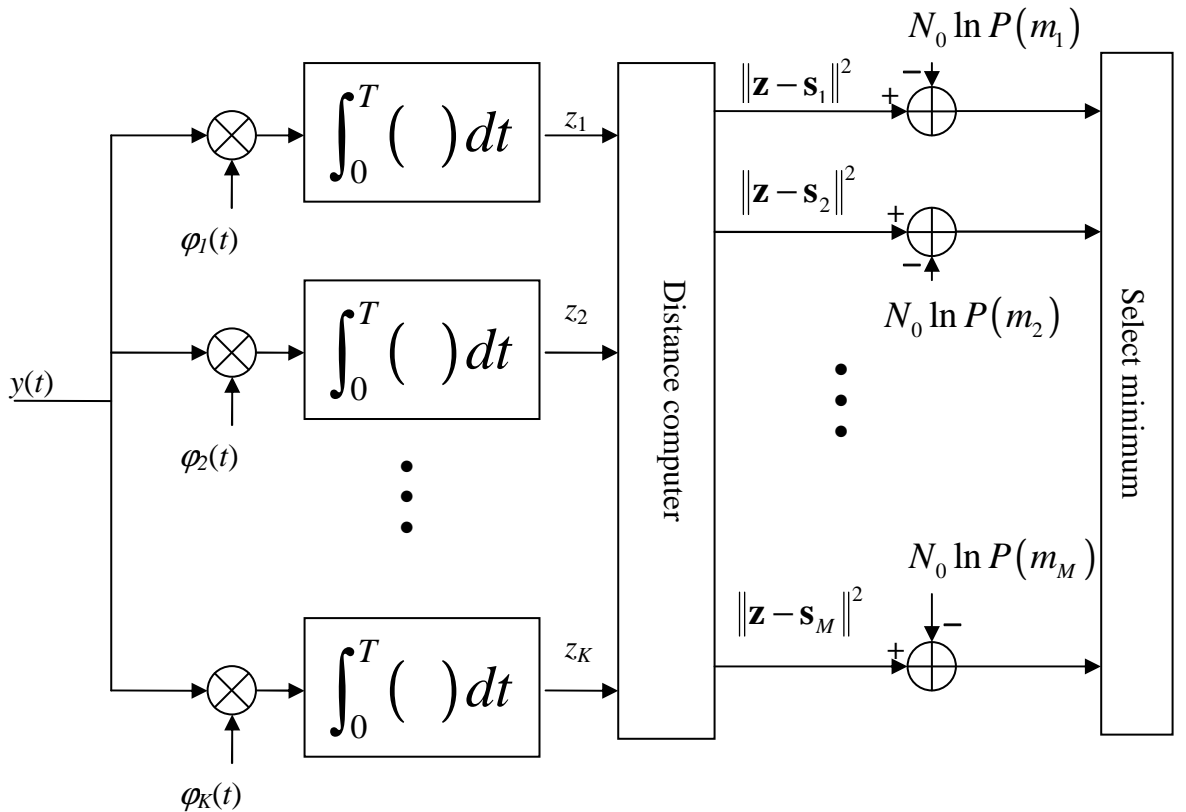


Fig.3- 2 Receiver structure using distance computer

The above structure uses K "correlators" that transform the random process $y(t)$ to the random vector \mathbf{z} . The distance computer has K inputs (coordinates of \mathbf{z}) and computes M distances (between the received vector \mathbf{z} and

the "prototypes" \mathbf{s}_i). If the symbols are equiprobable (this is often the case), the bias $-N_0 \ln P(m_i)$ is not required.

A different structure can be obtained if we develop the square of the distance.

$$\|\mathbf{z} - \mathbf{s}_i\|^2 = (\mathbf{z} - \mathbf{s}_i, \mathbf{z} - \mathbf{s}_i) = (\mathbf{z}, \mathbf{z}) + (\mathbf{s}_i, \mathbf{s}_i) - 2(\mathbf{z}, \mathbf{s}_i) \quad (1.19)$$

In the above expression, the squared norm of \mathbf{z} does not depend on the message. The same value $\|\mathbf{z}\|^2$ will appear for all symbols. It can be eliminated. The inner product of \mathbf{s}_i with itself is the energy of the signal $s_i(t)$.

$$(\mathbf{s}_i, \mathbf{s}_i) = \|\mathbf{s}_i\|^2 = \int_0^T s_i^2(t) dt = E_i$$

So, the expression (1.18) simplifies to:

$$\|\mathbf{z} - \mathbf{s}_i\|^2 - N_0 \ln P(m_i) = cste + E_i - 2(\mathbf{z}, \mathbf{s}_i) - N_0 \ln P(m_i)$$

The decision rule becomes:

$$\text{Decide } m_i \text{ if } \left\{ (\mathbf{z}, \mathbf{s}_i) - \frac{E_i}{2} + \frac{N_0}{2} \ln P(m_i) \right\} \max.$$

The inner product $(\mathbf{z}, \mathbf{s}_i)$ can be computed using the signal $y(t)$ present at the input of the receiver because $y(t) = z(t) + n_2(t)$ and $n_2(t) \perp z(t)$. So:

$$(\mathbf{z}, \mathbf{s}_i) = (\mathbf{y} - \mathbf{n}_2, \mathbf{s}_i) = (\mathbf{y}, \mathbf{s}_i) = \int_0^T y(t) s_i(t) dt$$

Finally, the decision rule is:

$$\text{Decide } m_i \text{ if } \left\{ \left[\int_0^T y(t) s_i(t) dt \right] - \frac{E_i}{2} + \frac{N_0}{2} \ln P(m_i) \right\} \max \quad (1.20)$$

The decision rule (1.20) can be implemented using the structure shown in Fig.3- 3. This structure uses M correlators and does not require a distance computer. If the symbols are equiprobable, the bias becomes only $\frac{E_i}{2}$. The correlation receiver has a simpler structure than the minimum distance one. However, it uses M correlators. In many cases, (PAM, MPSK and QAM) the dimension of the signal space is less or equal to two while the number of signals can be very high. In this case, the minimum distance receiver is better. However, when the dimensionality of the signal space is high and M is equal to K , it is better to use the correlation receiver structure.

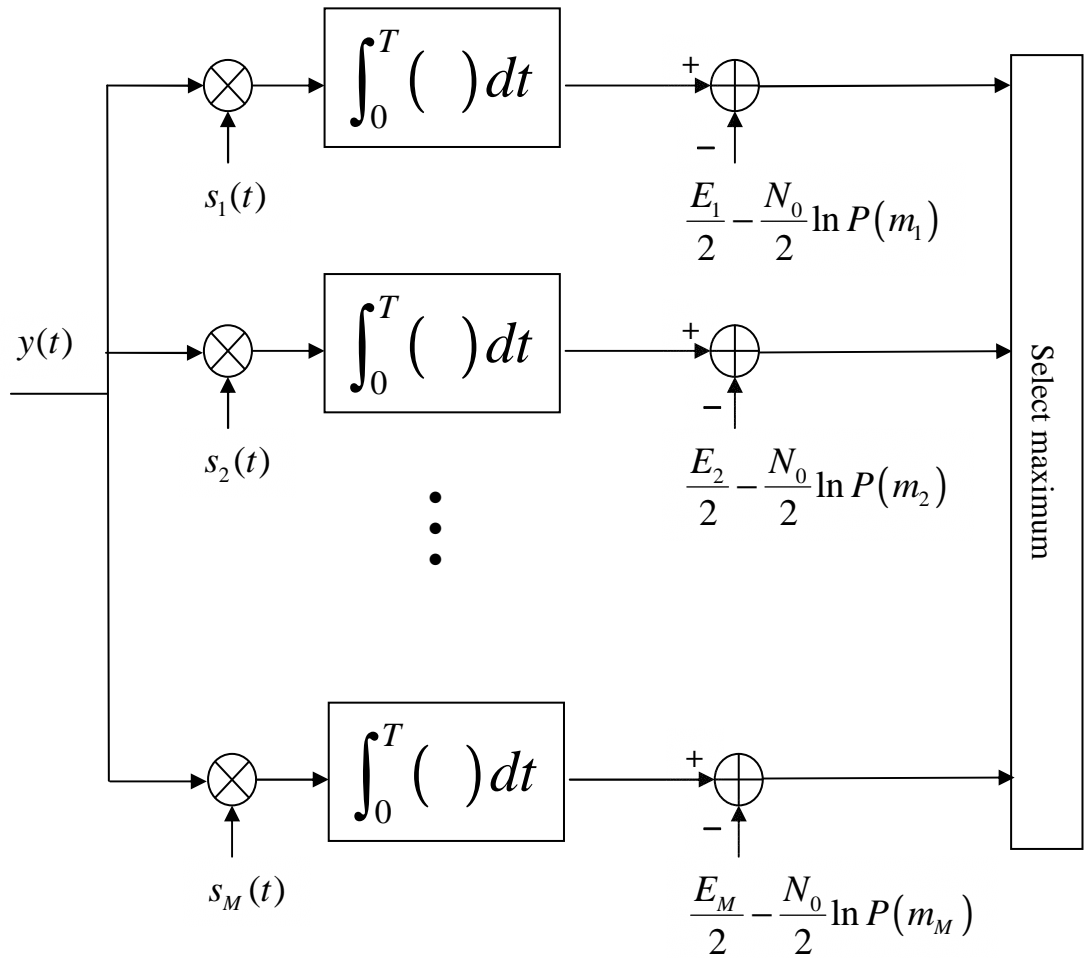


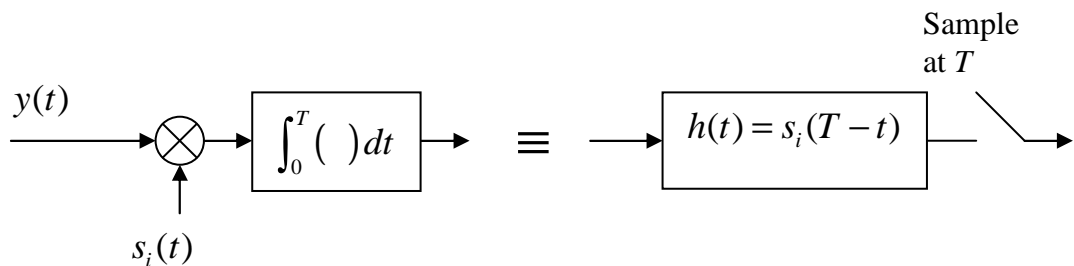
Fig.3- 3 Correlation receiver

The different correlators can be implemented by matched filters. We have already studied this equivalence in the matched filter course.

The correlator output is

$$\int_0^T y(t)s_i(t)dt = \int_0^T y(T-u)s_i(T-u)du = \left[\int_0^t s_i(T-u)y(t-u)du \right]_{t=T}$$

The last integral is the evaluation of the convolution of a filter with impulse response $h(t) = s_i(T-t)$ with the signal $y(t)$ at the time $t = T$. So, we obtain the following equivalence:



The decision rule divides the observation space into decision regions I_k such that $z \in I_k \Rightarrow$ decide m_k . If the symbols are equiprobable, the decision I_k is the set of points that are closer to the point s_k than to the other points.

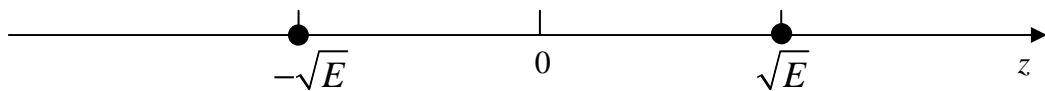
1.7 Some examples

For all examples, we assume equiprobable symbols.

Antipodal signaling

This communication system is a one-dimensional binary system. We use two signals:

$s_1(t) = \sqrt{E}\varphi(t)$ and $s_2(t) = -\sqrt{E}\varphi(t)$. The function $\varphi(t)$ has an energy equal to one. Using φ as a basis, the signal space is represented below.



We can use a minimum distance classifier. It uses only one correlator. Let the output of the correlator be called z . In this particular case, we don't have to compute the distance between the observed value z and the points $-\sqrt{E}$ and \sqrt{E} . The decision regions are separated by the origin. The optimum receiver is:

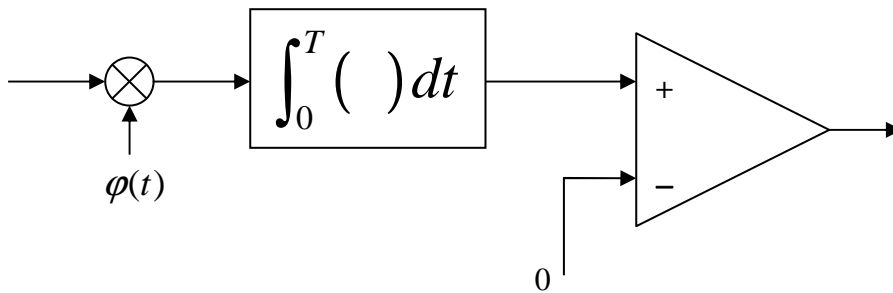


Fig.3- 4 Antipodal signalling receiver

Given that the distance between the mean of each distribution and the threshold is $A = \sqrt{E}$ and the variance is $\frac{N_0}{2}$, the probability of error is:

$$P(E) = \frac{1}{2} \operatorname{erfc}\left(\frac{A}{\sqrt{2}\sigma}\right) = \frac{1}{2} \operatorname{erfc}\sqrt{\frac{E}{N_0}} \quad (1.21)$$

This is the same result as the one deduced using the matched filter derived in the previous chapter.

Pulse amplitude modulation (PAM)

The PAM system is also a one dimensional system. However, in this system, we use M different amplitudes. The amplitudes are equally spaced and are such that the average value is zero. Furthermore, we assume that $M = 2^N$ in order to be able to encode the different levels in binary. So, M levels correspond to $N = \log_2 M$ binary digits. The level separation is A .

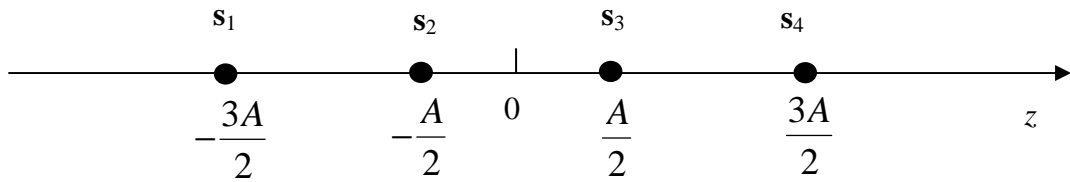


Fig.3- 5 PAM with $M = 4$

The decision regions are intervals separated by thresholds occurring midway between the signal levels. For example, the region I_2 corresponding to the signal s_2 is the interval $[-A, 0]$. The region I_1 corresponding to s_1 is the interval $]-\infty, -A]$. For the general case of M different signal levels, we are going to have $M - 1$ thresholds.

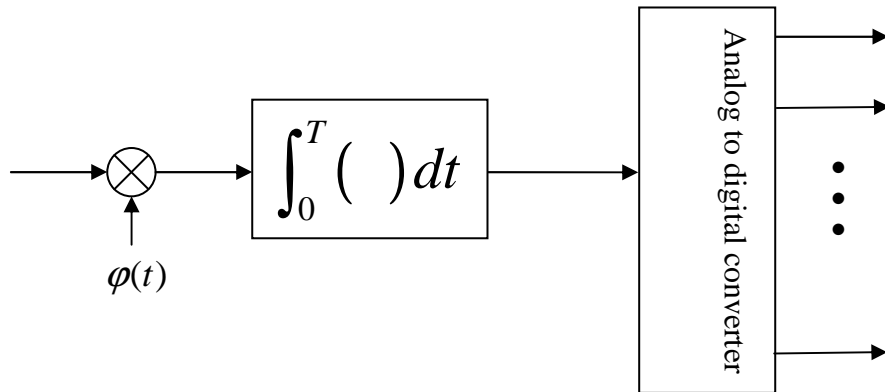


Fig.3- 6 PAM receiver

The above figure shows a typical PAM receiver. The analog to digital converter encodes directly the signal levels into binary. It can be implemented using $M - 1$ comparators followed by a priority encoder or it can be implemented using a successive approximation analog to digital converter. The

different levels are usually encoded in Gray code, so that an error in one level will cause only one bit to be in error.

To compute the probability of error, we have to take into consideration whether the level is an intermediate level (between two thresholds) or an extreme one (bounded by one threshold only). We have two extreme levels: \mathbf{s}_1 and \mathbf{s}_M . The conditional probability of error is the area of a tail of a Gaussian with a threshold away from the mean by $\frac{A}{2}$ and a variance $\frac{N_0}{2}$. So:

$$P(E | m_1) = P(E | m_M) = \frac{1}{2} \operatorname{erfc} \left(\frac{A}{2\sqrt{2}\sqrt{\frac{N_0}{2}}} \right) = \frac{1}{2} \operatorname{erfc} \left(\frac{A}{2\sqrt{N_0}} \right)$$

For the intermediate levels, we have two thresholds, so for the levels \mathbf{s}_2 up to \mathbf{s}_{M-1} , the conditional probability of error is the probability of being away from the mean by $\frac{A}{2}$ on both sides. So:

$$P(E | m_2) = \dots = P(E | m_{M-1}) = \operatorname{erfc} \left(\frac{A}{2\sqrt{2}\sqrt{\frac{N_0}{2}}} \right) = \operatorname{erfc} \left(\frac{A}{2\sqrt{N_0}} \right)$$

Finally, the probability of error is the average:

$$P(E) = \sum_{k=1}^M P(m_k) P(E | m_k) = \left(\frac{M-1}{M} \right) \operatorname{erfc} \left(\frac{A}{2\sqrt{N_0}} \right) \quad (1.22)$$

We can express the amplitude A as a function of the average energy of the signals:

$$E = \frac{1}{M} \sum_{k=1}^M \|\mathbf{s}_k\|^2 = \frac{1}{M} \sum_{k=1}^M |\alpha_k|^2$$

where the coefficients α_k are the amplitudes of the signals.

For the PAM signaling where the spacing between the points has a value of A , the different amplitudes are: $\pm \frac{A}{2}(2k-1)$ for $k = 1$ to $\frac{M}{2}$. The average energy is then:

$$E = \frac{2}{M} \sum_{k=1}^{\frac{M}{2}} \frac{A^2}{4} (2k-1)^2 = \frac{M^2-1}{12} A^2$$

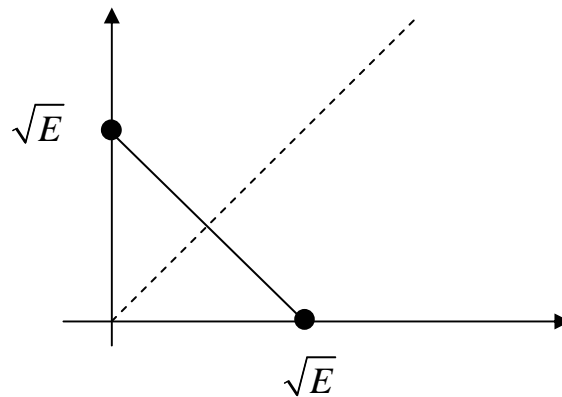
Replacing in (1.22), we obtain:

$$P(E) = \left(\frac{M-1}{M} \right) \operatorname{erfc} \sqrt{\frac{3}{M^2-1} \frac{E}{N_0}} \quad (1.23)$$

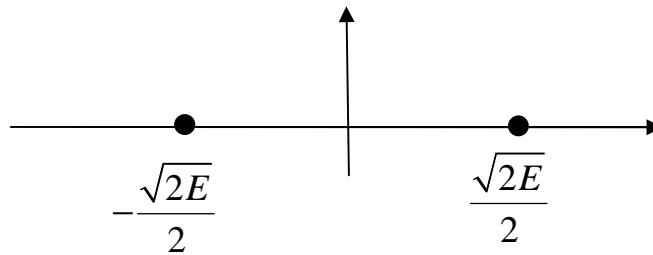
Orthogonal signaling

This is a two dimensional signaling scheme. The two signals are:

$s_1(t) = \sqrt{E}\varphi_1(t)$ and $s_2(t) = \sqrt{E}\varphi_2(t)$. The two basis functions are orthonormal.



The coordinates of the two vectors are $(\sqrt{E}, 0)$ and $(0, \sqrt{E})$. We can show that the probability of error is independent on translations and rotations in the signal space. So, under proper translation and rotation, the above signal space is equivalent to the following one:



The above figure is the same as the antipodal case. The probability of error is:

$$P(E) = \frac{1}{2} \operatorname{erfc} \left(\frac{A}{\sqrt{2}\sigma} \right)$$

where $A = \frac{\sqrt{2E}}{2} = \sqrt{\frac{E}{2}}$ and $\sigma^2 = \frac{N_0}{2}$

The probability of error is then:

$$P(E) = \frac{1}{2} \operatorname{erfc} \sqrt{\frac{E}{2N_0}} \quad (1.24)$$

An example of orthogonal signaling is the orthogonal FSK system. The receiver can be implemented as follows:

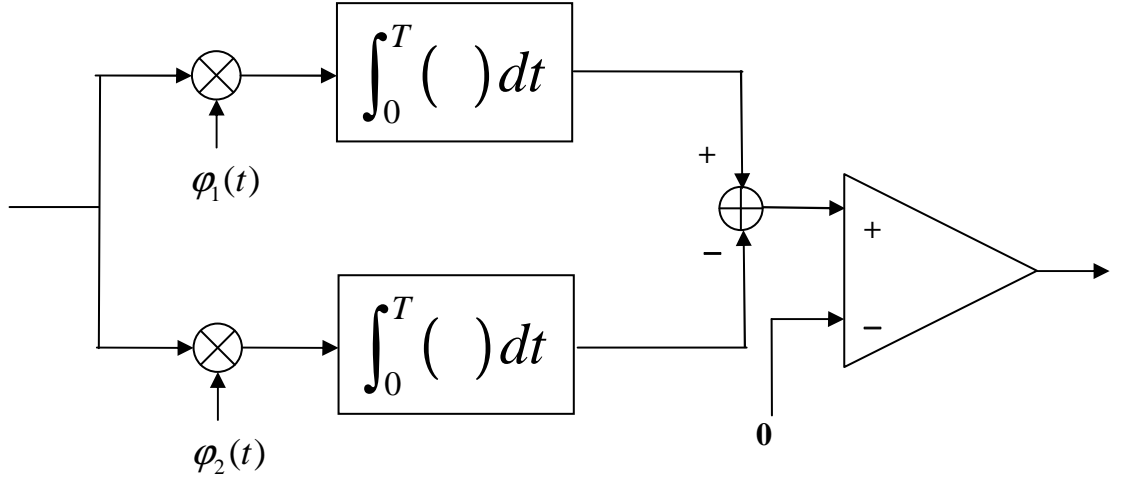


Fig.3- 7 orthogonal receiver structure

Non coherent FSK

In this communication system, the phase of the received waveform is unknown. It will be modeled as a uniformly distributed random variable over $[0, 2\pi]$. We assume also that the frequencies are selected so that the two carriers are orthogonal. The receiver observes one of the following two signals:

$$s_1(t) = A \cos(\omega_1 t + \Theta) \text{ and } s_2(t) = A \cos(\omega_2 t + \Theta)$$

Θ is a random variable uniformly distributed over $[0, 2\pi]$. The above representation corresponds to an infinite number of signals. Consider the following four orthonormal functions:

$$\varphi_1(t) = \sqrt{\frac{2}{T}} \cos \omega_1 t, \quad \varphi_2(t) = \sqrt{\frac{2}{T}} \sin \omega_1 t, \quad \varphi_3(t) = \sqrt{\frac{2}{T}} \cos \omega_2 t, \quad \varphi_4(t) = \sqrt{\frac{2}{T}} \sin \omega_2 t$$

and let θ be a realization of the random variable Θ . We can represent the above signals in the four dimensional space spanned by $\varphi_1, \varphi_2, \varphi_3, \varphi_4$. The corresponding vectors are:

$$\mathbf{s}_1 = \left(A \sqrt{\frac{T}{2}} \cos \theta, -A \sqrt{\frac{T}{2}} \sin \theta, 0, 0 \right)$$

$$\mathbf{s}_2 = \left(0, 0, A\sqrt{\frac{T}{2}} \cos \theta, -A\sqrt{\frac{T}{2}} \sin \theta \right)$$

The vector observed by the receiver is one of the above signals plus noise. We have seen that the components of noise are independent Gaussian random variables with zero mean and a variance equal to $\frac{N_0}{2}$. So, the four dimensional observed vector is the following Gaussian random vector: $\mathbf{z} = (z_1, z_2, z_3, z_4)$.

When ω_1 is transmitted and $\Theta = \theta$ is observed, \mathbf{z} has the following pdf:

$$f_{\mathbf{z}|\omega_1, \theta}(\mathbf{z} | \omega_1, \theta) = \frac{1}{(\pi N_0)^2} \exp \left\{ -\frac{1}{N_0} \left[\left(z_1 - A\sqrt{\frac{T}{2}} \cos \theta \right)^2 + \left(z_2 + A\sqrt{\frac{T}{2}} \sin \theta \right)^2 + z_3^2 + z_4^2 \right] \right\}$$

Using $E = \frac{A^2 T}{2}$, the above expression becomes:

$$f_{\mathbf{z}|\omega_1, \theta}(\mathbf{z} | \omega_1, \theta) = \frac{1}{(\pi N_0)^2} \exp \left\{ -\frac{1}{N_0} \left[\left(z_1 - \sqrt{E} \cos \theta \right)^2 + \left(z_2 + \sqrt{E} \sin \theta \right)^2 + z_3^2 + z_4^2 \right] \right\}$$

This expression becomes:

$$f_{\mathbf{z}|\omega_1, \theta}(\mathbf{z} | \omega_1, \theta) = \frac{1}{(\pi N_0)^2} \exp \left\{ -\frac{1}{N_0} \left[z_1^2 + z_2^2 + z_3^2 + z_4^2 + E - 2\sqrt{E} (z_1 \cos \theta - z_2 \sin \theta) \right] \right\}$$

When ω_2 is transmitted and $\Theta = \theta$ is observed, \mathbf{z} has the following pdf:

$$f_{\mathbf{z}|\omega_2, \theta}(\mathbf{z} | \omega_2, \theta) = \frac{1}{(\pi N_0)^2} \exp \left\{ -\frac{1}{N_0} \left[z_1^2 + z_2^2 + z_3^2 + z_4^2 + E - 2\sqrt{E} (z_3 \cos \theta - z_4 \sin \theta) \right] \right\}$$

A change of variable (from rectangular to polar) produces an expression of the pdf that is more informative.

So, let $z_1 = r_1 \cos \psi_1$, $z_2 = r_1 \sin \psi_1$, $z_3 = r_2 \cos \psi_2$ and $z_4 = r_2 \sin \psi_2$. The above pdf's become:

$$f(r_1, \psi_1, r_2, \psi_2 | \omega_1, \theta) = \frac{r_1 r_2}{(\pi N_0)^2} \exp \left\{ -\frac{1}{N_0} \left[r_1^2 + r_2^2 + E - 2\sqrt{E} r_1 \cos(\theta + \psi_1) \right] \right\}$$

$$f(r_1, \psi_1, r_2, \psi_2 | \omega_2, \theta) = \frac{r_1 r_2}{(\pi N_0)^2} \exp \left\{ -\frac{1}{N_0} \left[r_1^2 + r_2^2 + E - 2\sqrt{E} r_2 \cos(\theta + \psi_2) \right] \right\}$$

If we can measure the phase of the carrier, we can use the above expressions for an MAP receiver design. However, we don't have that

information. The solution is to take a decision after averaging the above pdf's over θ

$$\begin{aligned} f(r_1, \psi_1, r_2, \psi_2 | \omega_1) &= \frac{1}{2\pi} \int_0^{2\pi} f(r_1, \psi_1, r_2, \psi_2 | \omega_1, \theta) d\theta \\ &= \frac{r_1 r_2}{(\pi N_0)^2} \exp\left\{-\frac{1}{N_0} [r_1^2 + r_2^2 + E]\right\} \frac{1}{2\pi} \int_0^{2\pi} \exp\left[\frac{2\sqrt{E}}{N_0} r_1 \cos(\theta + \psi_1)\right] d\theta \end{aligned}$$

We use the fact that:

$$\frac{1}{2\pi} \int_0^{2\pi} \exp(z \cos \lambda) d\lambda = I_0(z), \quad I_0 \text{ is the modified Bessel function of the}$$

first kind of order zero. The pdf is:

$$f(r_1, \psi_1, r_2, \psi_2 | \omega_1) = \frac{r_1 r_2}{(\pi N_0)^2} \exp\left[-\frac{1}{N_0} (r_1^2 + r_2^2 + E)\right] I_0\left(\frac{2\sqrt{E}}{N_0} r_1\right) \quad (1.25)$$

When ω_2 is transmitted, the pdf becomes:

$$f(r_1, \psi_1, r_2, \psi_2 | \omega_2) = \frac{r_1 r_2}{(\pi N_0)^2} \exp\left[-\frac{1}{N_0} (r_1^2 + r_2^2 + E)\right] I_0\left(\frac{2\sqrt{E}}{N_0} r_2\right) \quad (1.26)$$

The maximum likelihood decision rule is then:

$$\text{"Decide } \omega_1 \text{ if } I_0\left(\frac{2\sqrt{E}}{N_0} r_1\right) > I_0\left(\frac{2\sqrt{E}}{N_0} r_2\right), \text{ decide } \omega_2 \text{ otherwise.}"$$

The function I_0 is a monotone increasing function of its argument. The decision rule can be simplified:

$$\text{"Decide } \omega_1 \text{ if } r_1 > r_2, \text{ decide } \omega_2 \text{ otherwise.}"$$

The receiver should compute r_1 and r_2 and compare them.

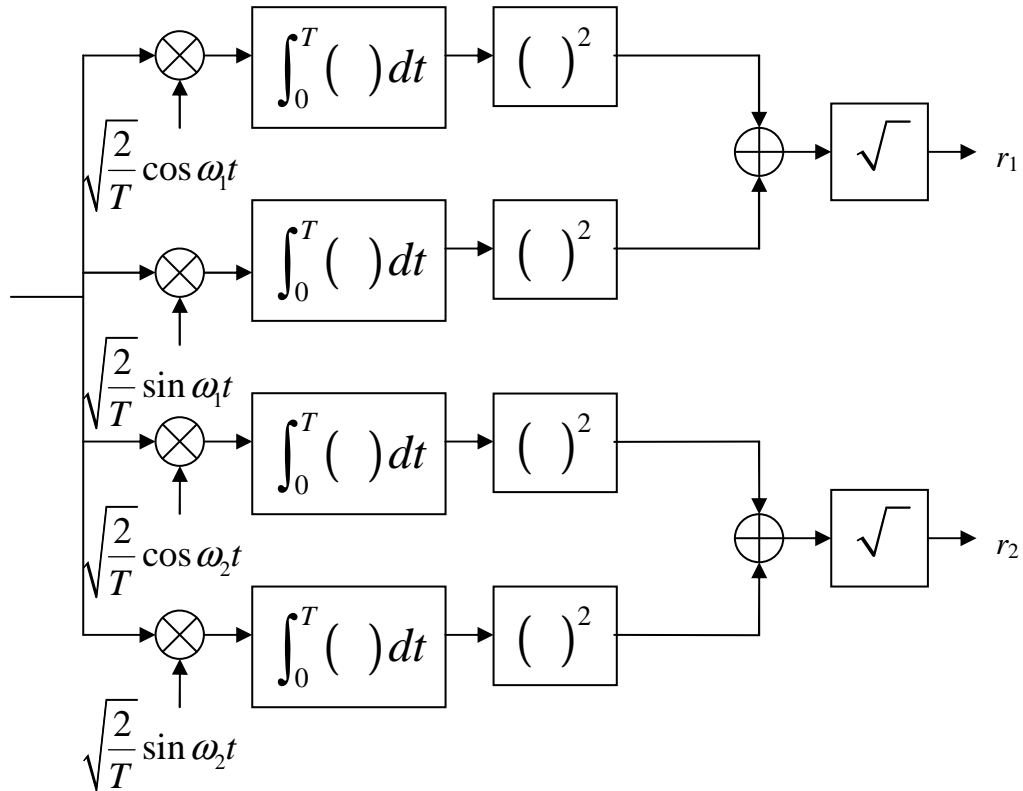


Fig.3- 8 Non coherent FSK receiver

We can show that the above structure can be implemented using two bandpass matched filters tuned respectively to ω_1 and ω_2 followed by envelop detectors. The impulse response of the bandpass filter matched to ω_1 is $h(t) = p(t)\cos(\omega_1 t + \theta)$ where $p(t) = 1$ for $0 \leq t \leq T$ and zero elsewhere and θ is arbitrary.

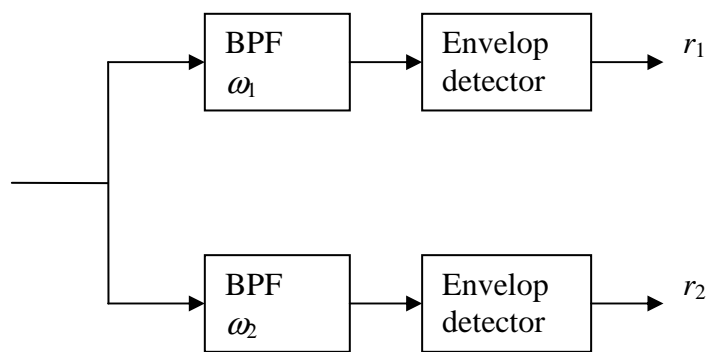
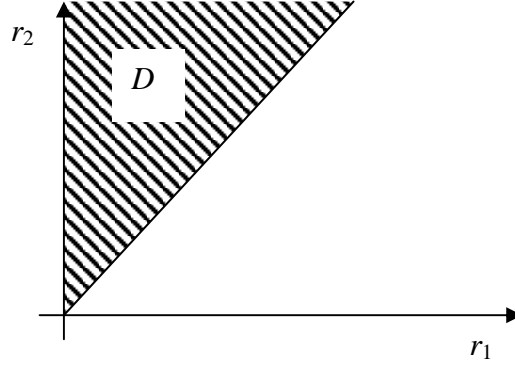


Fig.3- 9 Non coherent FSK receiver

The decision depends only on the amplitude of the output of the bandpass filters. The conditional probability of error can be computed as follows:

Given that ω_1 is transmitted, we have an error if r_2 is larger than r_1 . So:

$$P(E | \omega_1) = \iint_D f(r_1, r_2 | \omega_1) dr_1 dr_2$$



We need of course to compute the joint pdf of the two envelopes r_1 and r_2 .

$$\begin{aligned} f(r_1, r_2 | \omega_1) &= \int_0^{2\pi} \int_0^{2\pi} f(r_1, \psi_1, r_2, \psi_2) d\psi_1 d\psi_2 \\ &= \int_0^{2\pi} \int_0^{2\pi} \frac{r_1 r_2}{(\pi N_0)^2} \exp\left[-\frac{r_1^2 + r_2^2 + E}{N_0}\right] I_0\left(\frac{2\sqrt{E}}{N_0} r_1\right) d\psi_1 d\psi_2 \end{aligned}$$

Finally:

$$f(r_1, r_2 | \omega_1) = \frac{4r_1 r_2}{N_0^2} \exp\left[-\frac{r_1^2 + r_2^2 + E}{N_0}\right] I_0\left(\frac{2\sqrt{E}}{N_0} r_1\right) \quad (1.27)$$

From the above expression, we can see that the two variables (r_1 and r_2) are independent and their pdf's are:

$$f(r_1 | \omega_1) = \frac{2r_1}{N_0} \exp\left[-\frac{r_1^2 + E}{N_0}\right] I_0\left(\frac{2\sqrt{E}}{N_0} r_1\right) \quad (1.28)$$

This means that r_1 is Rician.

$$f(r_2 | \omega_1) = \frac{2r_2}{N_0} \exp\left[-\frac{r_2^2}{N_0}\right] \quad (1.29)$$

This means that r_2 is Rayleigh.

The conditional probability of error is:

$$P(E | \omega_1) = \int_0^{+\infty} \left[\int_{r_1}^{+\infty} f(r_1 | \omega_1) f(r_2 | \omega_2) dr_2 \right] dr_1$$

After some simple manipulations, we obtain:

$$P(E | \omega_1) = \frac{1}{2} \exp\left[-\frac{E}{2N_0}\right]$$

The other probability of error has the same expression. Since the two symbols are equiprobable, the probability of error has the same expression also.

$$P(E) = \frac{1}{2} \exp\left[-\frac{E}{2N_0}\right] \quad (1.30)$$