

# Communications Toolbox

For Use with MATLAB®

Computation  
└─

Visualization  
└─

Programming  
└─



User's Guide

*Version 2*

## How to Contact The MathWorks:



508-647-7000

Phone



508-647-7001

Fax



The MathWorks, Inc.  
3 Apple Hill Drive  
Natick, MA 01760-2098

Mail



<http://www.mathworks.com>  
<ftp.mathworks.com>  
<comp.soft-sys.matlab>

Web  
Anonymous FTP server  
Newsgroup



[support@mathworks.com](mailto:support@mathworks.com)  
[suggest@mathworks.com](mailto:suggest@mathworks.com)  
[bugs@mathworks.com](mailto:bugs@mathworks.com)  
[doc@mathworks.com](mailto:doc@mathworks.com)  
[subscribe@mathworks.com](mailto:subscribe@mathworks.com)  
[service@mathworks.com](mailto:service@mathworks.com)  
[info@mathworks.com](mailto:info@mathworks.com)

Technical support  
Product enhancement suggestions  
Bug reports  
Documentation error reports  
Subscribing user registration  
Order status, license renewals, passcodes  
Sales, pricing, and general information

### *Communications Toolbox User's Guide*

© COPYRIGHT 1996 - 2000 by The MathWorks, Inc.

The software described in this document is furnished under a license agreement. The software may be used or copied only under the terms of the license agreement. No part of this manual may be photocopied or reproduced in any form without prior written consent from The MathWorks, Inc.

**FEDERAL ACQUISITION:** This provision applies to all acquisitions of the Program and Documentation by or for the federal government of the United States. By accepting delivery of the Program, the government hereby agrees that this software qualifies as "commercial" computer software within the meaning of FAR Part 12.212, DFARS Part 227.7202-1, DFARS Part 227.7202-3, DFARS Part 252.227-7013, and DFARS Part 252.227-7014. The terms and conditions of The MathWorks, Inc. Software License Agreement shall pertain to the government's use and disclosure of the Program and Documentation, and shall supersede any conflicting contractual terms or conditions. If this license fails to meet the government's minimum needs or is inconsistent in any respect with federal procurement law, the government agrees to return the Program and Documentation, unused, to MathWorks.

MATLAB, Simulink, Stateflow, Handle Graphics, and Real-Time Workshop are registered trademarks, and Target Language Compiler is a trademark of The MathWorks, Inc.

Other product or brand names are trademarks or registered trademarks of their respective holders.

Printing History:	April 1996	First printing	New
	May 1997	Second printing	Revised for MATLAB 5
	September 2000	Third printing	Revised for Version 2 (Release 12)

## Preface

<b>What Is the Communications Toolbox? .....</b>	<b>viii</b>
<b>Related Products .....</b>	<b>ix</b>
<b>Using This Guide .....</b>	<b>xi</b>
Expected Background .....	xi
Supplementing This Guide with Command-Line Help .....	xii
<b>Configuration Information .....</b>	<b>xiii</b>
<b>Technical Conventions .....</b>	<b>xiv</b>
Polynomials as Vectors .....	xiv
Matrices .....	xiv
<b>Typographical Conventions .....</b>	<b>xv</b>

## Getting Started with the Communications Toolbox

1

<b>A Detailed Example .....</b>	<b>1-2</b>
What the Example Does .....	1-2
Where to Find the Example .....	1-3
How the Example Works .....	1-3
Output from the Example .....	1-6

<b>Random Signals and Error Analysis</b> .....	<b>2-3</b>
Error Analysis Features of the Toolbox .....	<b>2-3</b>
Random Signals .....	<b>2-3</b>
Error Rates .....	<b>2-7</b>
Eye Diagrams .....	<b>2-8</b>
Scatter Plots .....	<b>2-11</b>
 <b>Source Coding</b> .....	 <b>2-14</b>
Source Coding Features of the Toolbox .....	<b>2-14</b>
Representing Quantization Parameters .....	<b>2-14</b>
Quantizing a Signal .....	<b>2-15</b>
Optimizing Quantization Parameters .....	<b>2-18</b>
Implementing Differential Pulse Code Modulation .....	<b>2-19</b>
Optimizing DPCM Parameters .....	<b>2-21</b>
Companding a Signal .....	<b>2-22</b>
Selected Bibliography for Source Coding .....	<b>2-23</b>
 <b>Block Coding</b> .....	 <b>2-24</b>
Block Coding Features of the Toolbox .....	<b>2-25</b>
Block Coding Terminology .....	<b>2-26</b>
Representing Messages and Codewords .....	<b>2-26</b>
Representing Block Coding Parameters .....	<b>2-30</b>
Creating and Decoding Block Codes .....	<b>2-35</b>
Performing Other Block Code Tasks .....	<b>2-40</b>
Selected Bibliography for Block Coding .....	<b>2-42</b>
 <b>Convolutional Coding</b> .....	 <b>2-43</b>
Convolutional Coding Features of the Toolbox .....	<b>2-43</b>
Polynomial Description of a Convolutional Encoder .....	<b>2-43</b>
Trellis Description of a Convolutional Encoder .....	<b>2-46</b>
Creating and Decoding Convolutional Codes .....	<b>2-50</b>
Examples of Convolutional Coding .....	<b>2-52</b>
Selected Bibliography for Convolutional Coding .....	<b>2-55</b>
 <b>Modulation</b> .....	 <b>2-56</b>
Modulation Features of the Toolbox .....	<b>2-57</b>
Modulation Terminology .....	<b>2-58</b>

Representing Analog Signals .....	2-59
Simple Analog Modulation Example .....	2-61
Other Options in Analog Modulation .....	2-62
Filter Design Issues .....	2-62
Digital Modulation Overview .....	2-66
Representing Digital Signals .....	2-67
Significance of Sampling Rates .....	2-70
Representing Signal Constellations .....	2-70
Simple Digital Modulation Example .....	2-74
Customizing the Modulation Process .....	2-75
Other Options in Digital Modulation .....	2-77
Selected Bibliography for Modulation .....	2-77
<b>Special Filters .....</b>	<b>2-78</b>
Special Filter Features of the Toolbox .....	2-78
Noncausality and the Group Delay Parameter .....	2-78
Designing Hilbert Transform Filters .....	2-80
Filtering with Raised Cosine Filters .....	2-81
Designing Raised Cosine Filters .....	2-87
Selected Bibliography for Special Filters .....	2-88
<b>Galois Field Computations .....</b>	<b>2-89</b>
Galois Field Features of the Toolbox .....	2-89
Galois Field Terminology .....	2-89
Representing Elements of Galois Fields .....	2-90
Default Primitive Polynomials .....	2-93
Converting and Simplifying Element Formats .....	2-94
Arithmetic in Galois Fields .....	2-97
Polynomials over Prime Fields .....	2-99
Other Galois Field Functions .....	2-103
Selected Bibliography for Galois Fields .....	2-103

<b>Functions by Category</b> .....	<b>3-3</b>
<b>Alphabetical List of Functions</b> .....	<b>3-9</b>
ademod .....	<b>3-12</b>
ademodce .....	<b>3-16</b>
amod .....	<b>3-20</b>
amodce .....	<b>3-25</b>
apkconst .....	<b>3-28</b>
awgn .....	<b>3-32</b>
bchdeco .....	<b>3-34</b>
bchenco .....	<b>3-36</b>
bchpoly .....	<b>3-37</b>
bi2de .....	<b>3-41</b>
biterr .....	<b>3-43</b>
compand .....	<b>3-49</b>
convenc .....	<b>3-51</b>
cyclgen .....	<b>3-53</b>
cyclpoly .....	<b>3-55</b>
ddemod .....	<b>3-57</b>
ddemodce .....	<b>3-62</b>
de2bi .....	<b>3-67</b>
decode .....	<b>3-69</b>
demodmap .....	<b>3-73</b>
dmod .....	<b>3-78</b>
dmodce .....	<b>3-82</b>
dpcmdeco .....	<b>3-86</b>
dpcmenco .....	<b>3-87</b>
dpcmopt .....	<b>3-88</b>
encode .....	<b>3-89</b>
eyediagram .....	<b>3-95</b>
gen2par .....	<b>3-97</b>
gfadd .....	<b>3-99</b>
gfconv .....	<b>3-101</b>
gfcosets .....	<b>3-103</b>
gfdeconv .....	<b>3-105</b>
gfdiv .....	<b>3-108</b>
gffilter .....	<b>3-110</b>

gflneq	3-112
gfminpol	3-114
gfmul	3-116
gfplus	3-117
gfpretty	3-118
gfprimck	3-120
gfprimdf	3-121
gfprimfd	3-122
gfrank	3-124
gfrepcov	3-125
gfroots	3-126
gfsub	3-128
gftrunc	3-130
gftuple	3-131
gfweight	3-134
hamngen	3-135
hank2sys	3-138
hilbiir	3-140
istrellis	3-143
lloyds	3-145
marcumq	3-148
modmap	3-149
oct2dec	3-154
poly2trellis	3-155
qaskdeco	3-158
qaskenco	3-160
quantiz	3-163
randerr	3-165
randint	3-167
randsrc	3-168
rcosfir	3-170
rcosflt	3-172
rcosiir	3-175
rcosine	3-177
rsdeco	3-179
rsdecode	3-182
rsdecof	3-184
rsenco	3-185
rsencode	3-188
rsencof	3-190

rspoly .....	<b>3-191</b>
scatterplot .....	<b>3-193</b>
symerr .....	<b>3-195</b>
syndtable .....	<b>3-198</b>
vec2mat .....	<b>3-199</b>
vitdec .....	<b>3-201</b>
wgn .....	<b>3-205</b>



# Preface

---

<b>What Is the Communications Toolbox?</b>	viii
<b>Related Products</b>	ix
<b>Using This Guide</b>	xi
Expected Background	xi
Supplementing This Guide with Command-Line Help	xii
<b>Configuration Information</b>	xiii
<b>Technical Conventions</b>	xiv
Polynomials as Vectors	xiv
Matrices	xiv
<b>Typographical Conventions</b>	xv

## What Is the Communications Toolbox?

The Communications Toolbox is a set of MATLAB<sup>®</sup> functions that can help you design and analyze advanced communication systems. Functions in the toolbox can accomplish these tasks:

- Random signal production
- Error analysis, including eye diagrams and scatter plots
- Source coding, including scalar quantization, differential pulse code modulation, and companders
- Error-control coding, including convolutional and linear block coding
- Analog and digital modulation/demodulation
- Filtering of data using special filters
- Computations in Galois fields

## Related Products

The MathWorks provides several products that are especially relevant to the kinds of tasks you can perform with the Communications Toolbox. They are listed in the table below. In particular, the Communications Toolbox *requires* these products:

- MATLAB
- Signal Processing Toolbox

For more information about any of these products, see either:

- The online documentation for that product, if it is installed or if you are reading the documentation from the CD
- The MathWorks Web site, at <http://www.mathworks.com>; see the “products” section

---

**Note** The toolboxes listed below all include functions that extend MATLAB’s capabilities. The blocksets all include blocks that extend the capabilities of Simulink®.

---

Product	Description
CDMA Reference Blockset	Simulink block libraries for the design and simulation of the IS-95A wireless communications standard
Communications Blockset	Simulink block libraries for modeling the physical layer of communications systems
DSP Blockset	Simulink block libraries for the design, simulation, and prototyping of digital signal processing systems

Product	Description
Signal Processing Toolbox	Tool for algorithm development, signal and linear system analysis, and time-series data modeling
Simulink	Interactive, graphical environment for modeling, simulating, and prototyping dynamic systems

## Using This Guide

This guide describes and illustrates the capabilities of the Communications Toolbox. The table below matches sections of this guide with your possible learning goals.

Goal	Section
Examine an example in detail, to begin learning about the toolbox	Chapter 1, “Getting Started with the Communications Toolbox”
Learn how this toolbox implements a particular category of functionality, such as source coding	Chapter 2, “Using the Communications Toolbox”
Learn about particular functions in this toolbox	Chapter 3, “Reference”

### Expected Background

This guide assumes that you already have background knowledge in the subject of communications. If you do not yet have this background, then you can acquire it using a standard communications text or the books listed in one of this guide’s sections entitled “Selected Bibliography for... .”

#### For New Users

Start with Chapter 1, “Getting Started with the Communications Toolbox”, which describes an example in detail. Then read those parts of Chapter 2, “Using the Communications Toolbox” that address the functionality that concerns you. When you find out from that chapter which functions you want to use, refer to the references pages in Chapter 3, “Reference” that describe those functions.

#### For Experienced Users

The reference descriptions in Chapter 3, “Reference” are probably the most relevant parts of this guide for you. Each reference description includes the function’s syntax as well as a complete explanation of its options and operation. Many reference descriptions also include examples, a description of the function’s algorithm, and references to additional reading material.

You might also want to browse through Chapter 1, “Getting Started with the Communications Toolbox” and Chapter 2, “Using the Communications Toolbox” based on your interests or needs.

## Supplementing This Guide with Command-Line Help

Command-line help is text that MATLAB displays in its command window. The table below lists two kinds of command-line help that are available for the Communications Toolbox, along with the command that you would type at the MATLAB prompt in order to display the help text.

Type of Command-Line Help	MATLAB Command
List of functions in the Communications Toolbox	<code>hel p comm</code>
Information about a particular function	<code>hel p <i>function</i></code> (for example, <code>hel p ademod</code> )

### Method-Specific Help

Some multipurpose functions also provide command-line help on specific methods. For example, `hel p encode` displays text that describes the use of the `encode` command for error-control encoding. One specific method of error-control encoding is BCH encoding. The command

`encode bch`

displays text that describes the use of the `encode` command for BCH encoding. The functions that provide method-specific help are: `amod`, `ademod`, `amodce`, `ademodce`, `ddemod`, `ddemodce`, `decode`, `demodmap`, `dmod`, `dmodce`, `encode`, and `modmap`. The general help text, displayed by the `hel p function` command, lists the available methods.

## Configuration Information

To determine if the Communications Toolbox is installed on your system, type

`ver`

at the MATLAB prompt. MATLAB displays information about the version of MATLAB you are running, including a list of all toolboxes installed on your system and their version numbers. Check the list to see if the Communications Toolbox appears.

For information about installing the toolbox, see the *MATLAB Installation Guide* for your platform.

# Technical Conventions

This section mentions some technical conventions that this guide uses.

## Polynomials as Vectors

MATLAB represents a polynomial in one variable  $x$  using a vector that lists the polynomial's coefficients, arranged according to the powers of  $x$ . *Descending order* means that the coefficient of the highest power of  $x$  appears first and that the polynomial's constant term appears last. *Ascending order* is the opposite. The table below illustrates the conventions for functions in this toolbox and for built-in MATLAB functions.

Category of Functions	Vector That Represents the Polynomial $1+2x+3x^2$
Error-control coding or Galois field computations	[1, 2, 3] (ascending order)
Modulation/demodulation, e.g., when using filters	[3, 2, 1] (descending order)
Built-in MATLAB, e.g., roots, poly, polyval	[3, 2, 1] (descending order)

## Matrices

Matrix dimensions are described by listing the number of rows and the number of columns of the matrix in that order, as below.

```
u = [ 1 2 3; 4 5 6] % A 2-by-3 matrix
```



# Typographical Conventions

This guide uses some or all of these conventions.

Item	Convention to Use	Example
Example code	Monospace font	To assign the value 5 to A, enter <code>A = 5</code>
Function names/syntax	Monospace font	The cos function finds the cosine of each array element. Syntax line example is <code>MLGetVar ML_var_name</code>
Keys	<b>Boldface</b> with an initial capital letter	Press the <b>Return</b> key.
Literal strings (in syntax descriptions in Reference chapters)	<b>Monospace bold</b> for literals.	<code>f = freqspace(n, 'whole')</code>
Mathematical expressions	Variables in <i>italics</i> Functions, operators, and constants in standard text.	This vector represents the polynomial $p = x^2 + 2x + 3$
MATLAB output	Monospace font	MATLAB responds with <code>A =</code> <code>5</code>
Menu names, menu items, and controls	<b>Boldface</b> with an initial capital letter	Choose the <b>File</b> menu.
New terms	<i>Italics</i>	An <i>array</i> is an ordered collection of information.
String variables (from a finite list)	<i>Monospace italics</i>	<code>sysc = d2c(sysd, 'method')</code>



# Getting Started with the Communications Toolbox

---

<b>A Detailed Example . . . . .</b>	<b>1-2</b>
What the Example Does . . . . .	1-2
Where to Find the Example . . . . .	1-3
How the Example Works . . . . .	1-3
Output from the Example . . . . .	1-6

# A Detailed Example

This chapter describes a particular example in detail, to help you get started using the Communications Toolbox. It uses several functions from the toolbox, as the table below indicates.

Function	Purpose in Example
randi nt	Generate a random signal
dmodce	Modulate signals
ddemodce	Demodulate signals
bi terr	Compute bit error rate
modmap	Plot a signal constellation

This chapter assumes very little about your prior knowledge of MATLAB, although it still assumes that you have a basic knowledge about communications subject matter.

## What the Example Does

The example creates a random digital signal consisting of integers between 0 and 8, and modulates it using two varieties of the 8-ary quadrature amplitude shift keying (QASK) technique. This technique associates each integer in the signal with some point in an eight-point signal constellation, and then uses the associations to create a modulated signal.

There are  $8!$ , that is, `factorial (8)`, ways to associate eight symbols with eight constellation points. One category of configurations implements what is called Gray coding. In a Gray coded constellation, the symbol associated with a given point and the symbol of any of the point's nearest neighbors differ in exactly one bit. Thus, the constellation point associated with the symbol 3 (= 011) can have as a nearest neighbor the point associated with the symbol 1 (= 001), 2 (= 010), or 7 (= 111), but not any other number.

In order to compare the behavior of different constellation configurations, the example modulates the message signal separately using two varieties of 8-QASK modulation. Both varieties use constellations with the same points,

but one variety labels the constellation points so as to implement Gray code while the other variety does not implement Gray code. After modulating, the example adds noise to both modulated signals, demodulates both noisy signals, and compares the bit error rates in the two cases.

The example outputs the two bit error rates. The expectation is that although noise might cause demodulation errors in both cases, the errors in the Gray coding case should involve fewer bits. When you execute the example, check to see whether the bit error rate from the Gray coding case is smaller than the bit error rate from the non-Gray coding case.

## Where to Find the Example

If you have already installed MATLAB and the Communications Toolbox, then the toolbox will be there whenever you start up MATLAB. The example is contained in a file called `commgettingstarted.m`, which is located in the `toolbox/comm/commdemos` directory within your MATLAB installation. You can view the contents of the example file by typing

```
type commgettingstarted
```

at the MATLAB prompt.

You can execute the example by typing

```
commgettingstarted
```

at the MATLAB prompt.

## How the Example Works

This section displays and explains the example code, piece by piece.

### Setting Up Parameters

The first part of the example defines variables that the rest of the example will use. The symbol alphabet has  $M$  different symbols, namely, the integers between 0 and  $M-1$ . The message will be a column vector having  $len$  entries, each of which is chosen from the symbol alphabet.

The variables  $F_d$  and  $F_s$  refer to the relative sampling rates for the modulation scheme. They would be more meaningful if the example were sampling a real signal that had a natural notion of time. However, since this example uses a random signal that does not have a built-in notion of time, the main purpose of

$F_d$  and  $F_s$  is to indicate that the modulated signal has three entries for every one entry of the original signal.

```
% Set up parameters.  
M = 8; % Number of symbols in alphabet  
len = 10000; % Number of symbols in the original message  
F_d = 1; % Assume the original message is sampled  
% at a rate of 1 sample per second.  
F_s = 3; % The modulated signal will be sampled  
% at a rate of 3 samples per second.
```

### Creating the Signal

The variable `signal` is a `len-by-1` matrix, that is, a column vector of length `len`, whose entries are randomly chosen integers between 0 and  $M-1$ . This is the signal that the example will modulate. The `randint` function is part of this toolbox.

```
% Create a signal.  
signal = randint(len, 1, M); % Random digital message  
% consisting of integers between 0 and M-1
```

### Modulating the Signal

This part of the example modulates the data in the column vector `signal` in two different ways. The `dmodce` function performs both modulations and puts the results into the two-column matrix `modsignal`.

The first call to `dmodce`, which creates the first column of `modsignal`, tells `dmodce` to use QASK modulation on  $M$  symbols. The string 'qask' indicates the QASK method as well as the default square constellation configuration. In this case, the configuration implements Gray code.

The second call to `dmodce`, which creates the second column of `modsignal`, tells `dmodce` to use QASK modulation with a signal constellation whose configuration is represented in the vectors `inphase` and `quad`. The variables `inphase` and `quad` are length- $M$  vectors that list the in-phase and quadrature components, respectively, of the points in the signal constellation. The points are listed in sequence, to associate a message symbol of  $k$  with the  $(k+1)$ st elements in `inphase` and `quad`. Whereas Gray code labels the constellation points in a special way, this configuration lists points in a sequence that is merely convenient for creating `inphase` and `quad`.

These lines also illustrate some common ways to manipulate matrices in MATLAB. If you are not familiar with MATLAB's colon notation or with functions like `ones` and `zeros`, then you should consult the MATLAB documentation set.

```
% Use M-ary QASK modulation with two different labeled
% square constellations.
modsi gnal (:, 1) = dmodce(sig nal, Fd, Fs, 'qask', M);
inphase = [-3:2:3 -3:2:3];
quad = [ones(1, 4), -1*ones(1, 4)];
modsi gnal (:, 2) = dmodce(sig nal, Fd, Fs, 'qask/arb', inphase, quad);
```

### Adding Noise

According to the definition of baseband QASK modulation, `modsi gnal` is a *complex* matrix having `len*Fs/Fd` rows and two columns. The command below adds normally distributed random numbers to the real and imaginary parts of `modsi gnal`, to produce a noisy signal `noi sy`. The `randn` function is a built-in MATLAB function.

Notice that the command adds to `modsi gnal` an entire real matrix of the appropriate size and an entire imaginary matrix of the appropriate size. Using a loop to add noise to individual scalar entries of `modsi gnal` would be less efficient, since MATLAB is optimized for matrix operations.

```
% Add noise to real and imaginary parts of the modulated signal.
noi sy = modsi gnal + .5*randn(len*Fs/Fd, 2) ...
+j*.5*randn(len*Fs/Fd, 2);
```

### Demodulating the Signal

This part of the example demodulates the noisy modulated signal, `noi sy`, in two different ways. The `ddemodce` function performs both demodulations by operating on each column of `noi sy` separately. In each case, `ddemodce` puts the results into the two-column matrix `newsi gnal`.

```
% Demodulate to recover the message.
newsi gnal (:, 1) = ddemodce(noi sy(:, 1), Fd, Fs, 'qask', M);
newsi gnal (:, 2) = ddemodce(noi sy(:, 2), Fd, Fs, ...
'qask/arb', inphase, quad);
```

## Computing and Displaying Bit Error Rates

The `biterr` function compares each demodulated signal (that is, each column of `newsignal`) to the original signal. Then `biterr` computes the number of bit errors, as well as the rate or fraction of bit errors. The built-in MATLAB function `disp` displays the two bit error rates in the command window.

```
% Check whether Gray code resulted in fewer bit errors.
% Compare signal with each column of newsignal.
[num, rate] = biterr(newsignal, signal);
disp('Bit error rates for the two constellations used here')
disp('-----')
disp(['Gray code constellation:      ', num2str(rate(1))])
disp(['Non-Gray code constellation: ', num2str(rate(2))])
```

## Plotting a Signal Constellation

The `modmap` function plots and labels the default square signal constellation having `M` points. The constellation that `inphase` and `quad` determine looks the same, except that the points are labeled from left to right across each row in the diagram, starting with the upper row.

```
% Plot signal constellations with Gray code labeling.
modmap('qask', M);
```

## Output from the Example

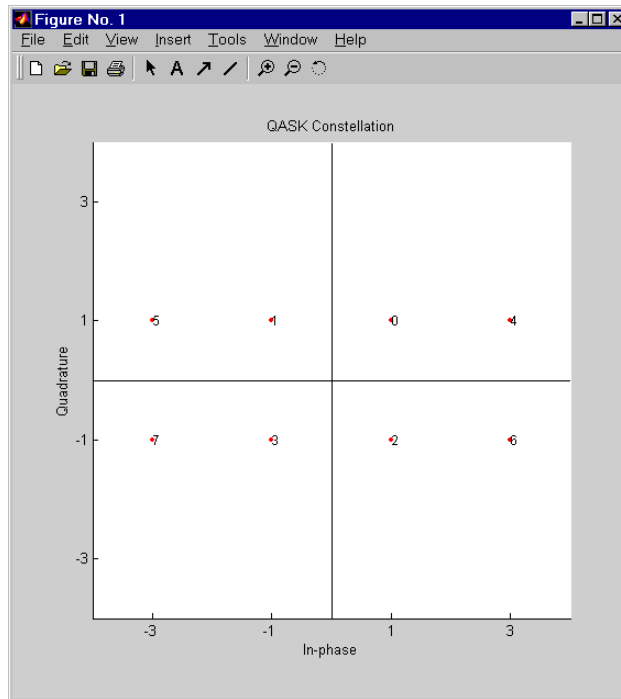
The example produces output in the command window like that shown below. Since the message signal and the noise are random, you will probably not get the exact numbers below. (For information about states and repeatable sequences of random numbers, see the reference page for the built-in MATLAB function `rand`.)

```
Bit error rates for the two constellations used here
-----
Gray code constellation:      0.0003
Non-Gray code constellation: 0.00036667
```

The example also produces a figure window containing the signal constellation plot in the figure below. The horizontal axis represents the in-phase components and the vertical axis represents the quadrature components. The dots are the constellation points. The number next to each dot is the message symbol associated with that dot. By considering the binary form of each



number from 0 to  $M-1$ , you can check that this constellation implements Gray code.



**Figure 1-1: Square 8-ary QASK Signal Constellation, Labeled for Gray Code**



# Using the Communications Toolbox

---

<b>Random Signals and Error Analysis</b>	2-3
<b>Source Coding</b>	2-14
<b>Block Coding</b>	2-24
<b>Convolutional Coding</b>	2-43
<b>Modulation</b>	2-56
<b>Special Filters</b>	2-78
<b>Galois Field Computations</b>	2-89

A typical communication system includes a signal source, sink, and channel, as well as processes for transmitting and receiving. This chapter describes and illustrates how to implement communication components using the functions provided in the Communications Toolbox. Each section in this chapter corresponds to a category of functionality within the Communications Toolbox. The sections are:

- “Random Signals and Error Analysis” on page 2-3
- “Source Coding” on page 2-14
- “Block Coding” on page 2-24
- “Convolutional Coding” on page 2-43
- “Modulation” on page 2-56
- “Special Filters” on page 2-78
- “Galois Field Computations” on page 2-89

## Random Signals and Error Analysis

Simulating a communication system often involves analyzing its response to the noise inherent in real-world components. Such analysis aims to illustrate the system's response and possibly to help design a system appropriate for the most likely kinds of noise.

### Error Analysis Features of the Toolbox

Error analysis tasks supported in the Communications Toolbox include:

- Simulating noise or signal sources using random signals
- Computing the error rate or number of errors
- Plotting an eye diagram
- Generating a scatter plot

This section describes these toolbox functions that accomplish error-analysis tasks: `biterr`, `eyediagram`, `randerr`, `randint`, `randsrc`, `scatterplot`, `symerr`, and `wgn`. Since error analysis is often a component of communication system simulation, other portions of this guide provide additional examples.

### Random Signals

Random signals are useful for simulating noise, errors, or signal sources. Besides built-in MATLAB functions like `rand` and `randn`, you can also use these functions from this toolbox:

- `wgn`, for generating white Gaussian noise
- `randsrc`, for generating random symbols
- `randint`, for generating uniformly distributed random integers
- `randerr`, for generating random bit error patterns

While `randsrc` and `randint` are suitable for representing sources, `randerr` is more appropriate for modeling channel errors.

#### White Gaussian Noise

The `wgn` function generates random matrices using a white Gaussian noise distribution. You specify the power of the noise in either dB (decibels), dBm, or linear units. You can generate either real or complex noise.

For example, the command below generates a column vector of length 50 containing real white Gaussian noise whose power is 2 dB. The function assumes that the load impedance is 1 Ohm.

```
y1 = wgn(50, 1, 2);
```

To generate complex white Gaussian noise whose power is 2 watts, across a load of 60 Ohms, use either of the commands below. Notice that the ordering of the string inputs does not matter.

```
y2 = wgn(50, 1, 2, 60, 'complex', 'linear');
y3 = wgn(50, 1, 2, 60, 'linear', 'complex');
```

To send a signal through an additive white Gaussian noise channel, use the `awgn` function.

### Random Symbol Matrices

The `randsrc` function generates random matrices whose entries are chosen independently from an alphabet that you specify, with a distribution that you specify. A special case generates bipolar matrices.

For example, the command below generates a 5-by-4 matrix whose entries are independently chosen and uniformly distributed in the set {1,3,5}. (Your results may vary because these are random numbers.)

```
a = randsrc(5, 4, [1, 3, 5])
```

```
a =
```

```

3     5     1     5
1     5     3     3
1     3     3     1
1     1     3     5
3     1     1     3
```

If you want 1 to be twice as likely to occur as either 3 or 5, then use the command below to prescribe the skewed distribution. Notice that the third input argument has two rows, one of which indicates the possible values of `b` and the other indicates the probability of each value.

```
b = randsrc(5, 4, [1, 3, 5; .5, .25, .25])
```

b =

3	3	5	1
1	1	1	1
1	5	1	1
1	3	1	3
3	1	3	1

### Random Integer Matrices

The `randint` function generates random integer matrices whose entries are in a range that you specify. A special case generates random binary matrices.

For example, the command below generates a 5-by-4 matrix containing random integers between 2 and 10.

```
c = randint(5, 4, [2, 10])
```

c =

2	4	4	6
4	5	10	5
9	7	10	8
5	5	2	3
10	3	4	10

If your desired range is [0,10] instead of [2,10] then you can use either of the commands below. They produce different numerical results, but use the same distribution.

```
d = randint(5, 4, [0, 10]);
```

```
e = randint(5, 4, 11);
```

### Random Bit Error Patterns

The `randerr` function generates matrices whose entries are either 0 or 1. However, its options are rather different from those of `randint`, since `randerr` is meant for testing error-control coding. For example, the command below generates a 5-by-4 binary matrix having the property that each row contains exactly one 1.

```
f = randerr(5, 4)
```

f =

0	0	1	0
0	0	1	0
0	1	0	0
1	0	0	0
0	0	1	0

You might use such a command to perturb a binary code that consists of five four-bit codewords. Adding the random matrix *f* to your code matrix (modulo 2) would introduce exactly one error into each codeword.

On the other hand, if you want to perturb each codeword by introducing one error with probability 0.4 and two errors with probability 0.6, then the command below should replace the one above.

```
% Each row has one '1' with probability 0.4, otherwise two '1's
g = randerr(5, 4, [1, 2; 0.4, 0.6])
```

g =

0	1	1	0
0	1	0	0
0	0	1	1
1	0	1	0
0	1	1	0

---

**Note** The probability matrix that is the third argument of `randerr` affects only the *number* of 1s in each row, not their placement.

---

As another application, you can generate an equiprobable binary 100-element column vector using any of the commands below. The three commands produce different numerical outputs, but use the same *distribution*. Notice that the third input arguments vary according to each function's particular way of specifying its behavior.

```
binarymatrix1 = randsrc(100, 1, [0 1]); % Possible values are 0, 1.
binarymatrix2 = randint(100, 1, 2); % Two possible values
binarymatrix3 = randerr(100, 1, [0 1; .5 .5]); % No 1s, or one 1
```



## Error Rates

Comparing messages before and after transmission can help you evaluate the quality of a communication system design or the performance of a special technique or algorithm. If your communication system uses several bits to represent a single symbol, then counting bit errors is different from counting symbol errors. In either the bit- or symbol-counting case, the error rate is the number of errors divided by the total number (of bits or symbols) transmitted.

The `biterr` function compares two messages and computes the number of bit errors and the bit error rate. The `symerr` function compares two messages and computes the number of symbol errors and the symbol error rate.

### Example: Computing Error Rates

The script below uses the `symerr` function to compute the symbol error rates for a noisy linear block code. After artificially adding noise to the encoded message, it compares the resulting noisy code to the original code. Then it decodes and compares the decoded message to the original one.

```
m = 3; n = 2^m-1; k = n-m; % Prepare to use Hamming code.
msg = randint(k*200, 1, 2); % 200 messages of k bits each
code = encode(msg, n, k, 'hamming');
codenoisy = rem(code+(rand(n*200, 1)>.95), 2); % Add noise.
% Decode and correct some errors.
newmsg = decode(codenoisy, n, k, 'hamming');
% Compute and display symbol error rates.
[codenum, coderate] = symerr(code, codenoisy);
[msgnum, msgrate] = symerr(msg, newmsg);
disp(['Error rate in the received code: ', num2str(coderate)])
disp(['Error rate after decoding: ', num2str(msgrate)])
```

The output is below. The error rate decreases after decoding because the Hamming decoder corrects some of the errors. Your results might vary because the example uses random numbers.

```
Error rate in the received code: 0.054286
Error rate after decoding: 0.03
```

### Comparison of Symbol Error Rate and Bit Error Rate

In the example above, the symbol errors and bit errors are the same because each symbol is a bit. The commands below illustrate the difference between symbol errors and bit errors in other situations.

```
a = [1 2 3]'; b = [1 4 4]';
format rat % Display fractions instead of decimals.
[snum, srate] = symerr(a, b)
```

```
snum =
```

```
2
```

```
srate =
```

```
2/3
```

```
[bnum, brate] = biterr(a, b)
```

```
bnum =
```

```
5
```

```
brate =
```

```
5/9
```

bnum is five because the second entries differ in two bits and the third entries differ in three bits. brate is 5/9 since the total number of bits is nine. The total number of bits is, by definition, the number of entries in a or b times the maximum number of bits among all entries of a and b.

### Eye Diagrams

An eye diagram is a simple and convenient tool for studying the effects of intersymbol interference and other channel impairments in digital transmission. To construct an eye diagram, plot the received signal against time on a fixed-interval axis. At the end of the fixed time interval, wrap around to the beginning of the time axis. Thus the diagram consists of many overlapping curves. One way to use an eye diagram is to look for the place

where the “eye” is most widely opened, and use that point as the decision point when demapping a demodulated signal to recover a digital message.

To produce an eye diagram from a signal, use the `eyediagram` function. The signal can have different formats, as the table below indicates.

**Table 2-1: Representing In-Phase and Quadrature Components of Signal**

Signal Format	Source of In-Phase Components	Source of Quadrature Components
Real matrix with two columns	First column	Second column
Complex vector	Real part	Imaginary part
Real vector	Vector contents	Quadrature component is always zero

### Example: Eye Diagrams

The code below illustrates the use of the eye diagram for finding the best decision point. It maps a random digital signal to a 16-QASK waveform, then uses a raised cosine filter to simulate a noisy transmission channel. Several commands manipulate the filtered data to isolate its steady-state behavior. Then the `eyediagram` command produces an eye diagram from the resulting signal.

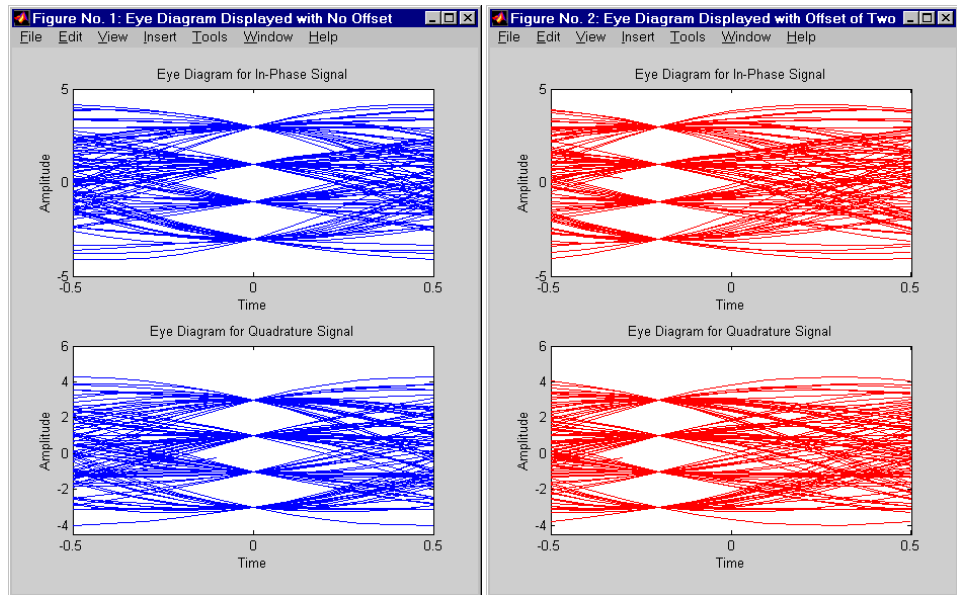
```
% Define the M-ary number and sampling rates.
M = 16; Fd = 1; Fs = 10;
Pd = 100; % Number of points in the calculation
msg_d = randint(Pd, 1, M); % Random integers in the range [0, M-1]
% Modulate using square constellation QASK method.
msg_a = modmap(msg_d, Fd, Fd, 'qask', M);
% Assume the channel is equivalent to a raised cosine filter.
delay = 3; % Delay of the raised cosine filter
rcv = rcosflt(msg_a, Fd, Fs, 'fir/normal', .5, delay);

% Truncate the output of rcosflt to remove response tails.
propdelay = delay .* Fs/Fd + 1; % Propagation delay of filter
```

```
rcv1 = rcv(propdelay:end-(propdelay-1),:); % Truncated version on
N = Fs/Fd;
```

```
% Plot the eye diagram of the resulting signal sampled and
% displayed with no offset.
offset1 = 0;
h1 = eyediagram(rcv1, N, 1/Fd, offset1);
set(h1, 'Name', 'Eye Diagram Displayed with No Offset');
```

Notice that a vertical line down the center of the diagram would cross the “eye” at its most widely opened point, as in the left-hand side below.



In the right-hand diagram above, a similar vertical line would *not* cross the eye at the most widely opened point. This diagram results from the commands

```
offset2 = 2;
h2 = eyediagram(rcv1, N, 1/Fd, offset2, 'r-');
set(h2, 'Name', 'Eye Diagram Displayed with Offset of Two');
```

This example continues by using the information gathered from the eye diagrams to choose the decision-timing offset in the demodmap command.

(Notice that the actual offset value in `demodmap` is `offset1+1` because `eyediagram` and `demodmap` express offsets in a different way.)

```
% Continue, using the offset information for digital demapping.
newmsg1 = demodmap(rcv1, [Fd offset1+1], Fs, 'qask', 16);
s1 = symerr(msg_d, newmsg1) % Number of symbol errors

s1 =

    0
```

By contrast, an offset value based on `offset2` leads to errors in the recovered digital signal. Your exact number of errors might vary because the message `msg_d` consists of random numbers.

```
newmsg2 = demodmap(rcv1, [Fd offset2+1], Fs, 'qask', 16);
s2 = symerr(msg_d, newmsg2)

s2 =

    8
```

As an additional example of using the `eyediagram` function, the commands below display the eye diagram with no offset, but based on data that is sampled with an offset of two samples. This sampling offset simulates errors in timing that result from being two samples away from perfect synchronization.

```
h3 = eyediagram(rcv1(1+offset2:end,:), N, 1/Fd, 0);
set(h3, 'Name', 'Eye Diagram Sampled with Offset of Two');
```

## Scatter Plots

A scatter plot of a signal shows the signal's value at a given decision point. In the best case, the decision point should be at the time when the eye of the signal's eye diagram is the most widely open.

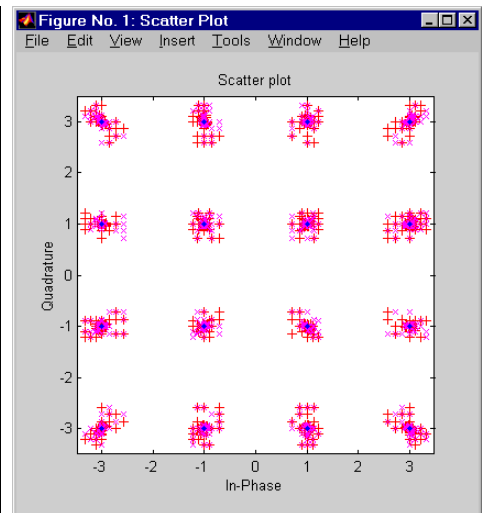
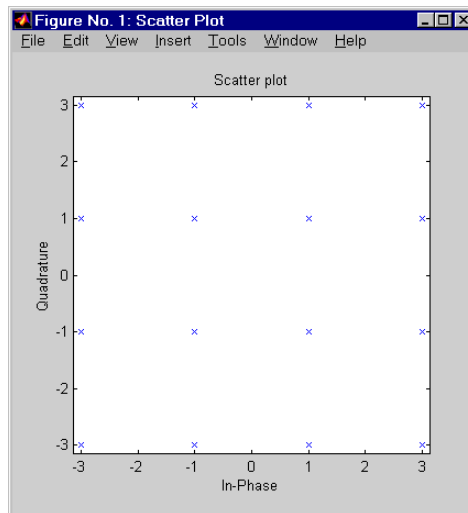
To produce a scatter plot from a signal, use the `scatterplot` function. The signal can have different formats, as in the case of the `eyediagram` function. See Table 2-1, Representing In-Phase and Quadrature Components of Signal, on page 2-9 for details.

## Example: Scatter Plots

The code below is similar to the example from the section, “Example: Eye Diagrams” on page 2-9. It produces a scatter plot from the received analog signal, instead of an eye diagram.

```
% Define the M-ary number and sampling rates.
M = 16; Fd = 1; Fs = 10;
Pd = 200; % Number of points in the calculation
msg_d = randint(Pd, 1, M); % Random integers in the range [0, M-1]
% Modulate using square constellation QASK method.
msg_a = modmap(msg_d, Fd, Fs, 'qask', M);
% Assume the channel is equivalent to a raised cosine filter.
rcv = rcosflt(msg_a, Fd, Fs);
% Create the scatter plot of the received signal,
% ignoring the first three and the last four symbols.
N = Fs/Fd;
rcv_a = rcv(3*N+1:end-4*N, :);
h = scatterplot(rcv_a, N, 0, 'bx');
```

Varying the third parameter in the scatterplot command changes the offset. An offset of zero yields optimal results, shown on the left below.



The diagram on the right results from the commands below. The x's and +'s reflect two offsets that are not optimal because they are too late and too early, respectively. Notice that in the diagram, the dots are the actual constellation points, while the other symbols are perturbations of those points.

```
hold on;  
scatterplot(rcv_a, N, N+1, 'r+', h); % Plot +'s  
scatterplot(rcv_a, N, N-1, 'mx', h); % Plot x's  
scatterplot(rcv_a, N, 0, 'b.', h); % Plot dots
```

## Source Coding

Source coding, also known as *quantization* or *signal formatting*, is a way of processing data in order to reduce redundancy or prepare it for later processing. Analog-to-digital conversion and data compression are two categories of source coding.

Source coding divides into two basic procedures: *source encoding* and *source decoding*. Source encoding converts a source signal into a digital signal using a quantization method. The symbols in the resulting signal are nonnegative integers in some finite range. Source decoding recovers the original information from the source coded signal.

### Source Coding Features of the Toolbox

This toolbox supports two source coding quantization methods: scalar quantization and predictive quantization. It does not support vector quantization. Functions in the toolbox can accomplish these tasks:

- Quantize a signal according to a partition and codebook that you specify
- Optimize partition and codebook parameters for a set of training data
- Encode or decode a signal using the differential pulse code modulation (DPCM) technique
- Optimize DPCM parameters for a set of training data
- Perform  $\mu$ -law or A-law compressor or expander calculations

### Representing Quantization Parameters

Scalar quantization is a process that maps all inputs within a specified range to a common value. It maps inputs in a different range of values to a different common value. In effect, scalar quantization digitizes an analog signal. Two parameters determine a quantization: a partition and a codebook. This section describes how toolbox functions represent these parameters.

#### Partitions

A quantization partition defines several contiguous, nonoverlapping ranges of values within the set of real numbers. To specify a partition in MATLAB, list the distinct endpoints of the different ranges in a vector.



For example, if the partition separates the real number line into the four sets:

1  $\{x: x \leq 0\}$

2  $\{x: 0 < x \leq 1\}$

3  $\{x: 1 < x \leq 3\}$  and

4  $\{x: 3 < x\}$

then you can represent the partition as the three-element vector

```
partition = [0, 1, 3];
```

Notice that the length of the partition vector is one less than the number of partition intervals.

### Codebooks

A codebook tells the quantizer which common value to assign to inputs that fall into each range of the partition. Represent a codebook as a vector whose length is the same as the number of partition intervals. For example, the vector

```
codebook = [-1, 0.5, 2, 3];
```

is one possible codebook for the partition  $[0, 1, 3]$ .

## Quantizing a Signal

The previous section described how you can represent the partition and codebook that determine your scalar quantization process. This section shows how to use these parameters in the `quantiz` function.

### Scalar Quantization Example 1

The code below shows how the `quantiz` function uses `partition` and `codebook` to map a real vector, `sample`, to a new vector, `quantized`, whose entries are either -1, 0.5, 2, or 3.

```
partition = [0, 1, 3];
codebook = [-1, 0.5, 2, 3];
sample = [-2.4, -1, -.2, 0, .2, 1, 1.2, 1.9, 2, 2.9, 3, 3.5, 5];
[index, quantized] = quantiz(sample, partition, codebook);
quantized
```

quantized =

Columns 1 through 7

-1.0000 -1.0000 -1.0000 -1.0000 0.5000 0.5000 2.0000

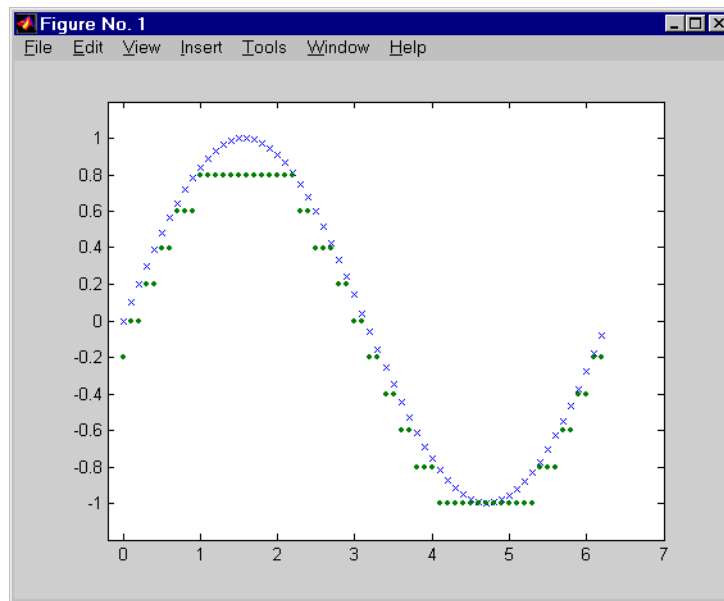
Columns 8 through 13

2.0000 2.0000 2.0000 2.0000 3.0000 3.0000

### Scalar Quantization Example 2

This example illustrates the nature of scalar quantization more clearly. After quantizing a sampled sine wave, it plots the original and quantized signals. The plot contrasts the x's that make up the sine curve with the dots that make up the quantized signal. The vertical coordinate of each dot is a value in the vector codebook.

```
t = [0:.1:2*pi]; % Times at which to sample the sine function
sig = sin(t); % Original signal, a sine wave
partition = [-1:.2:1]; % Length 11, to represent 12 intervals
codebook = [-1.2:.2:1]; % Length 12, one entry for each interval
[index, quants] = quantiz(sig, partition, codebook); % Quantize.
plot(t, sig, 'x', t, quants, '.')
axis([-1.2 7 -1.2 1.2])
```



### Determining Which Interval Each Input Is in

The `quantiz` function also returns a vector that tells which interval each input is in. For example, the output below says that the input entries lie within the intervals labeled 0, 6, and 5, respectively. Here, the 0th interval consists of real numbers less than or equal to 3; the 6th interval consists of real numbers greater than 8 but less than or equal to 9; and the 5th interval consists of real numbers greater than 7 but less than or equal to 8.

```
partition = [3, 4, 5, 6, 7, 8, 9];
index = quantiz([2 9 8], partition)
index =

0
6
5
```

If you continue this example by defining a codebook vector such as

```
codebook = [3, 3, 4, 5, 6, 7, 8, 9];
```

then the equation below relates the vector `index` to the quantized signal `quants`.

```
quants = codebook(index+1);
```

This formula for `quants` is exactly what the `quantiz` function uses if you instead phrase the example more concisely as below.

```
partition = [3, 4, 5, 6, 7, 8, 9];
codebook = [3, 3, 4, 5, 6, 7, 8, 9];
[index, quants] = quantiz([2 9 8], partition, codebook);
```

## Optimizing Quantization Parameters

Quantization distorts a signal. You can lessen the distortion by choosing appropriate partition and codebook parameters. However, testing and selecting parameters for large signal sets with a fine quantization scheme can be tedious. One way to produce partition and codebook parameters easily is to optimize them according to a set of so-called training data.

---

**Note** The training data that you use should be typical of the kinds of signals that you will actually be quantizing.

---

### Example: Optimizing Scalar Quantization Parameters

The `lloyd` function optimizes the partition and codebook according to the Lloyd algorithm. The code below optimizes the partition and codebook for one period of a sinusoidal signal, starting from a rough initial guess. Then it uses these parameters to quantize the original signal using the initial guess parameters as well as the optimized parameters. The output shows that the mean square distortion after quantizing is much less for the optimized parameters. Notice that the `quantiz` function automatically computes the mean square distortion and returns it as the third output parameter.

```
% Start with the setup from 2nd example in "Quantizing a Signal."
t = [0: .1: 2*pi];
sig = sin(t);
partition = [-1: .2: 1];
codebook = [-1.2: .2: 1];
% Now optimize, using codebook as an initial guess.
```

```

[partition2, codebook2] = lloyds(sig, codebook);
[index, quants, distort] = quantiz(sig, partition, codebook);
[index2, quant2, distort2] = quantiz(sig, partition2, codebook2);
% Compare mean square distortions from initial and optimized
[distort, distort2] % parameters.

```

```
ans =
```

```
0.0148    0.0024
```

## Implementing Differential Pulse Code Modulation

The quantization in the section “Quantizing a Signal” on page 2-15 requires no *a priori* knowledge about the transmitted signal. In practice, you can often make educated guesses about the present signal based on past signal transmissions. Using such educated guesses to help quantize a signal is known as *predictive quantization*. The most common predictive quantization method is differential pulse code modulation (DPCM).

The functions `dpcmenco`, `dpcmdeco`, and `dpcmopt` can help you implement a DPCM predictive quantizer with a linear predictor.

### DPCM Terminology

To determine an encoder for such a quantizer, you must supply not only a partition and codebook as described in “Representing Quantization Parameters” on page 2-14, but also a *predictor*. The predictor is a function that the DPCM encoder uses to produce the educated guess at each step. A linear predictor has the form

$$y(k) = p(1)x(k-1) + p(2)x(k-2) + \dots + p(m-1)x(k-m+1) + p(m)x(k-m)$$

where  $x$  is the original signal,  $y(k)$  attempts to predict the value of  $x(k)$ , and  $p$  is an  $m$ -tuple of real numbers. Instead of quantizing  $x$  itself, the DPCM encoder quantizes the *predictive error*,  $x - y$ . The integer  $m$  above is called the *predictive order*. The special case when  $m = 1$  is called *delta modulation*.

### Representing Predictors

If the guess for the  $k$ th value of the signal  $x$ , based on earlier values of  $x$ , is

$$y(k) = p(1)x(k-1) + p(2)x(k-2) + \dots + p(m-1)x(k-m+1) + p(m)x(k-m)$$

then the corresponding predictor vector for toolbox functions is

```
predictor = [0, p(1), p(2), p(3), ..., p(m-1), p(m)]
```

---

**Note** The initial zero in the predictor vector makes sense if you view the vector as the polynomial transfer function of a finite impulse response (FIR) filter.

---

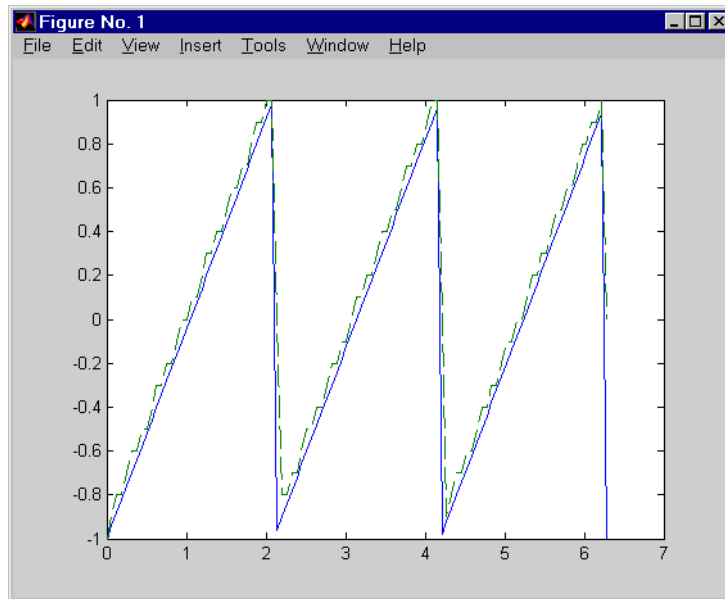
### Example: DPCM Encoding and Decoding

A simple special case of DPCM quantizes the difference between the signal's current value and its value at the previous step. Thus the predictor is just  $y(k) = x(k-1)$ . The code below implements this scheme. It encodes a sawtooth signal, decodes it, and plots both the original and decoded signals. The solid line is the original signal, while the dashed line is the recovered signals. The example also computes the mean square error between the original and decoded signals.

```
predictor = [0 1]; % y(k)=x(k-1)
partition = [-1:.1:.9];
codebook = [-1:.1:1];
t = [0:pi/50:2*pi];
x = sawtooth(3*t); % Original signal
% Quantize x using DPCM
encodedx = dpcmenco(x, codebook, partition, predictor);
% Try to recover x from the modulated signal.
decodedx = dpcmdeco(encodedx, codebook, predictor);
plot(t, x, t, decodedx, '- -')
distor = sum((x-decodedx).^2)/length(x) % Mean square error

distor =

    0.0327
```



## Optimizing DPCM Parameters

The section “Optimizing Quantization Parameters” on page 2-18 describes how you can use training data with the `lloydys` function to help find quantization parameters that will minimize signal distortion. This section describes similar procedures for using the `dpcmopt` function in conjunction with the two functions `dpcmenco` and `dpcmdeco`, which first appear in the previous section.

---

**Note** The training data that you use with `dpcmopt` should be typical of the kinds of signals that you will actually be quantizing with `dpcmenco`.

---

### Example: Comparing Optimized and Nonoptimized DPCM Parameters

This example is similar to the one in the last section. However, whereas the last example created `predictor`, `partition`, and `codebook` in a straightforward but haphazard way, this example uses the same codebook (now called `initcodebook`) as an initial guess for a new *optimized* codebook parameter. This example also uses the predictive order, 1, as the desired order of the new

optimized predictor. The `dpcmopt` function creates these optimized parameters, using the sawtooth signal `x` as training data. The example goes on to quantize the training data itself; in theory, the optimized parameters are suitable for quantizing other data that is similar to `x`. Notice that the mean square distortion here is much less than the distortion in the previous example.

```
t = [0: pi /50: 2*pi ];
x = sawtooth(3*t); % Original signal
initcodebook = [-1:.1:1]; % Initial guess at codebook
% Optimize parameters, using initial codebook and order 1.
[predictor, codebook, partition] = dpcmopt(x, 1, initcodebook);
% Quantize x using DPCM.
encodedx = dpcmenco(x, codebook, partition, predictor);
% Try to recover x from the modulated signal.
decodedx = dpcmdeco(encodedx, codebook, predictor);
distor = sum((x-decodedx).^2)/length(x) % Mean square error

distor =

    0.0063
```

## Companding a Signal

In certain applications, such as speech processing, it is common to use a logarithm computation, called a *compressor*, before quantizing. The inverse operation of a compressor is called an *expander*. The combination of a compressor and expander is called a *comparer*.

The `compand` function supports two kinds of companders:  $\mu$ -law and A-law companders. Its reference page lists both compressor laws.

### Example: A $\mu$ -Law Comparer

The code below quantizes an exponential signal in two ways and compares the resulting mean square distortions. First, it simply uses the `quantiz` function with a partition consisting of length-one intervals. In the second trial, `compand` implements a  $\mu$ -law compressor, `quantiz` quantizes the compressed data, and finally `compand` expands the quantized data. The output shows that the distortion is smaller for the second scheme. This is because equal-length intervals are well-suited to the logarithm of `sig`, but not well-suited to `sig` itself.



```

mu = 255; % Parameter for mu-law compander
sig = -4: .1: 4;
sig = exp(sig); % Exponential signal to quantize
V = max(sig);
% 1. Quantize using equal-length intervals and no compander.
[index, quants, distort] = quantiz(sig, 0: floor(V), 0: ceil(V));

% 2. Use same partition and codebook, but compress
% before quantizing and expand afterwards.
compsig = compand(sig, mu, V, 'mu/compressor');
[index, quants] = quantiz(compsig, 0: floor(V), 0: ceil(V));
newsig = compand(quants, mu, max(quants), 'mu/expander');
distort2 = sum((newsig - sig).^2) / length(sig);
[distort, distort2] % Display both mean square distortions.

ans =

    0.5348    0.0397

```

## Selected Bibliography for Source Coding

- [1] Kondo, A. M. *Digital Speech*. Chichester, England: John Wiley & Sons, 1994.
- [2] Sklar, Bernard. *Digital Communications: Fundamentals and Applications*. Englewood Cliffs, N.J.: Prentice-Hall, 1988.

## Block Coding

Error-control coding techniques detect and possibly correct errors that occur when messages are transmitted in a digital communication system. To accomplish this, the encoder transmits not only the information symbols but also extra redundant symbols. The decoder interprets what it receives, using the redundant symbols to detect and possibly correct whatever errors occurred during transmission. You might use error-control coding if your transmission channel is very noisy or if your data is very sensitive to noise. Depending on the nature of the data or noise, you might choose a specific type of error-control coding.

Block coding is a special case of error-control coding. Block coding techniques maps a fixed number of message symbols to a fixed number of code symbols. A block coder treats each block of data independently and is a memoryless device.

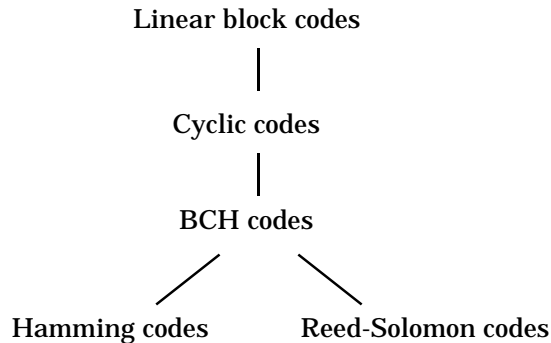
This section discusses these topics:

- “Block Coding Features of the Toolbox” on page 2-25
- “Block Coding Terminology” on page 2-26
- “Representing Messages and Codewords” on page 2-26
- “Representing Block Coding Parameters” on page 2-30
- “Creating and Decoding Block Codes” on page 2-35
- “Performing Other Block Code Tasks” on page 2-40

For background information about block coding, see the works listed in “Selected Bibliography for Block Coding” on page 2-42.

## Block Coding Features of the Toolbox

The class of linear block coding techniques includes categories shown below.



The Communications Toolbox supports general linear block codes. It also includes functions to process cyclic, BCH, Hamming, and Reed-Solomon codes (which are all special kinds of linear block codes). Functions in the toolbox can accomplish these tasks:

- Encode or decode a message using one of the techniques mentioned above
- Determine characteristics of a technique, such as error-correction capability or valid message length
- Perform lower-level computations associated with a technique, such as:
  - Compute a decoding table
  - Compute a generator or parity-check matrix
  - Convert between generator and parity-check matrices
  - Compute a generator polynomial

---

**Note** The functions in this toolbox are designed for block codes that use an alphabet having 2 or  $2^m$  symbols.

---

The table below lists the functions that are related to each supported block coding technique.

Table 2-2: Functions Related to Block Coding Techniques

Block Coding Technique	Toolbox Functions
Linear block	encode, decode, gen2par, syndtabl e
Cyclic	encode, decode, cycl pol y, cycl gen, gen2par, syndtabl e
BCH	encode, decode, bchenco, bchdeco, bchpol y, cycl gen, gen2par, syndtabl e
Hamming	encode, decode, hammgen, gen2par, syndtabl e
Reed-Solomon	encode, decode, rsenco, rsdeco, rsencode, rsdecode, rspol y, rsencof, rsdecof, syndtabl e

Block Coding Terminology

Throughout this section, the information to be encoded consists of a sequence of *message* symbols and the code that is produced consists of a sequence of *codewords*.

Each block of *k* message symbols is encoded into a codeword that consists of *n* symbols; in this context, *k* is called the message length, *n* is called the codeword length, and the code is called an [*n,k*] code.

Representing Messages and Codewords

Each message or codeword is an ordered grouping of symbols. The next few subsections illustrate the various ways that these symbols may be organized or interpreted as input and output.

Binary Vector Format

One straightforward MATLAB format for messages and codewords is a vector of 0s and 1s. That is, messages and codes might look like msg and code in the lines below.

```

n = 6; k = 4; % Set codeword length and message length
% for a [6, 4] code.
msg = [1 0 0 1 1 0 1 0 1 0 1 1]'; % Message is a binary column.
code = encode(msg, n, k, 'cyclic'); % Code will be a binary column.
msg'

ans =

    1     0     0     1     1     0     1     0     1     0     1     1

code'

ans =

Columns 1 through 12

    0     0     1     0     0     1     1     0     1     0     1     0

Columns 13 through 18

    0     1     1     0     1     1

```

In this example, `msg` consists of 12 entries, which are interpreted as three four-digit (since  $k = 4$ ) messages. The resulting vector `code` comprises three six-digit (since  $n = 6$ ) codewords, which are concatenated to form a vector of length eighteen.

### Binary Matrix Format

You can also organize coding information so as to emphasize the grouping of digits in a single message or codeword. The code below illustrates this by listing each four-digit message on a separate row in `msg` and each six-digit codeword on a separate row in `code`.

```

n = 6; k = 4; % Set codeword length and message length.
msg = [1 0 0 1; 1 0 1 0; 1 0 1 1]; % Message is a binary matrix.
code = encode(msg, n, k, 'cyclic'); % Code will be a binary matrix.
msg

```

```

msg =

      1      0      0      1
      1      0      1      0
      1      0      1      1

code

code =

      0      0      1      0      0      1
      1      0      1      0      1      0
      0      1      1      0      1      1

```

For all coding techniques *except* Reed-Solomon, the message matrix must have  $k$  columns. The corresponding code matrix has  $n$  columns.

**Reed-Solomon Coding Using Binary Matrix Format.** For Reed-Solomon codes, the message matrix must have  $m$  columns, where  $m$  is an integer greater than or equal to 3 that satisfies  $n = 2^m - 1$ .

### Decimal Format

Another way to process the same information is to regard each of the three rows of `msg` and `code` above as binary representations of decimal integers. MATLAB then accepts the corresponding decimal integers as valid messages, and returns decimal integers as codewords.

---

**Note** If  $2^n$  or  $2^k$  is large, then you should use the default binary format instead of the decimal format. This is because the function uses a binary format internally, while the round-off error associated with converting many bits to large decimal numbers and back might be substantial.

---



---

**Note** In this context, MATLAB expects the *leftmost* bit to be the least significant bit.

---

The syntax for the encode command must mention the decimal format explicitly, as in the example below. Notice that `/decimal` is appended to the fourth argument in the encode command.

```
n = 6; k = 4; % Set codeword length and message length.
msg = [9;5;13]; % Message is a decimal column vector.
% Code will be a decimal vector.
code = encode(msg, n, k, 'cyclic/decimal')

code =

    36
    21
    54
```

---

**Note** The three examples above used cyclic coding. The formats for messages and codes are similar for Hamming, generic linear, and BCH codes.

---

**Reed-Solomon Coding Using Decimal Format.** For Reed-Solomon coding using decimal formats, the message matrix must have  $k$  columns. Each entry in the matrix must be an integer between 0 and  $n$ . The example below illustrates the decimal format for Reed-Solomon coding using the encode command.

```
m = 3;
n = 2^m-1; k = 4; % Set codeword length and message length.
msgdec = [1 6 4 1; 0 0 4 3]; % Message is a decimal matrix.
% Code will be a decimal vector.
codedec = encode(msgdec, n, k, 'rs/decimal')

codedec =
```

```
    0    4    3    1    6    4    1
    3    7    5    0    0    4    3
```

### Exponential Format (Reed-Solomon Code Only)

For Reed-Solomon coding using exponential formats, the message matrix must have  $k$  columns. Each entry of the matrix must be an integer between -1 and

n-1. The example below is the exponential-form counterpart of the Reed-Solomon example from the previous section.

```
m = 3;
n = 2^m-1; k = 4; % Set codeword length and message length.
msg = [0 5 3 0; -1 -1 3 2];
% Message is an exponential-form matrix.
% Code will be an exponential-form matrix.
code = encode(msg, n, k, 'rs/power');
```

The name “exponential format” comes from one of MATLAB’s standard formats for elements of  $GF(2^m)$ . This format uses integers from -1 to  $2^m-2$ , where the symbol -Inf is sometimes substituted for -1. See “Exponential Format” on page 2-90 for definitions.

Representing Block Coding Parameters

This subsection describes the items that you might need in order to process  $[n,k]$  linear block codes. The table below lists the items and the coding techniques for which they are most relevant.

Table 2-3: Parameters Used in Block Coding Techniques

Parameter	Block Coding Technique
Generator Matrix	Generic linear block
Parity-Check Matrix	Generic linear block
Generator Polynomial	Cyclic, BCH, Reed-Solomon
Primitive Polynomial and List of Galois Field Elements	Hamming, Reed-Solomon
Decoding Table	Generic linear block, Hamming

Generator Matrix

The process of encoding a message into an  $[n,k]$  linear block code is determined by a  $k$ -by- $n$  generator matrix  $G$ . Specifically, the 1-by- $k$  message vector  $v$  is encoded into the 1-by- $n$  codeword vector  $vG$ . If  $G$  has the form  $[I_k \ P]$  or  $[P \ I_k]$ , where  $P$  is some  $k$ -by- $(n-k)$  matrix and  $I_k$  is the  $k$ -by- $k$  identity matrix, then  $G$  is said to be in *standard form*. (Some authors, e.g., Clark and Cain [1], use the



first standard form, while others, e.g., Lin and Costello [2], use the second.) Most functions in this toolbox assume that a generator matrix is in standard form when you use it as an input argument.

Some examples of generator matrices are in the next section, “Parity-Check Matrix.”

### Parity-Check Matrix

Decoding an  $[n,k]$  linear block code requires an  $(n-k)$ -by- $n$  parity-check matrix  $H$ . It satisfies  $GH^{\text{tr}} = 0 \pmod{2}$ , where  $H^{\text{tr}}$  denotes the matrix transpose of  $H$ ,  $G$  is the code's generator matrix, and this zero matrix is  $k$ -by- $(n-k)$ . If  $G = [I_k \ P]$  then  $H = [-P^{\text{tr}} \ I_{n-k}]$ . Most functions in this toolbox assume that a parity-check matrix is in standard form when you use it as an input argument.

The table below summarizes the standard forms of the generator and parity-check matrices for an  $[n,k]$  binary linear block code.

Type of Matrix	Standard Form	Dimensions
Generator	$[I_k \ P]$ or $[P \ I_k]$	$k$ -by- $n$
Parity-check	$[-P^{\text{tr}} \ I_{n-k}]$ or $[I_{n-k} \ -P^{\text{tr}}]$	$(n-k)$ -by- $n$

$I_k$  is the identity matrix of size  $k$  and the ' symbol indicates matrix transpose. (For *binary* codes, the minus signs in the parity-check form listed above are irrelevant; that is,  $-1 = 1$  in the binary field.)

**Examples.** In the command below, `parmat` is a parity-check matrix and `genmat` is a generator matrix for a Hamming code in which  $[n,k] = [2^3-1, n-3] = [7,4]$ . Notice that `genmat` has the standard form  $[P \ I_k]$ .

```
[parmat, genmat] = hamngen(3)
```

```
parmat =
```

```

1     0     0     1     0     1     1
0     1     0     1     1     1     0
0     0     1     0     1     1     1
```

genmat =

1	1	0	1	0	0	0
0	1	1	0	1	0	0
1	1	1	0	0	1	0
1	0	1	0	0	0	1

The next example finds parity-check and generator matrices for a [7,3] cyclic code. The `cyclpoly` function is mentioned below in “Generator Polynomial.”

```
genpoly = cyclpoly(7, 3);
[parmat, genmat] = cyclgen(7, genpoly)
```

parmat =

1	0	0	0	1	1	0
0	1	0	0	0	1	1
0	0	1	0	1	1	1
0	0	0	1	1	0	1

genmat =

1	0	1	1	1	0	0
1	1	1	0	0	1	0
0	1	1	1	0	0	1

The example below converts a generator matrix for a [5,3] linear block code into the corresponding parity-check matrix.

```
genmat = [1 0 0 1 0; 0 1 0 1 1; 0 0 1 0 1];
parmat = gen2par(genmat)
```

parmat =

1	1	0	1	0
0	1	1	0	1

The same function `gen2par` can also convert a parity-check matrix into a generator matrix.

## Generator Polynomial

Cyclic codes, including the special cases of BCH and Reed-Solomon codes, have special algebraic properties that allow a polynomial to determine the coding process completely. This so-called generator polynomial is a degree- $(n-k)$  divisor of the polynomial  $x^n-1$ . Van Lint [4] explains how a generator polynomial determines a cyclic code.

The functions in this toolbox that produce generator polynomials are `bchpoly`, `cyclpoly`, and `rspoly`. They represent a generator polynomial using a row vector that lists the polynomial's coefficients in order of *ascending* powers of the variable. Functions dealing with BCH and generic cyclic codes use binary digits as coefficients, as in the first example below. Functions dealing with Reed-Solomon codes express the coefficients (which are elements of  $\text{GF}(2^m)$ ) in exponential format, as in the second example below. See “Representing Elements of Galois Fields” on page 2-90 for a description of this exponential format for elements of Galois fields.

**Examples.** The command

```
genpoly = cyclpoly(7, 3)
```

```
genpoly =
```

```
1      0      1      1      1
```

finds that one valid generator polynomial for a [7,3] cyclic code is  $1 + x^2 + x^3 + x^4$ .

A second example finds that a generator polynomial for a [15,13] Reed-Solomon code is  $\alpha^3 + \alpha^5x + \alpha^0x^2$ , where  $\alpha$  is a root of MATLAB's default primitive polynomial for  $\text{GF}(15+1)$ .

```
r = rspoly(15, 13)
```

```
r =
```

```
3      5      0
```

## Primitive Polynomial and List of Galois Field Elements

Hamming and Reed-Solomon codes rely on algebraic fields that have  $2^m$  elements (or, more generally,  $p^m$  elements for a prime number  $p$ ). Elements of

such fields are named *relative to* a distinguished element of the field that is called a primitive element. Some functions in this toolbox use a primitive polynomial or a list of elements in the field as a way to determine the primitive element and, consequently, as a way to name elements of the field. See “Galois Field Computations” on page 2-89 and especially the subsection “Representing Elements of Galois Fields” for details about MATLAB’s use of primitive polynomials and lists of Galois field elements.

To reduce the mathematical background that you need to use the block coding functions, simply use the default parameters in commands that ask for primitive polynomials or lists of Galois field elements. For more specifics, see the reference pages for `encode`, `decode`, `hammgen`, `rsenco`, `rsencode`, `rsdeco`, `rsdecode`, and `rspoly`.

### Decoding Table

A decoding table tells a decoder how to correct errors that may have corrupted the code during transmission. Hamming codes can correct any single-symbol error in any codeword. Other codes can correct, or partially correct, errors that corrupt more than one symbol in a given codeword.

This toolbox represents a decoding table as a matrix with  $n$  columns and  $2^{n-k}$  rows. Each row gives a correction vector for one received codeword vector. A Hamming decoding table has  $n+1$  rows. The `syndtable` function generates a decoding table for a given parity-check matrix.

### Example: Using a Decoding Table

The script below shows how to use a Hamming decoding table to correct an error in a received message. The `hammgen` function produces the parity-check matrix, while the `syndtable` function produces the decoding table. The transpose of the parity-check matrix is multiplied on the left by the received codeword, yielding the *syndrome*. The decoding table helps determine the correction vector. The corrected codeword is the sum (modulo 2) of the correction vector and the received codeword.

```
% Use a [7, 4] Hamming code.
m = 3; n = 2^m - 1; k = n - m;
parmat = hammgen(m); % Produce parity-check matrix.
trt = syndtable(parmat); % Produce decoding table.
recd = [1 0 0 1 1 1 1] % Suppose this is the received vector.
syndrome = rem(recd * parmat', 2);
```

```

syndrome_de = bi2de(syndrome, 'left-msb'); % Convert to decimal.
disp([' Syndrome = ', num2str(syndrome_de), ...
      ' (decimal), ', num2str(syndrome), ' (binary)'])
corrvect = trt(1+syndrome_de,:) % Correction vector
% Now compute the corrected codeword.
correctedcode = rem(corrvect+recd, 2)

```

The output is below.

```

recd =

     1     0     0     1     1     1     1

Syndrome = 3 (decimal), 0 1 1 (binary)

corrvect =

     0     0     0     0     1     0     0

correctedcode =

     1     0     0     1     0     1     1

```

## Creating and Decoding Block Codes

The functions for encoding and decoding linear block codes are `encode`, `decode`, `bchenco`, `bchdeco`, `rsenco`, `rsdeco`, `rsencode`, `rsdecode`, `rsencof`, and `rsdecof`. The first two in this list are general-purpose functions that invoke other functions from the list when appropriate. This section discusses how to use these functions to create and decode generic linear block codes, cyclic codes, BCH codes, Hamming codes, and Reed-Solomon codes.

### Generic Linear Block Codes

Encoding a message using a generic linear block code requires a generator matrix. If you have defined variables `msg`, `n`, `k`, and `genmat`, then either of the commands

```

code = encode(msg, n, k, 'linear', genmat);
code = encode(msg, n, k, 'linear/decimal', genmat);

```

encodes the information in `msg` using the `[n,k]` code that the generator matrix `genmat` determines. The `/decimal` option, suitable when  $2^n$  and  $2^k$  are not very large, indicates that `msg` contains nonnegative decimal integers rather than their binary representations. See “Representing Messages and Codewords” on page 2-26 or the reference page for `encode` for a description of the formats of `msg` and `code`.

Decoding the code requires the generator matrix and possibly a decoding table. If you have defined variables `code`, `n`, `k`, `genmat`, and possibly also `trt`, then the commands

```
newmsg = decode(code, n, k, 'linear', genmat);
newmsg = decode(code, n, k, 'linear/decimal', genmat);
newmsg = decode(code, n, k, 'linear', genmat, trt);
newmsg = decode(code, n, k, 'linear/decimal', genmat, trt);
```

decode the information in `code`, using the `[n,k]` code that the generator matrix `genmat` determines. `decode` also corrects errors according to instructions in the decoding table that `trt` represents.

**Example: Generic Linear Block Coding.** The example below encodes a message, artificially adds some noise, decodes the noisy code, and keeps track of errors that the decoder detects along the way. Since the decoding table contains only zeros, the decoder does not correct any errors.

```
n = 4; k = 2;
genmat = [[1 1; 1 0], eye(2)]; % Generator matrix
msg = [0 1; 0 0; 1 0]; % Three messages, two bits each
% Create three codewords, four bits each.
code = encode(msg, n, k, 'linear', genmat);
noisycode = rem(code + randerr(3, 4, [0 1; .7 .3]), 2); % Add noise.
trt = zeros(2^(n-k), n); % No correction of errors
% Decode, keeping track of all detected errors.
[newmsg, err] = decode(noisycode, n, k, 'linear', genmat, trt);
err_words = find(err~=0) % Find out which words had errors.
```

The output indicates that errors occurred in the first and second words. Your results might vary since this example uses random numbers as errors.

```
err_words =
```

```
1
2
```

## Cyclic Codes

Encoding a message using a cyclic code requires a generator polynomial. If you have defined variables `msg`, `n`, `k`, and `genpoly`, then either of the commands

```
code = encode(msg, n, k, 'cyclic', genpoly);
code = encode(msg, n, k, 'cyclic/decimal', genpoly);
```

encodes the information in `msg` using the  $[n,k]$  code determined by the generator polynomial `genpoly`. `genpoly` is an optional argument for `encode`. The default generator polynomial is `cyclpoly(n, k)`. The `/decimal` option, suitable when  $2^n$  and  $2^k$  are not very large, indicates that `msg` contains nonnegative decimal integers rather than their binary representations. See “Representing Messages and Codewords” on page 2-26 or the reference page for `encode` for a description of the formats of `msg` and `code`.

Decoding the code requires the generator polynomial and possibly a decoding table. If you have defined variables `code`, `n`, `k`, `genpoly`, and `trt`, then the commands

```
newmsg = decode(code, n, k, 'cyclic', genpoly);
newmsg = decode(code, n, k, 'cyclic/decimal', genpoly);
newmsg = decode(code, n, k, 'cyclic', genpoly, trt);
newmsg = decode(code, n, k, 'cyclic/decimal', genpoly, trt);
```

decode the information in `code`, using the  $[n,k]$  code that the generator matrix `genmat` determines. `decode` also corrects errors according to instructions in the decoding table that `trt` represents. `genpoly` is an optional argument in the first two syntaxes above. The default generator polynomial is `cyclpoly(n, k)`.

There are no lower-level functions that provide alternative means to process cyclic codes.

**Example.** The example in the section “Generic Linear Block Codes” on page 2-35 can be modified so that it uses the cyclic coding technique, instead of the linear block code with the generator matrix `genmat`. Make the changes listed below:

- Replace the second line by  
`genpoly = [1 0 1]; % generator poly is 1 + x^2`
- In the fifth and ninth lines (encode and decode commands), replace `genmat` by `genpoly` and replace `'linear'` by `'cyclic'`.

Another example of encoding and decoding a cyclic code is on the reference page for `encode`.

### BCH Codes

BCH codes are a special case of cyclic codes, though the decoding algorithm for BCH codes is more complicated than that for generic cyclic codes. The discussion in the section “Cyclic Codes” above applies almost exactly to the case of BCH codes. The only differences are that:

- `bch` replaces `cyclic` in the syntax for encode and decode.
- `bchpoly(n, k)` replaces `cyclpoly(n, k)` as the default generator polynomial.
- `n` and `k` must be valid codeword and message lengths for BCH code.

Valid codeword lengths for BCH code are those integers of the form  $2^m - 1$  for some integer  $m$  greater than or equal to 3. Given a valid BCH codeword length, the corresponding valid BCH message lengths are those numbers in the second column of the output of the command below.

```
params = bchpoly(n); % Where n = 2^m-1 for some integer m >= 3
```

For example, the output of the command below shows that a BCH code with codeword length 15 may have message length 5, 7, or 11. No other message lengths are valid for this codeword length.

```
params = bchpoly(15)
```

```
params =
```

15	11	1
15	7	2
15	5	3

The third column of the output above represents the error-correction capability for each pair of codeword length and message length.



**Choice of Functions for BCH Coding.** To process BCH codes, you can use either the encode and decode functions, or the lower-level bchenco and bchdeco functions. The syntax of the lower-level functions is slightly different from that of the higher-level functions. The only difference in functionality is that the higher-level functions prepare the input data (including default values of options that you omit) before invoking the lower-level commands. The reference page for encode contains an example that uses encode and decode. The reference pages for bchenco and bchdeco contain other examples.

### Hamming Codes

The reference pages for encode and decode contain examples of encoding and decoding Hamming codes. Also, the section “Decoding Table” on page 2-34 illustrates error-correction in a Hamming code. There are no lower-level functions that provide alternative means to process Hamming codes.

### Reed-Solomon Codes

Reed-Solomon codes are useful for correcting errors that occur in bursts. The codeword length  $n$  of a Reed-Solomon code must have the form  $2^m - 1$ , where  $m$  is an integer greater than or equal to 3. The error correction capability of a Reed-Solomon code is  $\text{floor}((n - k) / 2)$ . Since  $n$  is an odd number, the coding is more efficient when the message length  $k$  is also odd.

One difference between Reed-Solomon codes and the other codes supported in this toolbox is that Reed-Solomon codes process symbols in  $\text{GF}(2^m)$  instead of  $\text{GF}(2)$ . Each such symbol is specified by  $m$  bits. That is why some parts of the section “Representing Messages and Codewords” on page 2-26 make exceptions for Reed-Solomon codes.

Encoding a message using a Reed-Solomon code requires a generator polynomial. The `rspoly` function finds generator polynomials. For example, the command

```
genpoly = rspoly(15, 12)
```

```
genpoly =
```

```
      6      13      11      0
```

shows that the generator polynomial for a [15,12] Reed-Solomon code is

$$\alpha^6 + \alpha^{13}x + \alpha^{11}x^2 + x^3$$

where  $\alpha$  is a root of MATLAB's default primitive polynomial for GF(16). In this example,  $m = 4$ ,  $n = 2^m - 1 = 15$ , and  $k = 12$ .

**Choice of Functions for Reed-Solomon Coding.** To process Reed-Solomon codes, you can use either the encode and decode functions, or the lower-level `rsenco`, `rsdeco`, `rsencode`, and `rsdecode` functions. The syntax of the lower-level functions is slightly different from that of the higher-level functions. The only difference in functionality is that the higher-level functions prepare the input data (including default values of options that you omit) before invoking the lower-level functions. The reference pages for the lower-level functions contain examples that illustrate their use.

## Performing Other Block Code Tasks

This section describes functions that compute typical parameters associated with block codes and functions that convert information from one format to another. Specific tasks are:

- Finding a generator polynomial
- Finding generator and parity-check matrices
- Converting between parity-check and generator matrices
- Finding the error-correction capability

### Finding a Generator Polynomial

To find a generator polynomial for cyclic, BCH, and Reed-Solomon codes, use the functions `cyclpoly`, `bchpoly`, and `rspoly`, respectively. The commands

```
genpolyCyclic = cyclpoly(7, 4);
genpolyBCH = bchpoly(7, 4);
genpolyRS = rspoly(7, 4);
```

all represent valid ways to find one generator polynomial for a [7,4] code of the respective coding method. The result is suitable for use in other block coding functions, such as `encode`.

For generic cyclic coding, there might be more than one generator polynomial consistent with a given codeword length and message length. The `cyclpoly` command syntax includes ways to retrieve all of them or those that satisfy certain constraints that you specify. For example, the command

```
genpolys = cyclpoly(7, 4, 'all')
```

```
genpolys =
```

```
      1      0      1      1
      1      1      0      1
```

shows that  $1 + x^2 + x^3$  and  $1 + x + x^3$  are two possible generator polynomials for a [7,4] cyclic code.

See the reference pages for `cyclpoly`, `bchpoly`, and `rspoly` for details about other options.

### Finding Generator and Parity-Check Matrices

To find a parity-check and generator matrix for a Hamming code with codeword length  $2^m-1$ , use the `hamngen` function as below. `m` must be at least three.

```
[parmat, genmat] = hamngen(m); % Hammi ng
```

To find a parity-check and generator matrix for a cyclic code, use the `cyclgen` function. You must provide the codeword length and a valid generator polynomial. You can use the `cyclpoly` command to produce one possible generator polynomial after you provide the codeword length and message length. For example,

```
[parmat, genmat] = cyclgen(7, cyclpoly(7, 4)); % Cycl ic
```

To find a parity-check and generator matrix for a BCH code, use the same `cyclgen` function mentioned above. Since the generator polynomial must now be valid for BCH code, the `bchpoly` function replaces `cyclpoly`.

```
[parmat, genmat] = cyclgen(7, bchpoly(7, 4)); % BCH
```

### Converting Between Parity-Check and Generator Matrices

The `gen2par` function converts a generator matrix into a parity-check matrix, and vice-versa. Examples to illustrate this are on the reference page for `gen2par`.

### Finding the Error-Correction Capability

The error-correction capability of BCH codes and Reed-Solomon codes depends on the codeword length and message length. The functions `bchpoly` and `rspoly`

perform such computations. To retrieve the error-correction capability  $t$  of BCH and Reed-Solomon codes, respectively, use the commands below.

```
[temp1, temp2, temp3, temp4, t] = bchpoly(n, k); % BCH
[temp1, t] = rspoly(n, k); % Reed-Solomon
```

For Reed-Solomon codes, the error-correction capability is  $\text{floor}((n-k)/2)$ ; for BCH codes, there is no easy formula.

### Selected Bibliography for Block Coding

- [1] Clark, George C. Jr. and J. Bibb Cain. *Error-Correction Coding for Digital Communications*. New York: Plenum Press, 1981.
- [2] Lin, Shu and Daniel J. Costello, Jr. *Error Control Coding: Fundamentals and Applications*. Englewood Cliffs, N.J.: Prentice-Hall, 1983.
- [3] Peterson, W. Wesley and E. J. Weldon, Jr. *Error-correcting Codes*, 2nd ed. Cambridge, Mass.: MIT Press, 1972.
- [4] van Lint, J. H. *Introduction to Coding Theory*. New York: Springer-Verlag, 1982.

## Convolutional Coding

Convolutional coding is a special case of error-control coding. Unlike a block coder, a convolutional coder is not a memoryless device. Even though a convolutional coder accepts a fixed number of message symbols and produces a fixed number of code symbols, its computations depend not only on the current set of input symbols but on some of the previous input symbols.

This section:

- Outlines the convolutional coding features of the Communications Toolbox
- Defines the two supported ways to describe a convolutional encoder:
  - Polynomial description
  - Trellis description
- Describes how to encode and decode using the `convenc` and `vitdec` functions
- Gives additional examples of convolutional coding

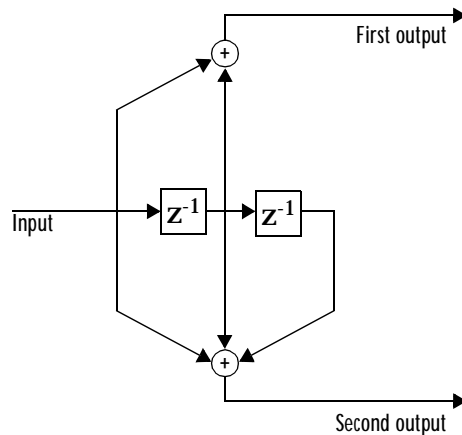
### Convolutional Coding Features of the Toolbox

The Communications Toolbox supports feedforward or feedback convolutional codes that can be described by a trellis structure or a set of generator polynomials. It uses the Viterbi algorithm to implement hard-decision and soft-decision decoding.

For background information about convolutional coding, see the works listed in “Selected Bibliography for Convolutional Coding” on page 2-55.

### Polynomial Description of a Convolutional Encoder

A polynomial description of a convolutional encoder describes the connections among shift registers and modulo-2 adders. For example, the figure below depicts a feedforward convolutional encoder that has one input, two outputs, and two shift registers.



**Figure 2-1: Example of a Convolutional Encoder Diagram with Shift Registers**

A polynomial description of a convolutional encoder has either two or three components, depending on whether the encoder is a feedforward or feedback type:

- Constraint lengths
- Generator polynomials
- Feedback connection polynomials (for feedback encoders only)

### Constraint Lengths

The constraint lengths of the encoder form a vector whose length is the number of inputs in the encoder diagram. The elements of this vector indicate the number of bits stored in each shift register, *including* the current input bits.

In the figure above, the constraint length is three. It is a scalar because the encoder has one input stream, and its value is one plus the number of shift registers for that input.

### Generator Polynomials

If the encoder diagram has  $k$  inputs and  $n$  outputs, then the code generator matrix is a  $k$ -by- $n$  matrix. The element in the  $i$ th row and  $j$ th column indicates how the  $i$ th input contributes to the  $j$ th output.

For *systematic* bits of a systematic feedback encoder, match the entry in the code generator matrix with the corresponding element of the feedback connection vector. See “Feedback Connection Polynomials” below for details.

In other situations, you can determine the (i,j) entry in the matrix as follows:

- 1 Build a binary number representation by placing a 1 in each spot where a connection line from the shift register feeds into the adder, and a zero elsewhere. The leftmost spot in the binary number represents the current input, while the rightmost spot represents the oldest input that still remains in the shift register.
- 2 Convert this binary representation into an octal representation by considering consecutive triplets of bits, starting from the rightmost bit. The rightmost bit in each triplet is the least significant. If the number of bits is not a multiple of three, then place zero bits at the left end as necessary. (For example, interpret 1101010 as 001 101 010 and convert it to 152.)

For example, the binary numbers corresponding to the upper and lower adders in the figure above are 110 and 111, respectively. These binary numbers are equivalent to the octal numbers 6 and 7, respectively. Thus the generator polynomial matrix is [6 7].

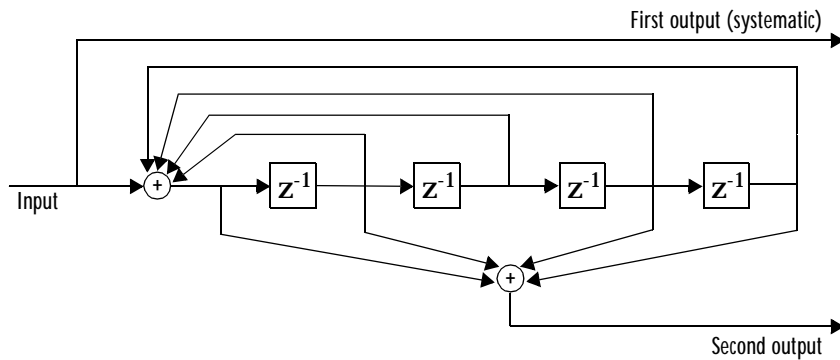
For a table of some good convolutional code generators, refer to [1] in the section “Selected Bibliography for Block Coding” on page 2-42, especially that book’s appendices.

### Feedback Connection Polynomials

If you are representing a feedback encoder, then you need a vector of feedback connection polynomials. The length of this vector is the number of inputs in the encoder diagram. The elements of this vector indicate the feedback connection for each input, using an octal format. First build a binary number representation as in step 1 above. Then convert the binary representation into an octal representation as in step 2 above.

If the encoder has a feedback configuration and is also systematic, then the code generator and feedback connection parameters corresponding to the systematic bits must have the same values.

For example, the diagram below shows a rate 1/2 systematic encoder with feedback.



This encoder has a constraint length of 5, a generator polynomial matrix of  $\begin{bmatrix} 37 & 33 \end{bmatrix}$ , and a feedback connection polynomial of 37. The first generator polynomial matches the feedback connection polynomial because the first output corresponds to the systematic bits.

### Using the Polynomial Description in MATLAB

To use the polynomial description with the functions `convenc` and `vitdec`, first convert it into a trellis description using the `poly2trellis` function. For example, the command below computes the trellis description of the encoder in Figure 2-1, Example of a Convolutional Encoder Diagram with Shift Registers, on page 2-44.

```
trellis = poly2trellis(3, [6 7]);
```

The MATLAB structure `trellis` is a suitable input argument for `convenc` and `vitdec`.

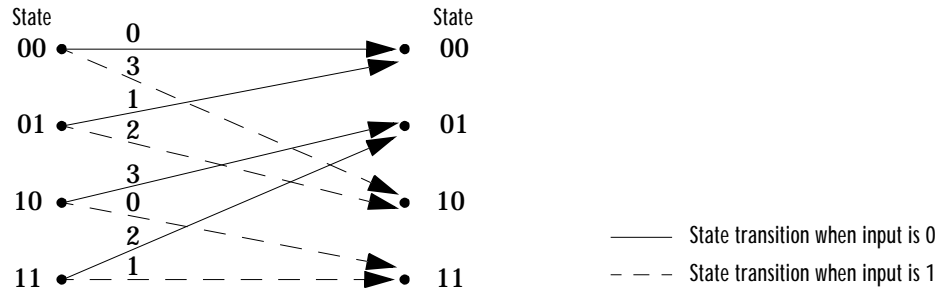
### Trellis Description of a Convolutional Encoder

A trellis description of a convolutional encoder shows how each possible input to the encoder influences both the output and the state transitions of the encoder. This section describes trellises, describes how to represent trellises in MATLAB, and gives an example of a MATLAB trellis.

The figure below depicts a trellis for the convolutional encoder from the previous section. The encoder has four states (numbered in binary from 00 to 11), a one-bit input, and a two-bit output. (The ratio of input bits to output bits makes this encoder a rate-1/2 encoder.) Each solid arrow shows how the



encoder changes its state if the current input is zero, and each dashed arrow shows how the encoder changes its state if the current input is one. The octal numbers above each arrow indicate the current output of the encoder.



**Figure 2-2: A Trellis for a 4-State Rate-1/2 Convolutional Encoder**

As an example of interpreting this trellis diagram, if the encoder is in the 10 state and receives an input of zero, then it outputs the code symbol 3 and changes to the 01 state. If it is in the 10 state and receives an input of one, then it outputs the code symbol 0 and changes to the 11 state.

Note that any polynomial description of a convolutional encoder is equivalent to some trellis description, although some trellises have no corresponding polynomial descriptions.

### Specifying a Trellis in MATLAB

To specify a trellis in MATLAB, use a specific form of a MATLAB structure called a trellis structure. A trellis structure must have five fields, as in the table below.

**Table 2-4: Fields of a Trellis Structure for a Rate  $k/n$  Code**

Field in Trellis Structure	Dimensions	Meaning
numInputSymbols	Scalar	Number of input symbols to the encoder: $2^k$
numOutputSymbols	Scalar	Number of output symbols from the encoder: $2^n$
numStates	Scalar	Number of states in the encoder

Table 2-4: Fields of a Trellis Structure for a Rate  $k/n$  Code (Continued)

Field in Trellis Structure	Dimensions	Meaning
nextStates	numStates-by- $2^k$ matrix	Next states for all combinations of current state and current input
outputs	numStates-by- $2^k$ matrix	Outputs (in decimal) for all combinations of current state and current input

**Note** While your trellis structure can have any name, its fields must have the *exact* names as in the table. Field names are case-sensitive.

In the nextStates matrix, each entry is an integer between 0 and numStates-1. The element in the  $i$ th row and  $j$ th column denotes the next state when the starting state is  $i-1$  and the input bits have decimal representation  $j-1$ . To convert the input bits to a decimal value, use the first input bit as the most significant bit (MSB). For example, the second column of the nextStates matrix stores the next states when the current set of input values is  $\{0, \dots, 0, 1\}$ . To learn how to assign numbers to states, see the reference page for `istrellis`.

In the outputs matrix, the element in the  $i$ th row and  $j$ th column denotes the encoder's output when the starting state is  $i-1$  and the input bits have decimal representation  $j-1$ . To convert to decimal value, use the first output bit as the MSB.

How to Create a MATLAB Trellis Structure

Once you know what information you want to put into each field, you can create a trellis structure in any of these ways:

- Define each of the five fields individually, using `structurename.fieldname` notation. For example, set the first field of a structure called `s` using the command below. Use additional commands to define the other fields.  
`s.numInputSymbols = 2;`  
The reference page for the `istrellis` function illustrates this approach.

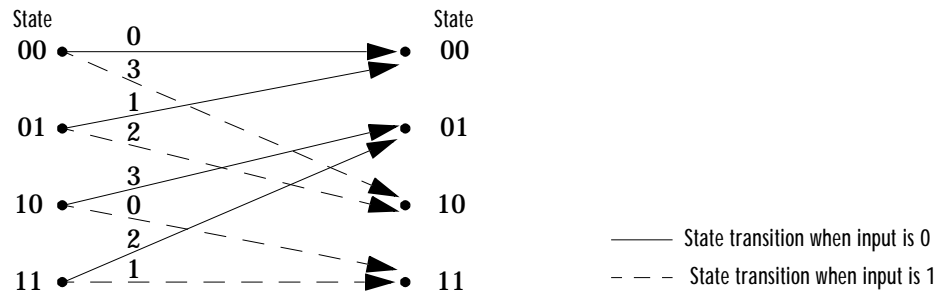
- Collect all field names and their values in a single struct command. For example:  

```
s = struct('numInputSymbols', 2, 'numOutputSymbols', 2, ...
          'numStates', 2, 'nextStates', [0 1; 0 1], 'outputs', [0 0; 1 1]);
```
- Start with a polynomial description of the encoder and use the `poly2trellis` function to convert it to a valid trellis structure. The polynomial description of a convolutional encoder is described in “Polynomial Description of a Convolutional Encoder” on page 2-43.

To check whether your structure is a valid trellis structure, use the `istrellis` function.

### Example: A MATLAB Trellis Structure

Reconsider the trellis shown in Figure 2-2, A Trellis for a 4-State Rate-1/2 Convolutional Encoder, which is repeated below.



To build a trellis structure that describes it, use the command below.

```
trellis = struct('numInputSymbols', 2, 'numOutputSymbols', 4, ...
                'numStates', 4, 'nextStates', [0 2; 0 2; 1 3; 1 3], ...
                'outputs', [0 3; 1 2; 3 0; 2 1]);
```

The number of input symbols is 2 because the trellis diagram has two types of input path, the solid arrow and the dashed arrow. The number of output symbols is 4 because the numbers above the arrows can be either 0, 1, 2, or 3. The number of states is 4 because there are four bullets on the left side of the trellis diagram (equivalently, four on the right side). To compute the matrix of next states, create a matrix whose rows correspond to the four current states on the left side of the trellis, whose columns correspond to the inputs of 0 and

1, and whose elements give the next states at the end of the arrows on the right side of the trellis. To compute the matrix of outputs, create a matrix whose rows and columns are as in the next states matrix, but whose elements give the octal outputs shown above the arrows in the trellis.

## Creating and Decoding Convolutional Codes

The functions for encoding and decoding convolutional codes are `convenc` and `vitdec`. This section discusses using these functions to create and decode convolutional codes.

### Encoding

A simple way to use `convenc` to create a convolutional code is shown in the commands below.

```
t = poly2trellis([4 3], [4 5 17; 7 4 2]); % Define trellis.
code = convenc(ones(100, 1), t); % Encode a string of ones.
```

The first command converts a polynomial description of a feedforward convolutional encoder to the corresponding trellis description. The second command encodes 100 bits, or 50 two-bit symbols. Since the code rate in this example is  $2/3$ , the output vector `code` contains 150 bits (that is, 100 input bits times  $3/2$ ).

### Hard-Decision Decoding

To decode using hard decisions, use the `vitdec` function with the flag `'hard'` and with *binary* input data. Since the output of `convenc` is binary, hard-decision decoding can use the output of `convenc` directly, without additional processing. This example extends the previous example and implements hard decision decoding.

```
t = poly2trellis([4 3], [4 5 17; 7 4 2]); % Define trellis.
code = convenc(ones(100, 1), t); % Encode a string of ones.
tb = 2; % Traceback length for decoding
decoded = vitdec(code, t, tb, 'trunc', 'hard'); % Decode.
```

## Soft-Decision Decoding

To decode using soft decisions, use the `vitdec` function with the flag 'soft'. You must also specify the number, `nsdec`, of soft-decision bits and use input data consisting of integers between 0 and

$$2^{nsdec} - 1$$

An input of 0 represents the most confident 0, while an input of  $2^{nsdec}-1$  represents the most confident 1. Other values represent less confident decisions. For example, the table below lists interpretations of values for 3-bit soft decisions.

**Table 2-5: Input Values for 3-bit Soft Decisions**

Input Value	Interpretation
0	Most confident 0
1	Second most confident 0
2	Third most confident 0
3	Least confident 0
4	Least confident 1
5	Third most confident 1
6	Second most confident 1
7	Most confident 1

**Example: Soft-Decision Decoding.** The script below illustrates decoding with 3-bit soft decisions. First it creates a convolutional code with `convenc` and adds white Gaussian noise to the code with `awgn`. Then, to prepare for soft-decision decoding, the example uses `quantiz` to map the noisy data values to appropriate decision-value integers between 0 and 7. The second argument in `quantiz` is a partition vector that determines which data values map to 0, 1, 2, etc. The partition is chosen so that values near 0 map to 0, and values near 1 map to 7. (You can refine the partition to obtain better decoding performance if your application requires it.) Finally, the example decodes the code and computes the bit error rate. Notice that when comparing the decoded data with the original message, the example must take the decoding delay into account.

The continuous operation mode of `vitdec` causes a delay equal to the traceback length, so `msg(1)` corresponds to `decoded(tblen+1)` rather than to `decoded(1)`.

```
msg = randi nt(4000, 1, 2, 139); % Random data
t = poly2trellis(7, [171 133]); % Define trellis.
code = convenc(msg, t); % Encode the data.
ncode = awgn(code, 6, 'measured', 244); % Add noise.

% Quantize to prepare for soft-decision decoding.
qcode = quantiz(ncode, [0.001, .1, .3, .5, .7, .9, .999]);

tblen = 48; delay = tblen; % Traceback length
decoded = vitdec(qcode, t, tblen, 'cont', 'soft', 3); % Decode.

% Compute bit error rate.
[number, ratio] = biterr(decoded(delay+1: end), msg(1: end-delay))
```

The output is below.

```
number =

    5

ratio =

    0.0013
```

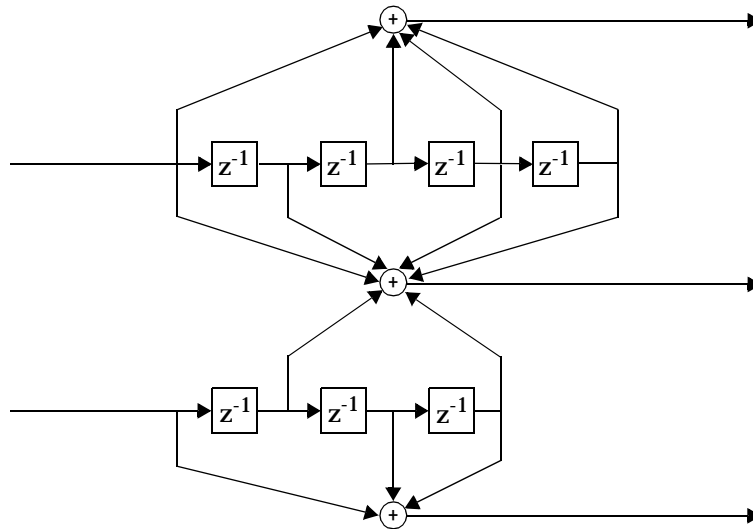
## Examples of Convolutional Coding

This section contains more examples of convolutional coding:

- The first example determines the correct trellis parameter for its encoder and then uses it to process a code. The decoding process uses hard decisions and the continuous operation mode. This operation mode causes a decoding delay, which the error rate computation takes into account.
- The second example processes a punctured convolutional code. The decoding process uses the unquantized decision type.

### Example: A Rate-2/3 Feedforward Encoder

The example below uses the rate 2/3 feedforward encoder depicted in the schematic below. The accompanying description explains how to determine the trellis structure parameter from a schematic of the encoder and then how to perform coding using this encoder.



**Figure 2-3: Schematic for a Rate 2/3 Feedforward Convolutional Encoder**

**Determining Coding Parameters.** The `convenc` and `vitdec` functions can implement this code if their parameters have the appropriate values.

The encoder's constraint length is a vector of length 2 since the encoder has two inputs. The elements of this vector indicate the number of bits stored in each shift register, including the current input bits. Counting memory spaces in each shift register in the diagram and adding one for the current inputs leads to a constraint length of [5 4].

To determine the code generator parameter as a 2-by-3 matrix of octal numbers, use the element in the  $i$ th row and  $j$ th column to indicate how the  $i$ th input contributes to the  $j$ th output. For example, to compute the element in the second row and third column, notice that the leftmost and two rightmost elements in the second shift register of the diagram feed into the sum that forms the third output. Capture this information as the binary number 1011,

which is equivalent to the octal number 13. The full value of the code generator matrix is `[27 33 0; 0 5 13]`.

To use the constraint length and code generator parameters in the `convenc` and `vitdec` functions, use the `poly2trellis` function to convert those parameters into a trellis structure. The command to do this is below.

```
trel = poly2trellis([5 4], [27 33 0; 0 5 13]); % Define trellis.
```

Using the Encoder. Below is a script that uses this encoder.

```
len = 1000;
msg = randint(2*len, 1); % Random binary message of 2-bit symbols
trel = poly2trellis([5 4], [27 33 0; 0 5 13]); % Trellis
code = convenc(msg, trel); % Encode the message.
ncode = rem(code + randerr(3*len, 1, [0 1; .96 .04]), 2); % Add noise.
decoded = vitdec(ncode, trel, 34, 'cont', 'hard'); % Decode.
[number, ratio] = biterr(decoded(68+1:end), msg(1:end-68));
```

Notice that `convenc` accepts a vector containing 2-bit symbols and produces a vector containing 3-bit symbols, while `vitdec` does the opposite. Also notice that `biterr` ignores the first 68 elements of `decoded`. That is, the decoding delay is 68, which is the number of bits per symbol (2) of the recovered message times the traceback depth value (34) in the `vitdec` function. The first 68 elements of `decoded` are zeros, while subsequent elements represent the decoded messages.

### Example: A Punctured Convolutional Code

This example processes a punctured convolutional code. It begins by generating 3000 random bits and encoding them using a rate-1/2 convolutional encoder. The resulting vector contains 6000 bits, which are mapped to values of -1 and 1 for transmission. The puncturing process removes every third value and results in a vector of length 4000. The punctured code, `punctcode`, passes through an additive white Gaussian noise channel. Afterwards, the example inserts values to reverse the puncturing process. While the puncturing process removed both -1s and 1s from `code`, the insertion process inserts zeros. Then `vitdec` decodes the vector of -1s, 1s, and 0s using the 'unquant' decision type. This unquantized decision type is appropriate here for these reasons:

- `tcode` uses -1 to represent the 1s in `code`.
- `tcode` uses 1 to represent the 0s in `code`.



- The inserted 0s are acceptable for the 'unquant' decision type, which allows any real values as input.

Finally, the example computes the bit error rate and the number of bit errors.

```
len = 3000; msg = randint(len, 1, 2, 94384); % Random data
t = poly2trellis(7, [171 133]); % Define trellis.
code = convenc(msg, t); % Length is 2*len.
tcode = -2*code+1; % Transmit -1s and 1s.

% Puncture by removing every third value.
punctcode = tcode;
punctcode(3:3:end)=[]; % Length is (2*len)*2/3.

ncode = awgn(punctcode, 8, 'measured', 1234); % Add noise.

% Insert zeros.
nicode = zeros(2*len, 1); % Zeros represent inserted data.
nicode(1:3:end) = ncode(1:2:end); % Write actual data.
nicode(2:3:end) = ncode(2:2:end); % Write actual data.

decoded = vitdec(nicode, t, 96, 'trunc', 'unquant'); % Decode.
[number, ratio]=biterr(decoded, msg); % Bit error rate
```

## Selected Bibliography for Convolutional Coding

- [1] Clark, George C. Jr. and J. Bibb Cain. *Error-Correction Coding for Digital Communications*. New York: Plenum Press, 1981.
- [2] Gitlin, Richard D., Jeremiah F. Hayes, and Stephen B. Weinstein. *Data Communications Principles*. New York: Plenum, 1992.

# Modulation

In most media for communication, only a fixed range of frequencies is available for transmission. One way to communicate a message signal whose frequency spectrum does not fall within that fixed frequency range, or one that is otherwise unsuitable for the channel, is to alter a transmittable signal according to the information in your message signal. This alteration is called *modulation*, and it is the modulated signal that you transmit. The receiver then recovers the original signal through a process called *demodulation*.

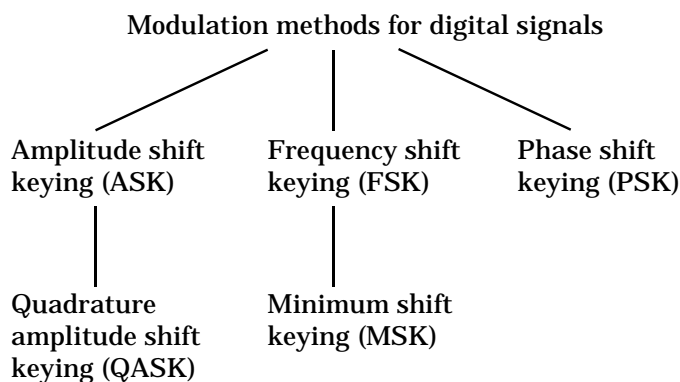
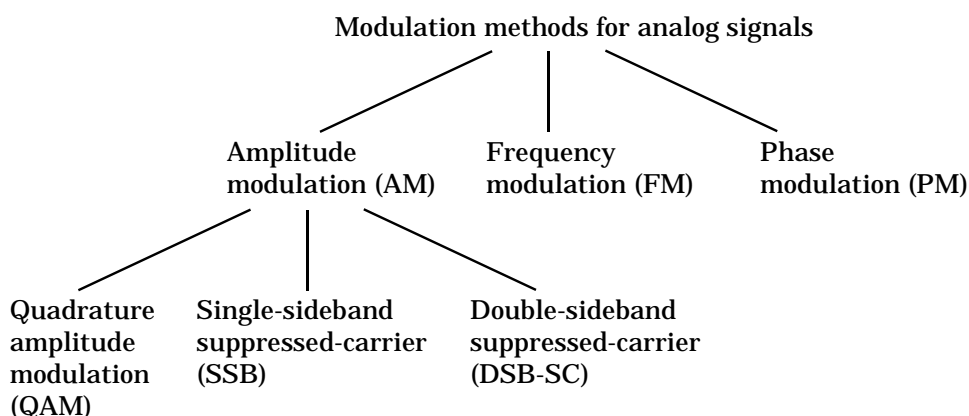
The table shows how this section is organized.

Subject	Topics
General modulation	“Modulation Features of the Toolbox” on page 2-57
	“Modulation Terminology” on page 2-58
Analog modulation	“Representing Analog Signals” on page 2-59
	“Simple Analog Modulation Example” on page 2-61
	“Other Options in Analog Modulation” on page 2-62
	“Filter Design Issues” on page 2-62
Digital modulation	“Digital Modulation Overview” on page 2-66
	“Representing Digital Signals” on page 2-67
	“Significance of Sampling Rates” on page 2-70
	“Representing Signal Constellations” on page 2-70
	“Simple Digital Modulation Example” on page 2-74
	“Customizing the Modulation Process” on page 2-75
	“Other Options in Digital Modulation” on page 2-77

For background information about modulation and demodulation, see the works listed in “Selected Bibliography for Modulation” on page 2-77.

## Modulation Features of the Toolbox

The available methods of modulation depend on whether the input signal is analog or digital. The figures below show the modulation techniques that the Communications Toolbox supports for analog and digital signals, respectively. As the figures suggest, some categories of techniques include named special cases.



### Baseband Versus Passband Simulation

For a given modulation technique, two ways to simulate modulation techniques are called *baseband* and *passband*. Baseband simulation, also known as the *lowpass equivalent method*, requires less computation. This toolbox supports both baseband and passband simulation. Since baseband simulation is more prevalent, this guide focuses more on baseband simulation.

---

**Note** To use this toolbox for passband simulation, see the reference pages for the functions `amod`, `ademod`, `dmod`, and `ddemod`.

---

### Supported Modulation Tasks

Functions in the toolbox can accomplish these tasks:

- Modulate a signal using one of the techniques shown in the figures above
- Demodulate a signal using one of the techniques shown in the figures above
- Map a digital signal to an analog signal, before modulation
- Demap an analog signal to a digital signal, after demodulation
- Map, demap, and plot constellations for QASK modulation

The modulation and demodulation functions also let you control such features as the initial phase of the modulated signal, post-demodulation filtering, and the decision timing for digital demodulation.

### Modulation Terminology

Modulation is a process by which a *carrier signal* is altered according to information in a *message signal*. The *carrier frequency*, denoted  $F_c$ , is the frequency of the carrier signal. The *sampling rate* is the rate at which the message signal is sampled during the simulation.

The frequency of the carrier signal is usually much greater than the highest frequency of the input message signal. The Nyquist sampling theorem requires that the simulation sampling rate  $F_s$  be greater than two times the highest frequency of the modulated signal, in order for the demodulator to recover the message correctly. The sampling rate  $F_s$  of a modulated digital signal is greater than or equal to the sampling rate  $F_d$  of the original message signal before modulation.

The table below lists the requirements in terms of the input arguments for this toolbox's modulation and demodulation functions. Note that the situations are not mutually exclusive.

Situation	Requirement
Passband simulation	$2 * (\text{highest frequency of modulated signal}) < F_s$
Digital signals	$F_d \leq F_s$
Passband simulation, digital signals	$F_d < F_c$

## Representing Analog Signals

To perform baseband modulation of an analog signal using this toolbox, start with a real message signal and a sampling rate  $F_s$  in Hertz. For modulation techniques *other than* quadrature amplitude modulation (QAM), represent the signal using a vector  $x$ , the entries of which give the signal's values in time increments of  $1/F_s$ . Baseband modulation (using a technique other than QAM) produces a complex vector.

For example, if  $t$  measures time in seconds, then the vector  $x$  below is the result of sampling a frequency-one sine wave 100 times per second for 2 seconds. The vector  $y$  represents the modulated signal. The output shows that  $y$  is complex.

```

Fs = 100; % Sampling rate is 100 samples per second.
t = [0:1/Fs:2]'; % Sampling times for 2 seconds
x = sin(2*pi*t); % Representation of the signal
y = amodce(x, Fs, 'pm'); % Modulate x to produce y.
whos
  Name      Size      Bytes  Class
  Fs        1x1         8    double array
  t         201x1      1608   double array
  x         201x1      1608   double array
  y         201x1     3216   double array (complex)

```

Grand total is 604 elements using 6440 bytes

### Baseband Modulated Signals Defined

This section explains the connection between this complex vector  $y$  and the real signal that you might expect to get after modulating a real signal. If the modulated signal has the waveform

$$Y_1(t) \cos(2\pi f_c t + \theta) - Y_2(t) \sin(2\pi f_c t + \theta)$$

where  $f_c$  is the carrier frequency and  $\theta$  is the carrier signal's initial phase, then a baseband simulation recognizes that this equals

$$\text{Re} \left[ (Y_1(t) + jY_2(t)) e^{j\theta} e^{j2\pi f_c t} \right]$$

and models only the part inside the curly brackets. Here  $j$  is the square root of -1. The complex vector  $y$  is a sampling of the complex signal  $(Y_1(t) + jY_2(t)) \exp(j\theta)$ .

---

**Note** You can also simultaneously process several signals of equal length. To do this, make  $x$  a matrix in which each signal occupies one column. The corresponding modulated signal  $y$  is a complex matrix whose  $k$ th column is the modulation of the  $k$ th column of  $x$ .

---

### Changes for QAM

The case for quadrature amplitude modulation (QAM) is similar, except that the message signal has in-phase and quadrature components. Represent the signal using a matrix  $x$  that has an even number of columns. The odd-indexed columns represent in-phase components of the signal and the even-indexed columns represent quadrature components. If the message signal is a  $2n$ -by- $m$  matrix, then the modulated signal is an  $n$ -by- $m$  matrix. As in the other methods, baseband modulation turns a real message signal into a complex modulated signal.

For example, the code below implements QAM on a set of sinusoidal input signals.

```
Fs = 100; % Sampling rate is 100 samples per second.
t = [0:1/Fs:2]'; % Sampling times
% Signal is a four column matrix.
```

```
% Each column models a sinusoidal signal, the frequencies
% of which are 1 Hz, 1.5 Hz, 2 Hz, 2.5 Hz respectively.
x = sin([2*pi*t, 3*pi*t, 4*pi*t, 5*pi*t]);
y = amodce(x, Fs, 'qam'); % Modulate x to produce y.
```

The output below shows the sizes and types of x and y.

```
whos
      Name      Size      Bytes  Class

      Fs         1x1          8  double array
      t        201x1        1608  double array
      x        201x4        6432  double array
      y        201x2        6432  double array (complex)
```

Grand total is 1408 elements using 14480 bytes

## Simple Analog Modulation Example

This example illustrates the basic format of the baseband modulation and demodulation commands, `amodce` and `ademodce`. Although the example uses the AMDSB-TC method, most elements of this example apply to other analog modulation techniques as well. The example samples an analog signal and modulates it. Then it demodulates it and displays the order of magnitude of the variance between the original and demodulated signals.

```
% Sample the signal for two seconds,
% at a rate of 100 samples per second.
Fs = 100;
t = [0:1/Fs:2]';
% The signal is a sum of sinusoids.
x = sin(2*pi*t) + sin(4*pi*t);
% Use AMDSB-TC modulation to produce y.
y = amodce(x, Fs, 'amdsb-tc');
% Demodulate y to recover the message.
z = ademodce(y, Fs, 'amdsb-tc');
v = floor(log10(var(x-z)))

v =
```

### Other Options in Analog Modulation

The table below lists a few ways in which you might vary the simple example in the previous section in order to perform the modulation and demodulation slightly differently. See the reference pages for full details about options.

Table 2-6: Substitutions in “Simple Analog Modulation Example”

Modification of Process	Modifications in the Code
Set the carrier signal's initial phase to phase, measured in radians	<code>y = amodce(x, [Fs phase], 'amdsb-tc');</code> <code>z = ademodce(y, [Fs phase], 'amdsb-tc');</code>
Use a lowpass filter after demodulating. num and den are row vectors that give the coefficients, in <i>descending</i> order, of the numerator and denominator of the filter's transfer function.	<code>z = ademodce(y, Fs, 'amdsb-tc', 0, num, den);</code>  (For other demodulation methods, the 0 in the statement above would be unnecessary. See the reference page for ademodce for details.)
(AM-SSB only) Use a Hilbert filter in the time domain. num and den are as above.	<code>y = amodce(x, Fs, 'amssb/time', num, den);</code> <code>z = ademodce(y, Fs, 'amssb');</code>
(AMDSB only) Use a Costas phase-locked loop	<code>z = ademodce(y, Fs, 'amdsb-tc/costas');</code>  <i>or</i>  <code>y = amodce(x, Fs, 'amdsb-sc');</code> <code>z = ademodce(y, Fs, 'amdsb-sc/costas');</code>
(AMDSB-TC only) Shift the signal values by offset before modulating and after demodulating	<code>y = amodce(x, Fs, 'amdsb-tc', offset);</code> <code>z = ademodce(y, Fs, 'amdsb-tc', offset);</code>

### Filter Design Issues

After demodulating, you might want to filter out the carrier signal, especially if you are using passband simulation. The Signal Processing Toolbox provides functions that can help you design your filter, such as `butter`, `cheby1`, `cheby2`, and `ellip`. Different demodulation methods have different properties, and you



might need to test your application with several filters before deciding which is most suitable. This subsection mentions two issues that relate to the use of filters: cutoff frequency and time lag.

### Example: Varying the Filter's Cutoff Frequency

In many situations, a suitable cutoff frequency is half the carrier frequency. Since the carrier frequency must be higher than the bandwidth of the message signal, a cutoff frequency chosen in this way limits the bandwidth of the message signal. If the cutoff frequency is too high, then the carrier frequency may not be filtered out. If the cutoff frequency is too low, then it might narrow the bandwidth of the message signal.

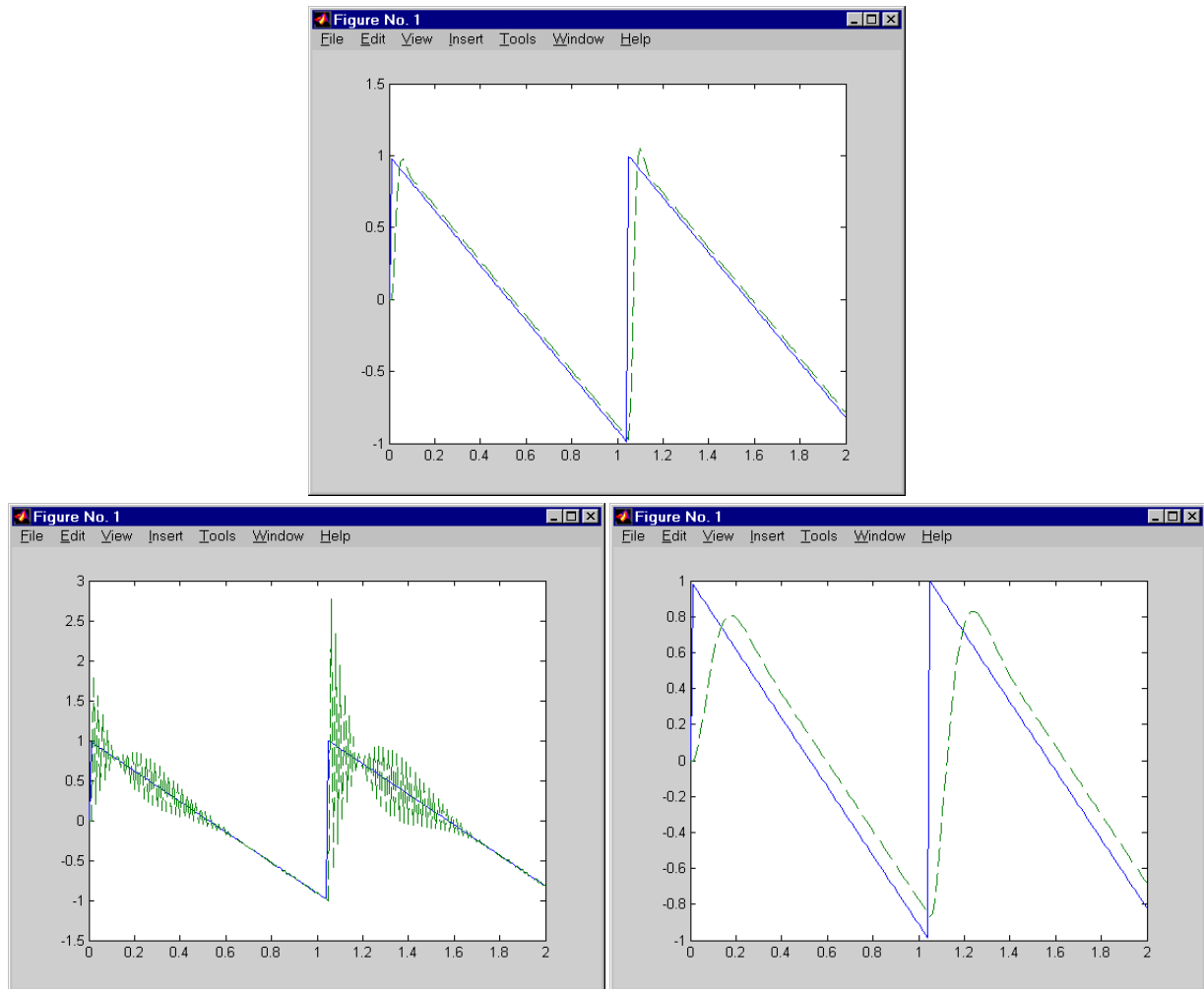
The code below modulates a sawtooth message signal, demodulates the resulting signal using a Butterworth filter, and plots the original and recovered signals. Note that the scaling in the butter function causes the cutoff frequency of the filter to be  $F \cdot F_s/2$ , not  $F$  itself.

```

Fc = 25; % Carrier frequency
Fs = 100; % Signal sampling rate
t = [0:1/Fs:2]'; % Times to sample the signal
x = sawtooth(6*t, 0); % Signal is a sawtooth.
y = amod(x, Fc, Fs, 'amssb'); % Modulate.
F = Fc/Fs; % Change F to vary the filter's cutoff frequency.
[num, den] = butter(2, F); % Design Butterworth filter.
z = ademod(y, Fc, Fs, 'amssb', num, den); % Demodulate and filter.
plot(t, x, '-', t, z, '--') % Plot original and recovered signals.

```

The plots below show the effects of three lowpass filters with different cutoff frequencies. In each plot, the dotted curve is the demodulated signal and the solid curve is the original message signal. The top plot uses the suggested cutoff frequency ( $F = F_c/F_s$ ). The lower left plot uses a higher cutoff frequency ( $F = 3.9 \cdot F_c/F_s$ ), which allows the carrier signal to interfere with the demodulated signal. The lower right plot uses a lower cutoff frequency ( $F = F_c/F_s/4$ ), which narrows the bandwidth of the demodulated signal.



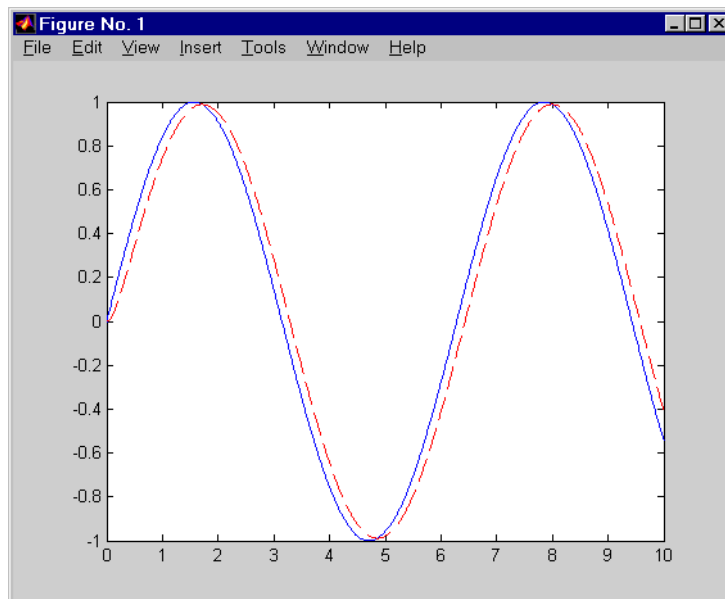
**Figure 2-4: Original and Recovered Signals, with Filter Cutoff  $F = F_c/F_s$ ,  $3.9F_c/F_s$ , and  $F_c/F_s/4$**

### Example: Time Lag From Filtering

There is invariably a time delay between a demodulated signal and the original received signal. Both the filter order and the filter parameters directly affect the length of this delay. The example below illustrates the time delay by

plotting a signal before and after the modulation, demodulation, and filtering processes. The solid curve is the original sine wave and the dashed curve is the recovered signal.

```
Fs = 100; % Sampling rate of signal
[num, den] = butter(2, 0.8); % Design Butterworth filter.
t = [0:1/Fs:10]'; % Times to sample the signal
x = sin(t); % Signal is a sine wave.
y = amodce(x, Fs, 'pm'); % Modulate.
z = ademodce(y, Fs, 'pm', num, den); % Demodulate and filter.
plot(t, x, t, z, 'r--') % Plot original signal and recovered signal.
```



### Digital Modulation Overview

Modulating a digital signal can be interpreted as a combination of two steps: mapping the digital signal to an analog signal and modulating the analog signal. These are depicted in the schematic below.

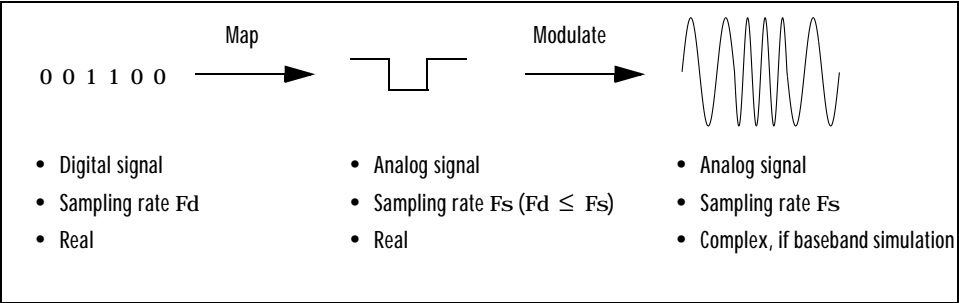


Figure 2-5: Two Steps of Digital Modulation

Except for FSK and MSK methods, when the receiver tries to recover a digital message from the analog signal that it receives, it performs two steps: demodulating the analog signal and demapping the demodulated analog signal to produce a digital message. These are depicted in the schematic below.

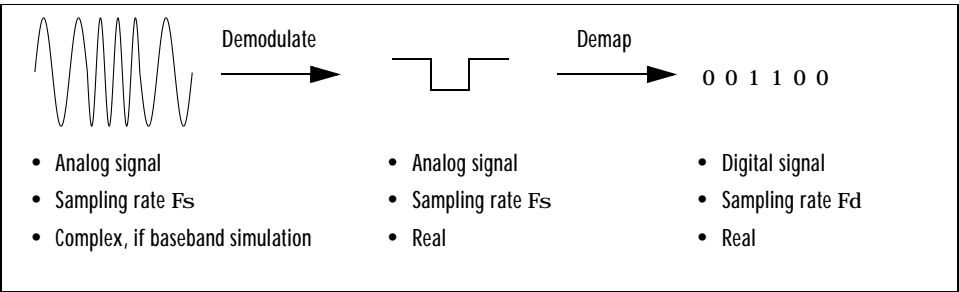


Figure 2-6: Two Steps of Digital Demodulation

For FSK and MSK methods, the demodulator uses correlation techniques instead of the two-stage process above.

The mapping process increases the sampling rate of the signal from  $F_d$  to  $F_s$ , whereas the demapping process decreases the sampling rate from  $F_s$  to  $F_d$ .

Functions in this toolbox can perform any of these steps, as summarized in the table below.

**Table 2-7: Functions for the Steps of Digital Modulation and Demodulation**

Step	Function
Mapping and modulation	dmodce or dmod
Mapping only	modmap
Modulation without mapping	dmodce or dmod, with <b>/nomap</b> flag
Demodulation and demapping	ddemodce or ddemod
Demodulation without demapping (ASK, PSK, or QASK)	ddemodce or ddemod, with <b>/nomap</b> flag
Demapping only	demodmap

The functions are described in more detail in the sections that follow.

## Representing Digital Signals

This section describes the formats for digital message signals, the analog signals to which they map, and the analog signals that result from the two-stage baseband digital modulation process. The last part, “Constellations and Mapped Signals (PSK, QASK),” discusses some special formats that apply to the PSK and QASK modulation methods.

### Message Signals

To perform M-ary baseband modulation of a digital signal using this toolbox, start with a message signal consisting of integers in the range [0, M-1]. Represent the signal using a vector **x**. Associate with the message signal a sampling rate **Fd**, which means that the entries of **x** give the signal’s values in time increments of 1/**Fd**.

### Mapped Signals

Mapping produces a real signal **y** whose sampling rate **Fs** must satisfy

$$F_s > F_d$$

(For passband simulation, in which the carrier frequency  $F_c$  appears explicitly, both of the relations  $F_s > F_c > F_d$  and  $F_s > 2F_c$  must hold.) If  $x$  consists of  $n$  samples, then  $y$  contains  $n \cdot F_s / F_d$  samples. The actual dimensions of  $y$  depend on the modulation scheme, as detailed in “To Map a Digital Signal (General Information)” on page 3-150.

For example, the vector  $x$  below samples a random digital signal 100 times per second for 2 seconds. The vector  $y$  represents the mapped signal, sampled three times as frequently. The output shows that  $y$  contains three times as many samples as  $x$ .

```
Fd = 100; % Sampling rate of x
M = 32; % Digital symbols are 0, 1, 2, ..., 31
x = randint(2*Fd, 1, M); % Representation of the digital signal
Fs = 3*Fd; % Sampling rate of mapped signal
y = modmap(x, Fd, Fs, 'ask', M); % Mapped signal
r = [size(x, 1) size(y, 1)] % Number of rows in x and y

r =

    200    600
```

Modulated Signals

Baseband modulation produces a complex signal with sampling rate  $F_s$ . Notice that this is the same sampling rate as the mapped signal. Baseband signals are explained briefly in the section, “Representing Analog Signals” on page 2-59; for more details, see the works listed in “Selected Bibliography for Modulation” on page 2-77. To illustrate the size and nature of the modulated signal, supplement the example in the paragraph above with these commands.

```
z = dmodce(x, Fd, [Fs pi/2], 'ask', M);
whos
```

Name	Size	Bytes	Class
Fd	1x1	8	double array
Fs	1x1	8	double array
M	1x1	8	double array
r	1x2	16	double array
x	200x1	1600	double array
y	600x1	4800	double array

```
z          600x1          9600 double array (complex)
```

Grand total is 1405 elements using 16040 bytes

### Constellations and Mapped Signals (PSK, QASK)

If you map a digital message using the phase shift keying (PSK) or quadrature amplitude shift keying (QASK) modulation method, then `modmap` describes the amplitude and phase of the resulting analog signal using an in-phase part and a quadrature part. For this reason, one column in the original message signal vector corresponds to *two* columns in the mapped signal matrix.

For example, compare the code below with the example in “Mapped Signals” above. The mapped signal `ypsk` is a two-column matrix, whereas the earlier ASK example produced a column vector. The first column of `ypsk` gives the in-phase components of the samples and the second column gives the quadrature components.

```
Fd = 100; % Sampling rate of x
M = 32; % Digital symbols are 0, 1, 2, ..., 31.
x = randint(2*Fd, 1, M); % Representation of the digital signal
Fs = 3*Fd; % Sampling rate of mapped signal
ypsk = modmap(x, Fd, Fs, 'psk', M); % PSK mapped signal
s = size(ypsk)

s =
```

```
600      2
```

**Using Signal Constellation Plots.** To understand the in-phase and quadrature description more easily, refer to a signal constellation plot. Each point in the constellation represents an analog signal to which `modmap` can map the digital message data. Each row of `y` in the example above gives the two rectangular coordinates of some point in the constellation. To produce a signal constellation plot that corresponds to the example above, use the command

```
modmap('psk', M) % Using M = 32 from before
```

More about creating signal constellation plots is in the section “Representing Signal Constellations” on page 2-70.

## Significance of Sampling Rates

The vectors and matrices that form the input and output of the modulation and demodulation functions do not have a built-in notion of time. That is, MATLAB does not know whether the digital signal `[0 1 2 3 4 5 6 7]` represents an 8-second signal sampled once per second, or a 1-second signal sampled eight times, or something else. However, many functions appearing in this “Modulation” section ask for one or more sampling rates. This subsection discusses the significance of these sampling rates.

If your application has a natural notion of time, then you are free to use it in the modulation and demodulation functions. For example, if you generate the digital signal `[0 1 2 3 4 5 6 7]` and know that it represents a 1-second signal sampled eight times, then set  $F_d = 8$ . On the other hand, if you know that the signal represents a 2-second signal sampled four times per second, then set  $F_d = 4$ . You can also use the formula

```
Fd = size(x,1) / (max(t)-min(t)); % if x=signal, t=sample times
```

for a signal  $x$  sampled at times  $t$ . Here  $x$  is a matrix or vector and  $t$  is a vector whose length is the number of rows of  $x$ .

For most digital modulation computations, MATLAB does not directly use the sampling rates  $F_d$  and  $F_s$  of digital message signals and mapped signals, respectively. What it uses is their *ratio*  $F_s/F_d$ . For example, the two commands below produce exactly the same result, because  $3/1$  equals  $6/2$ .

```
y13 = dmodce([0 1 2 3 4 5 6 7]', 1, 3, 'ask', 8);  
y26 = dmodce([0 1 2 3 4 5 6 7]', 2, 6, 'ask', 8);
```

One exceptional situation in which the individual value of  $F_d$  matters occurs in the MSK and M-ary FSK methods. The default separations between successive frequencies are  $F_d/2$  and  $F_d$  for these two methods, respectively.

## Representing Signal Constellations

The QASK method depends on a choice of a signal constellation. The QASK mapping and demapping functions in this toolbox can process two special types of signal constellations, as well as a general type of constellation that you can define as you choose. The special types are called square and circle constellations and the general type is called an arbitrary constellation. This section describes how you can tell MATLAB what signal constellation you want to use, and how you can plot signal constellations.



## Square Constellations

To use a square constellation, you only need to tell MATLAB the number of points in the constellation. This number,  $M$ , must be a power of two. For example, to map the digital signal [3 8 15 30 28] to a square constellation having 32 points, use the `qaskenco` function as below.

```
[inphase, quadr] = qaskenco([3 8 15 30 28], 32);
```

The returned vectors `inphase` and `quadr` give the in-phase and quadrature components, respectively, of the mapped signal. The command

```
msg = qaskdeco(inphase, quadr, 32);
```

demaps to recover the original message [3 8 15 30 28]. Notice that in both cases, the square constellation is described only by the number 32.

The modulation and demodulation functions use the  $M$ -ary number and the method string 'qask' to specify the square constellation. The command below implements QASK modulation on the message [3 8 15 30 28], using a 32-point square constellation. The command assumes that the sampling rates are 1 Hz before modulating and 2 Hz after modulating.

```
y = dmodce([3 8 5 30 28], 1, 2, 'qask', 32);
```

**Plotting Square Constellations.** To plot a square constellation with  $M$  points, use one of these commands.

```
qaskenco(M)
modmap('qask', M);
```

## Circle Constellations

To use a circle constellation having equally spaced points on each circle, you need to give MATLAB this information, in this order:

- 1 The number of points on each circle
- 2 The radius of each circle
- 3 The phase of one point on each circle

The three types of information occupy three vectors of the same length. The first entries of the three vectors determine one circle, the second entries of the three vectors determine another circle, and so on.

For example, the `apkconst` command below returns the complex coordinates of the points on a circle constellation that contains sixteen points on each of two circles. The inner circle has radius one, and one of the constellation points has zero phase. The outer circle has radius three and a constellation point at 10 degrees.

```
y = apkconst([16 16], [1 3], [0 10*pi/180]);
```

The constellation contains two circles because each vector has length two. The constellation has 32 points in total because the sum of entries in the first vector is 32.

The modulation and demodulation functions use three equal-length vectors and the method string 'qask/ci r' to specify the circle constellation. The command below implements QASK modulation on the message [3 8 15 30 28], using the circle constellation described above.

```
y = dmodce([3 8 5 30 28], 1, 2, 'qask/ci r', [16 16], [1 3], ...  
[0 10*pi/180]);
```

**Default Values.** If you do not provide the phase vector, then by default one constellation point on each circle will have zero phase. If you provide neither the phase vector nor the radius vector, then by default the  $k$ th circle will have radius  $k$ , and one of the constellation points will have zero phase. You must provide the vector that specifies how many points are on each circle.

**Plotting Circle Constellations.** To plot a circle constellation in which `numsig` gives the number of points on each circle, `amp` gives the radius of each circle, and `phs` gives the phase of one point on each circle, use one of these commands.

```
apkconst(numsig, amp, phs)  
modmap('qask/ci r', numsig, amp, phs);
```

To label the constellation points by number, use this syntax instead.

```
apkconst(numsig, amp, phs, 'n')
```

### Arbitrary Constellations

You can also use a signal constellation that does not fit into the categories above. To do this, you need to give MATLAB two real vectors of equal length, one that contains the in-phase components of the constellation point and one that contains the corresponding quadrature components. You also need to use

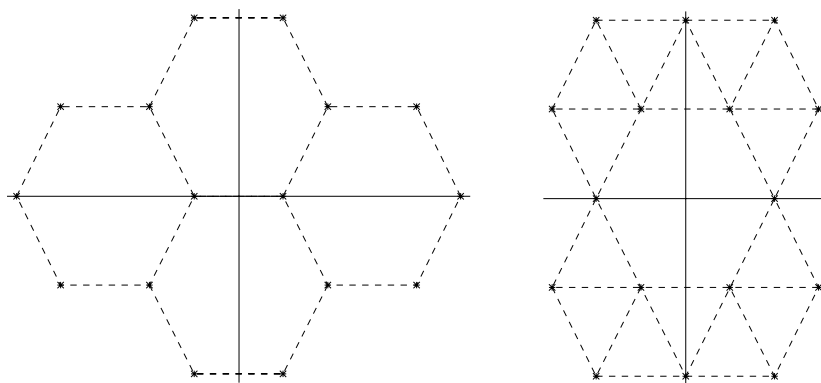
the method string 'qask/arb' in the modulation, demodulation, mapping, and demapping functions.

For example, the code examples below plot signal constellations that have a hexagonal and triangular structure, respectively. They use the `modmap` function.

```
% Example #1: A hexagonal constellation
inphase = [1/2 1 1 1/2 1/2 2 2 5/2];
quadr = [0 1 -1 2 -2 1 -1 0];
inphase = [inphase; -inphase]; inphase = inphase(:);
quadr = [quadr; quadr]; quadr = quadr(:);
modmap('qask/arb', inphase, quadr);
```

```
% Example #2: A triangular constellation
figure;
inphase = [1/2 -1/2 1 0 3/2 -3/2 1 -1];
quadr = [1 1 0 2 1 1 2 2];
inphase = [inphase; -inphase]; inphase = inphase(:);
quadr = [quadr; -quadr]; quadr = quadr(:);
modmap('qask/arb', inphase, quadr);
```

The figure below shows plots of the hexagonal and triangular signal constellations on the left and right, respectively. The dashed lines are not part of MATLAB's output, and appear below only to suggest the hexagonal and triangular structures.



The modulation and demodulation functions also use the method string 'qask/arb' and a pair of equal-length vectors like `inphase` and `quadr` to determine your constellation. For example, to modulate the message [3 8 5 10 7] using the QASK method with one of the constellations described in the examples above, supplement the example code with this command.

```
y = dmodce([3 8 5 10 7], 1, 2, 'qask/arb', inphase, quadr);
```

### Simple Digital Modulation Example

This example illustrates the basic format of the baseband modulation and demodulation commands, `dmodce` and `ddemodce`. Although the example uses the PSK method, most elements of this example apply to digital modulation techniques other than PSK.

The example generates a random digital signal, modulates it, and adds noise. Then it creates a scatter plot, demodulates the noisy signal, and computes the symbol error rate. The `ddemodce` function demodulates the analog signal `y` and then demaps to produce the digital signal `z`.

Notice that the scatter plot does not look exactly like a signal constellation. Whereas the signal constellation would have 16 precisely located points, the noise causes the scatter plot to have a small cluster of points approximately where each constellation point would be. However, the noise is sufficiently small that the signal can be recovered perfectly.

---

**Note** Since some options vary by method, you should check the reference pages before adapting the code here for other uses.

---

Below are the code and the scatter plot.

```
M = 16; % Use 16-ary modulation.
Fd = 1; % Assume the original message is sampled
% at a rate of 1 sample per second.
Fs = 3; % The modulated signal will be sampled
% at a rate of 3 samples per second.
x = randint(100, 1, M); % Random digital message
% Use M-ary PSK modulation to produce y.
y = dmodce(x, Fd, Fs, 'psk', M);
% Add some Gaussian noise.
```

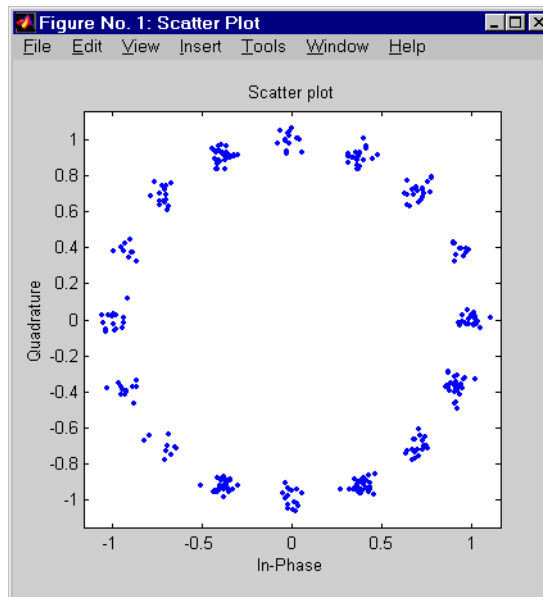
```

ynoi sy = y + .04*randn(300,1) + .04*j*randn(300,1);
% Create scatter plot from noisy data.
scatterplot(ynoi sy, 1, 0, 'b. ');
% Demodulate y to recover the message.
z = ddemodce(ynoi sy, Fd, Fs, 'psk', M);
s = symerr(x, z) % Check symbol error rate.

```

s =

0



## Customizing the Modulation Process

Recall from “Digital Modulation Overview” on page 2-66 that the modulation and demodulation processes each consist of two steps. You can tell the toolbox functions to carry out only selected steps in the processes. For example, this might be useful if you want to use standard mapping and demapping techniques along with unusual or proprietary modulation and demodulation techniques.

Mapping Without Modulating and Demapping Without Demodulating

To map the digital signal to an analog signal without modulating the analog signal, use the `modmap` function instead of the `dmodce` function. To demap the analog signal to a digital signal without demodulating the analog signal, use the `demodmap` function instead of the `ddemodce` function.

To alter the basic example so that it does not modulate or demodulate the analog signals at all, replace the “old commands” listed in the first column of the table below with the “new commands” listed in the second column.

Table 2-8: Changes in “Simple Digital Modulation Example” to Avoid Modulating

Old Command	New Command
<code>y = dmodce(x, Fd, Fs, 'psk', M);</code>	<code>y = modmap(x, Fd, Fs, 'psk', M);</code>
<code>ynoi sy = y + .04*randn(300, 1) + .04*j*randn(300, 1);</code>	<code>ynoi sy = y + .04*randn(300, 2) + .04*j*randn(300, 2);</code>
<code>z = ddemodce(y, Fd, Fs, 'psk', M);</code>	<code>z = demodmap(y, Fd, Fs, 'psk', M);</code>

Modulating Without Mapping and Demodulating Without Demapping

To carry out the analog modulation step on a signal that has already been mapped from a digital signal to an analog signal, use the `dmodce` function with the extra word `/nomap` appended to the method string. To carry out the analog demodulation step but avoid demapping the resulting signal to a digital signal, use the `ddemodce` function with the extra word `/nomap` appended to the method string.

If you substituted your own mapping and demapping steps into the basic example then it would look something like the code below. The lines in the second grouping differ from the original example.

```
M = 16; % Use 16-ary modulation.
Fd = 1; % Assume the original message is sampled
% at a rate of 1 sample per second.
Fs = 3; % The modulated signal will be sampled
% at a rate of 3 samples per second.
x = randint(100, 1, M); % Random digital message

% Important changes are below.
mapx = mymappingfunction(x); % Use your own function here.
```

```

y = dmodce(mapx, Fd, Fs, 'psk/nomap', M); % Modulate without mapping.
% Demodulate y without demapping.
demody = ddemodce(y, Fd, Fs, 'psk/nomap', M);
% Now demap.
z = mydemappingfunction(demody); % Use your own function here.

```

## Other Options in Digital Modulation

The table below lists a few ways in which you might vary the example in the section “Simple Digital Modulation Example” on page 2-74 in order to perform the modulation and demodulation slightly differently. See the reference pages for full details about options.

**Table 2-9: Substitutions in the Digital Example**

Modification of Process	Modifications in the Code in “Simple Digital Modulation Example” on page 2-74
Set the carrier signal's initial phase to phase, measured in radians	<pre> y = dmodce(x, Fd, [Fs phase], 'psk', M); z = ddemodce(y, Fd, [Fs phase], 'psk', M); </pre>
Use a lowpass filter after demodulating but before demapping. num and den are row vectors that give the coefficients, in <i>descending</i> order, of the numerator and denominator of the filter's transfer function.	<pre> z = ddemodce(y, Fd, Fs, 'psk', M, num, den); </pre> <p>(See also “Filter Design Issues” on page 2-62 if you plan to use filters.)</p>
(ASK only) Use a Costas phase-locked loop	<pre> y = dmodce(x, Fd, Fs, 'ask', M); z = ddemodce(y, Fd, Fs, 'ask/costas', M); </pre>
(FSK only) Use noncoherent demodulation	<pre> y = dmodce(x, Fd, Fs, 'fsk', M); z = ddemodce(y, Fd, Fs, 'fsk/noncoherence', M); </pre>

## Selected Bibliography for Modulation

- [1] Jeruchim, Michel C., Philip Balaban, and K. Sam Shanmugan. *Simulation of Communication Systems*. New York: Plenum Press, 1992.
- [2] Proakis, John G. *Digital Communications*, 3rd ed. New York: McGraw-Hill, 1995.
- [3] Sklar, Bernard. *Digital Communications: Fundamentals and Applications*. Englewood Cliffs, N.J.: Prentice-Hall, 1988.

## Special Filters

The Communications Toolbox includes several functions that can help you design and use filters. Other filtering capabilities are in the Signal Processing Toolbox.

### Special Filter Features of the Toolbox

Filtering tasks supported in the Communications Toolbox include:

- Designing a Hilbert transform filter
- Filtering data using a raised cosine filter
- Designing a raised cosine filter

After discussing an implementation issue relating to filters' group delays, this section describes the toolbox functions that accomplish these tasks: `hilb`, `rcosfl`, `rcosine`, and the lower-level functions `rcosfir` and `rcosir`.

For background information about Hilbert filters and raised cosine filters, see the works listed in "Selected Bibliography for Special Filters" on page 2-88. For a demonstration involving raised cosine filters, see `rcosdemo`.

### Noncausality and the Group Delay Parameter

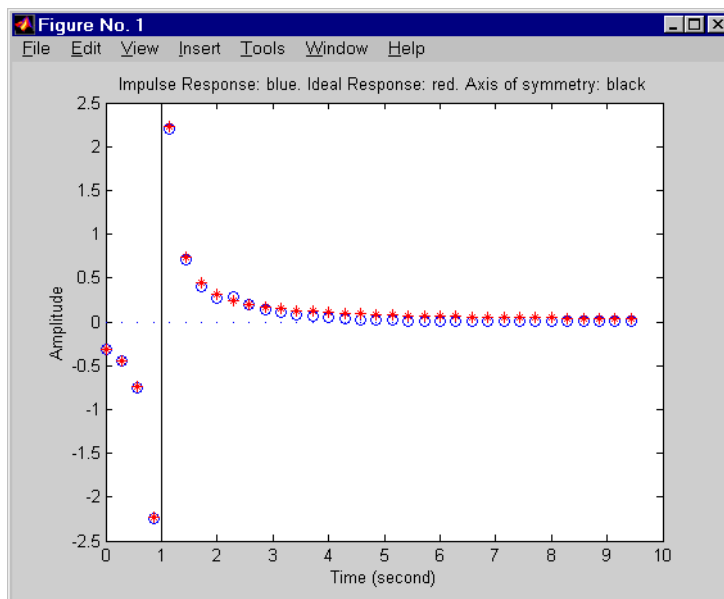
Without propagation delays, both Hilbert filters and raised cosine filters are noncausal. This means that the current output depends on the system's future input. In order to design only *realizable* filters, the `hilb`, `rcosine`, and `rcosfl` functions delay the input signal before producing an output. This delay, known as the filter's *group delay*, is the time between the filter's initial response and its peak response. The group delay is defined as

$$-\frac{d}{d\omega}\theta(\omega)$$

where  $\theta$  is the phase of the filter and  $\omega$  is the frequency in radians. This delay is set so that the impulse response before time zero is negligible and can safely be ignored by the function.

For example, the Hilbert filter whose impulse is shown below uses a group delay of 1 second. Notice in the figure that the impulse response near time 0 is small and that the large impulse response values occur near time 1.





**Figure 2-7: Impulse Response of a Hilbert Filter**

### Example: Compensating for Group Delays When Analyzing Data

Comparing filtered with unfiltered data might be easier if you delay the unfiltered signal by the filter's group delay. For example, suppose you use the code below to filter  $x$  and produce  $y$ .

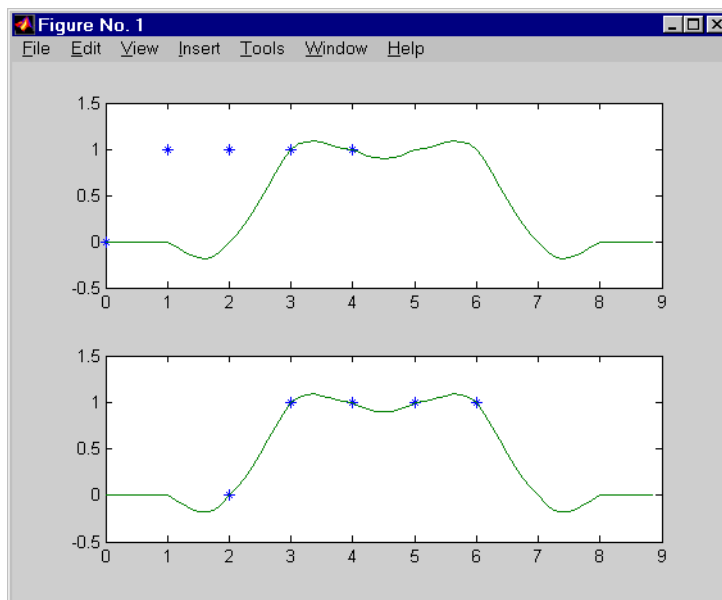
```
tx = 0:4; % Times for data samples
x = [0 1 1 1 1]'; % Binary data samples
% Filter the data and use a delay of 2 seconds.
delay = 2;
[y, ty] = rcosflt(x, 1, 8, 'fir', .3, delay);
```

Here, the elements of  $tx$  and  $ty$  represent the times of each sample of  $x$  and  $y$ , respectively. However,  $y$  is delayed relative to  $x$ , so corresponding elements of  $x$  and  $y$  do not have the same time values. Plotting  $y$  against  $ty$  and  $x$  against  $tx$  is less useful than plotting  $y$  against  $ty$  and  $x$  against a *delayed version* of  $tx$ .

```
% Top plot
subplot(2, 1, 1), plot(tx, x, ' * ', ty, y);
```

```
% Bottom plot delays tx.
subplot(2, 1, 2), plot(tx+delay, x, ' *', ty, y);
```

For another example of compensating for group delay, see the raised-cosine filter demo, `rcosdemo`.



## Designing Hilbert Transform Filters

The `hilb` function designs a Hilbert transform filter and produces either:

- A plot of the filter's impulse response, or
- A quantitative characterization of the filter, using either a transfer function model or a state-space model

### Example with Default Parameters

For example, typing simply

```
hilb
```

plots the impulse response of a fourth-order digital Hilbert transform filter having a 1-second group delay. The sample time is  $2/7$  seconds. In this

particular design, the tolerance index is 0.05. The plot also displays the impulse response of the ideal Hilbert transform filter having a 1-second group delay. The plot is in Figure 2-7, Impulse Response of a Hilbert Filter, on page 2-79.

To compute this filter's transfer function, use the command below.

```
[num, den] = hilbiir

num =

    -0.3183    -0.3041    -0.5160    -1.8453     3.3105

den =

    1.0000    -0.4459    -0.1012    -0.0479    -0.0372
```

Here, the vectors `num` and `den` contain the coefficients of the numerator and denominator, respectively, of the transfer function in ascending order of powers of  $z^{-1}$ .

The commands in this section used the function's default parameters. You can also control the filter design by specifying the sample time, group delay, bandwidth, and tolerance index. The reference entry for `hilbiir` explains these parameters. The group delay is also mentioned above in "Noncausality and the Group Delay Parameter" on page 2-78.

## Filtering with Raised Cosine Filters

The `rcosflt` function applies a raised cosine filter to data. Because `rcosflt` is a versatile function, you can:

- Use `rcosflt` to both design and implement the filter.
- Specify a raised cosine filter and use `rcosflt` only to filter the data.
- Design and implement either raised cosine filters or square-root raised cosine filters.
- Specify the rolloff factor and/or group delay of the filter, if `rcosflt` designs the filter.
- Design and implement either FIR or IIR filters.

This section discusses the use of sampling rates in filtering, and then covers these options. For additional examples, see `rcosdemo`.

## Sampling Rates

The basic `rcosflt` syntax

```
y = rcosflt(x, Fd, Fs...) % Basic syntax
```

assumes by default that you want to apply the filter to a digital signal `x` whose sampling rate is `Fd`. The filter's sampling rate is `Fs`. The ratio of `Fs` to `Fd` must be an integer. By default, the function upsamples the input data by a factor of `Fs/Fd` before filtering. It upsamples by inserting `Fs/Fd - 1` zeros between input data samples. The upsampled data consists of `Fs/Fd` samples per symbol and has sampling rate `Fs`.

An example using this syntax is below. The output sampling rate is four times the input sampling rate.

```
y1 = rcosflt([1; 0; 0], 1, 4, 'fir'); % Upsample by factor of 4/1.
```

**Maintaining the Input Sampling Rate.** You can also override the default upsampling behavior. In this case, the function assumes that the input signal already has sampling rate `Fs` and consists of `Fs/Fd` samples per symbol. You might want to maintain the sampling rate in a receiver's filter if the corresponding transmitter's filter has already upsampled sufficiently.

To maintain the sampling rate, modify the fourth input argument in `rcosflt` to include the string `Fs`. For example, in the first command below, `rcosflt` uses its default upsampling behavior and the output sampling rate is four times the input sampling rate. By contrast, the second command below uses `Fs` in the string argument and thus maintains the sampling rate throughout.

```
y1 = rcosflt([1; 0; 0], 1, 4, 'fir'); % Upsample by factor of 4/1.  
y2 = rcosflt([1; 0; 0], 1, 4, 'fir/Fs'); % Maintain sampling rate.
```

The second command assumes that the sampling rate of the input signal is 4, and that the input signal contains 4/1 samples per symbol.

An example that uses the '`Fs`' option at the receiver is in "Combining Two Square-Root Raised Cosine Filters" on page 2-85.

## Designing Filters Automatically

The simplest syntax of `rcosflt` assumes that the function should both design and implement the raised cosine filter. For example, the command below designs an FIR raised cosine filter and then filters the input vector `[1;0;0]` with it. The second and third input arguments indicate that the function should upsample the data by a factor of 8 (that is,  $8/1$ ) during the filtering process.

```
y = rcosflt([1; 0; 0], 1, 8);
```

**Types of Raised Cosine Filters.** You can have `rcosflt` design other types of raised cosine filters by using a fourth input argument. Variations on the previous example are below.

```
y = rcosflt([1; 0; 0], 1, 8, 'fir'); % Same as original example
y = rcosflt([1; 0; 0], 1, 8, 'fir/sqrt'); % FIR square-root RC filter
y = rcosflt([1; 0; 0], 1, 8, 'iir'); % IIR raised cosine filter
y = rcosflt([1; 0; 0], 1, 8, 'iir/sqrt'); % IIR square-root RC filter
```

## Specifying Filters Using Input Arguments

If you have a transfer function for a raised cosine filter, then you can provide it as an input to `rcosflt` so that `rcosflt` does not design its own filter. This is useful if you want to use `rcosine` to design the filter once and then use the filter many times. For example, the `rcosflt` command below uses the `'filter'` flag to indicate that transfer function is an input argument. The input `num` is a vector that represents the FIR transfer function by listing its coefficients.

```
num = rcosine(1, 8); y = rcosflt([1; 0; 0], 1, 8, 'filter', num);
```

This syntax for `rcosflt` works whether `num` represents the transfer function for a square-root raised cosine FIR filter or an ordinary raised cosine FIR filter. For example, the code below uses a square-root raised cosine FIR filter. Only the definition of `num` is different.

```
num = rcosine(1, 8, 'sqrt'); y = rcosflt([1; 0; 0], 1, 8, 'filter', num);
```

You can also use a raised cosine IIR filter. To do this, modify the fourth input argument of the `rcosflt` command above so that it contains the string `'iir'` and provide a denominator argument. An example is below.

```
delay = 8;
[num, den] = rcosine(1, 8, 'iir', .5, delay);
y = rcosflt([1; 0; 0], 1, 8, 'iir/filter', num, den, delay);
```

### Controlling the Rolloff Factor

If `rcosflt` designs the filter automatically, then you can control the rolloff factor of the filter, as described below. If you specify your own filter, then `rcosflt` does not need to know its rolloff factor.

The rolloff factor determines the excess bandwidth of the filter. For example, a rolloff factor of .5 means that the bandwidth of the filter is 1.5 times the input sampling frequency,  $F_d$ . This also means that the transition band of the filter extends from  $.5 * F_d$  to  $1.5 * F_d$ .

The default rolloff factor is .5, but if you want to use a value of .2, then you can use a command such as the one below. Typical values for the rolloff factor are between .2 and .5.

```
y = rcosflt([1;0;0], 1, 8, 'fir', .2); % Rolloff factor is .2.
```

### Controlling the Group Delay

If `rcosflt` designs the filter automatically, then you can control the group delay of the filter, as described below. If you specify your own FIR filter, then `rcosflt` does not need to know its group delay.

The filter's group delay is the time between the filter's initial response and its peak response. The default group delay in the implementation is three input samples. To specify a different value, measure it in input symbol periods and provide it as the sixth input argument. For example, the command below specifies a group delay of six input samples, which is equivalent to  $6*8/1$  output samples.

```
y = rcosflt([1;0;0], 1, 8, 'fir', .2, 6); % Delay is 6 input samples.
```

The group delay influences the size of the output, as well as the order of the filter if `rcosflt` designs the filter automatically. See the reference page for `rcosflt` for details that relate to the syntax you want to use.

**Example: Raised Cosine Filter Delays.** The code below filters a signal using two different group delays. A larger delay results in a smaller error in the frequency response of the filter. The plot shows how the two filtered signals differ, and the output `pt` indicates that the first peak occurs at different times for the two filtered signals.

```
[y, t] = rcosflt(ones(10, 1), 1, 8, 'fir', .5, 6); % Delay = 6 samples  
[y1, t1] = rcosflt(ones(10, 1), 1, 8, 'fir', .5, 8); % Delay = 8 samples
```

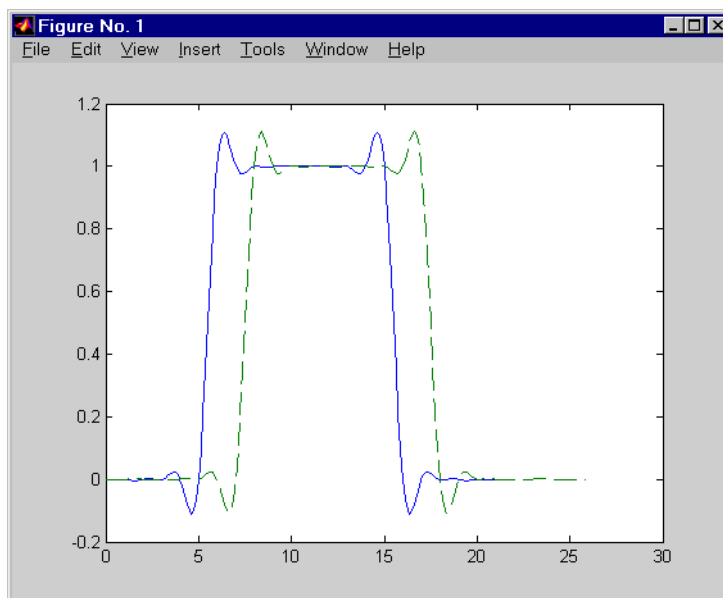
```

plot(t, y, t1, y1, '--') % Two curves indicate the different delays.
peak = t(find(y == max(y))); % Times where first curve peaks
peak1 = t1(find(y1 == max(y1))); % Times where second curve peaks
pt = [min(peak), min(peak1)] % First peak time for both curves

```

```
pt =
```

```
14.6250    16.6250
```



**Figure 2-8: Delays of Three Samples (Dashed) and Five Samples (Solid)**

If  $F_s/F_d$  is at least 4, then a group delay value of at least 8 works well in many cases. In the examples of this section,  $F_s/F_d$  is 8.

### Combining Two Square-Root Raised Cosine Filters

If you want to split the filtering equally between the transmitter's filter and the receiver's filter, then you can use a pair of square-root raised cosine filters. In theory, the combination of two square-root raised cosine filters is equivalent to a single normal raised cosine filter. However, the limited impulse response of

practical square-root raised cosine filters causes a slight difference between the response of two successive square-root raised cosine filters and the response of one raised cosine filter.

Using `rcosine` and `rcosflt` to Implement Square-Root Raised Cosine Filters. One way to implement the pair of square-root raised cosine filters is to follow these steps:

- 1 Use `rcosine` with the `'sqrt'` flag to design a square-root raised cosine filter.
- 2 Use `rcosflt` in the transmitter section of code to upsample and filter the data.
- 3 Use `rcosflt` in the receiver section of code to filter the received data *without* upsampling it. Use the `'Fs'` flag to avoid upsampling.

An example of this approach is below. Notice that the syntaxes for `rcosflt` use the `'filter'` flag to indicate that you are providing the filter's transfer function as an input.

```
% First approach
x = randint(100, 1, 2, 1234); % Data
num = rcosine(1, 8, 'sqrt'); % Transfer function of filter
y = rcosflt(x, 1, 8, 'filter', num); % Filter the data.
z = rcosflt(y, 1, 8, 'Fs/filter', num); % Filter the received data
% but do not upsample it.
```

Using `rcosflt` Alone. Another way to implement the pair of square-root raised cosine filters is to have `rcosflt` both design and use the square-root raised cosine filter. This approach avoids using `rcosine`. The corresponding example code is below. Notice that the syntaxes for `rcosflt` use the `'sqrt'` flag to indicate that you want it to design a square-root raised cosine filter.

```
% Second approach
x = randint(100, 1, 2, 1234); % Data (again)
y1 = rcosflt(x, 1, 8, 'sqrt'); % Design and use a filter.
z1 = rcosflt(y1, 1, 8, 'sqrt/Fs'); % Design and use a filter
% but do not upsample the data.
```

Because these two approaches are equivalent, `y` is the same as `y1` and `z` is the same as `z1`.



## Designing Raised Cosine Filters

The `rcosine` function designs (but does not apply) filters of these types:

- Finite impulse response (FIR) raised cosine filter
- Infinite impulse response (IIR) raised cosine filter
- FIR square-root raised cosine filter
- IIR square-root raised cosine filter

The function returns the transfer function as output. To learn about applying raised cosine filters, see “Filtering with Raised Cosine Filters” on page 2-81.

### Sampling Rates

The `rcosine` function assumes that you want to apply the filter to a digital signal whose sampling rate is  $F_d$ . The function also requires you to provide the filter’s sampling rate,  $F_s$ . The ratio of  $F_s$  to  $F_d$  must be an integer.

### Example Designing a Square-Root Raised Cosine Filter

For example, the command below designs a square-root raised cosine FIR filter with a sampling rate of 2, for use with a digital signal whose sampling rate is 1.

```
num = rcosine(1, 2, 'fir/sqrt')

num =

Columns 1 through 7

    0.0021   -0.0106    0.0300   -0.0531   -0.0750    0.4092    0.8037

Columns 8 through 13

    0.4092   -0.0750   -0.0531    0.0300   -0.0106    0.0021
```

Here, the vector `num` contains the coefficients of the filter, in ascending order of powers of  $z^{-1}$ .

### Other Options in Filter Design

You can also control the filter design by specifying the rolloff factor, group delay, and (for IIR filters) tolerance index explicitly, instead of having `rcosine` use its default values. The reference entry for `rcosine` explains these

parameters. The group delay is also mentioned above in “Noncausality and the Group Delay Parameter” on page 2-78.

### **Selected Bibliography for Special Filters**

[1] Korn, Israel. *Digital Communications*. New York: Van Nostrand Reinhold, 1985.

[2] Oppenheim, Alan V. and Ronald W. Schaffer. *Discrete-Time Signal Processing*. Englewood Cliffs, N.J.: Prentice Hall, 1989.

## Galois Field Computations

A *Galois field* is an algebraic field that has a finite number of elements. The number of elements is always of the form  $p^m$ , where  $p$  is a prime number and  $m$  is a positive integer. Galois fields are used in error-control coding.

### Galois Field Features of the Toolbox

The Communications Toolbox provides functions for manipulating elements of Galois fields, working with polynomials over Galois fields, and performing other tasks related to Galois fields. This section discusses these topics:

- “Galois Field Terminology” on page 2-89
- “Representing Elements of Galois Fields” on page 2-90
- “Default Primitive Polynomials” on page 2-93
- “Converting and Simplifying Element Formats” on page 2-94
- “Arithmetic in Galois Fields” on page 2-97
- “Polynomials over Prime Fields” on page 2-99

For background information about Galois fields or their use in error-control coding, see the works listed in “Selected Bibliography for Galois Fields” on page 2-103.

### Galois Field Terminology

Throughout this section,  $p$  is a prime number and  $m$  is a positive integer.

Also, this document uses a few terms that are not used consistently in the literature. The definitions adopted here appear in van Lint [4].

- A *primitive element* of  $\text{GF}(p^m)$  is a cyclic generator of the group of nonzero elements of  $\text{GF}(p^m)$ . This means that every nonzero element of the field can be expressed as the primitive element raised to some integer power. Primitive elements are called  $\alpha$  throughout this section.
- A *primitive polynomial* for  $\text{GF}(p^m)$  is the minimal polynomial of some primitive element of  $\text{GF}(p^m)$ . As a consequence, it has degree  $m$  and is irreducible.

## Representing Elements of Galois Fields

This section discusses how to represent Galois field elements using this toolbox's exponential format and polynomial format. It also describes a way to list all elements of the Galois field, because some functions use such a list as an input argument. Finally, it discusses the nonuniqueness of representations of Galois field elements.

The elements of  $\text{GF}(p)$  can be represented using the integers from 0 to  $p-1$ .

When  $m$  is at least two,  $\text{GF}(p^m)$  is called an extension field. Integers alone cannot represent the elements of  $\text{GF}(p^m)$  in a straightforward way. MATLAB uses two main conventions for representing elements of  $\text{GF}(p^m)$ : the exponential format and the polynomial format.

**Note** Both the exponential format and the polynomial format are relative to your choice of a particular primitive element  $\alpha$  of  $\text{GF}(p^m)$ .

### Exponential Format

This format uses the property that every nonzero element of  $\text{GF}(p^m)$  can be expressed as  $\alpha^c$  for some integer  $c$  between 0 and  $p^m-2$ . Higher exponents are not needed, since the theory of Galois fields implies that every nonzero element of  $\text{GF}(p^m)$  satisfies the equation  $x^{q-1} = 1$  where  $q = p^m$ .

MATLAB's use of the exponential format is shown in the table below.

Table 2-10: Exponential Format in MATLAB

Element of $\text{GF}(p^m)$	MATLAB Representation of the Element
0	-Inf
$\alpha^0 = 1$	0
$\alpha^1$	1
...	...
$\alpha^{q-2}$ where $q = p^m$	$q-2$

Although `-Inf` is the standard exponential representation of the zero element, all negative integers are equivalent to `-Inf` when used as *input* arguments in exponential format. This equivalence can be useful; for example, see the concise line of code at the end of the section “Default Primitive Polynomials” on page 2-93.

---

**Note** The equivalence of all negative integers and `-Inf` as exponential formats means that, for example, `-1` does *not* represent  $\alpha^{-1}$ , the multiplicative inverse of  $\alpha$ . Instead, `-1` represents the zero element of the field.

---

### Polynomial Format

The polynomial format uses the property that every element of  $\text{GF}(p^m)$  can be expressed as a polynomial in  $\alpha$  with exponents between 0 and  $m-1$ , and coefficients in  $\text{GF}(p)$ . In the polynomial format, the element

$$A(1) + A(2) \alpha + A(3) \alpha^2 + \dots + A(m) \alpha^{m-1}$$

is represented in MATLAB by the vector

$$[A(1) \ A(2) \ A(3) \ \dots \ A(m)]$$

---

**Note** The Galois field functions in this toolbox represent a polynomial as a vector that lists the coefficients in order of *ascending* powers of the variable. This is the opposite of the order that other MATLAB functions use.

---

### List of All Elements of a Galois Field

Some Galois field functions in this toolbox require an argument that lists all elements of an extension field  $\text{GF}(p^m)$ . This is again relative to a particular primitive element  $\alpha$  of  $\text{GF}(p^m)$ . The proper format for the list of elements is that of a matrix having  $p^m$  rows, one for each element of the field. The matrix has  $m$  columns, one for each coefficient of a power of  $\alpha$  in the polynomial format shown in “Polynomial Format” above. The first row contains only zeros because it corresponds to the zero element in  $\text{GF}(p^m)$ . If  $k$  is between 2 and  $p^m$ , then the  $k$ th row specifies the polynomial format of the element  $\alpha^{k-2}$ .

The minimal polynomial of  $\alpha$  aids in the computation of this matrix, since it tells how to express  $\alpha^m$  in terms of lower powers of  $\alpha$ . For example, the table below lists the elements of  $\text{GF}(3^2)$ , where  $\alpha$  is a root of the primitive polynomial  $2 + 2x + x^2$ . This polynomial allows repeated use of the substitution

$$\alpha^2 = -2 - 2\alpha = 1 + \alpha$$

when performing the computations in the middle column of the table.

Table 2-11: Elements of  $\text{GF}(9)$

Exponential Format	Polynomial Format	Row of MATLAB Matrix of Elements
$\alpha^{-\text{Inf}}$	0	0 0
$\alpha^0$	1	1 0
$\alpha^1$	$\alpha$	0 1
$\alpha^2$	$1+\alpha$	1 1
$\alpha^3$	$\alpha + \alpha^2 = \alpha + 1 + \alpha = 1 + 2\alpha$	1 2
$\alpha^4$	$\alpha + 2\alpha^2 = \alpha + 2 + 2\alpha = 2$	2 0
$\alpha^5$	$2\alpha$	0 2
$\alpha^6$	$2\alpha^2 = 2 + 2\alpha$	2 2
$\alpha^7$	$2\alpha + 2\alpha^2 = 2\alpha + 2 + 2\alpha = 2 + \alpha$	2 1

An automatic way to generate the matrix whose rows are in the third column of the table above is to use the code below.

```
p = 3; m = 2;
% Use the primitive polynomial 2 + 2x + x^2 for GF(9).
primpoly = [2 2 1];
field = gftuple([-1: p^m-2]', primpoly, p);
```

The `gftuple` function is discussed in more detail in “Converting and Simplifying Element Formats” on page 2-94.

## Nonuniqueness of Representations

A given field has more than one primitive element. If two primitive elements have different minimal polynomials, then the corresponding matrices of elements will have their rows in a different order. If the two primitive elements share the same minimal polynomial, then the matrix of elements of the field is the same.

---

**Note** You may use whatever primitive element you want, as long as you understand how the inputs and outputs of Galois field functions depend on the choice of *some* primitive polynomial. It is usually best to use the same primitive polynomial throughout a given script or function.

---

Other ways in which representations of elements are not unique arise from the equations that Galois field elements satisfy. For example, an exponential format of 8 in GF(9) is really the same as an exponential format of 0, since  $\alpha^8 = 1 = \alpha^0$  in GF(9). As another example, the substitution mentioned just before Table 2-11, Elements of GF(9), shows that the polynomial format [0 0 1] is really the same as the polynomial format [1 1].

## Default Primitive Polynomials

This toolbox provides a *default* primitive polynomial for each extension field. You can retrieve this polynomial using the `gfpri mdf` function. The command

```
primpoly = gfpri mdf(m, p); % If m and p are already defined
```

produces the standard row-vector representation of the default minimal polynomial for GF( $p^m$ ).

For example, the command below shows that the default primitive polynomial for GF(9) is  $2 + x + x^2$ , *not* the polynomial used in the section, “List of All Elements of a Galois Field” on page 2-91.

```
gfpri mdf(2, 3)
```

```
ans =
```

```
2      1      1
```

To generate a list of elements of  $GF(p^m)$  using the default primitive polynomial, use the command

```
field = gftuple([-1:p^m-2]', m, p);
```

Converting and Simplifying Element Formats

This section describes how to convert between the exponential and polynomial formats for Galois field elements, as well as how to simplify a given representation.

Converting to Simplest Polynomial Format

The `gftuple` function produces the simplest polynomial representation of an element of  $GF(p^m)$ , given either an exponential representation or a polynomial representation of that element. This can be useful for generating the list of elements of  $GF(p^m)$  that other functions require.

The simplest use of `gftuple` requires two arguments: one representing an element of  $GF(p^m)$  and the other indicating the primitive polynomial that MATLAB should use when computing the output. An optional third argument is the prime  $p$ ; if it is omitted, then the default is 2. The table below indicates how `gftuple` behaves when given the first two arguments in various formats.

Table 2-12: Behavior of `gftuple` Depending on Format of Inputs

How to Specify Element	How to Indicate Primitive Polynomial	What <code>gftuple</code> Produces
Exponential format; $c$ = any integer	Integer $m > 1$	Polynomial format of $\alpha^c$ , where $\alpha$ is a root of the <i>default</i> primitive polynomial for $GF(p^m)$
Example: <code>tp = gftuple(6, 2, 3); % c = 6 here</code>		
Exponential format; $c$ = any integer	Vector of coefficients of primitive polynomial	Polynomial format of $\alpha^c$ , where $\alpha$ is a root of the <i>given</i> primitive polynomial
Example: <code>polynomial = gfpri MDF(2, 3); tp = gftuple(6, polynomial, 3); % c = 6 here</code>		



Table 2-12: Behavior of `gftuple` Depending on Format of Inputs (Continued)

How to Specify Element	How to Indicate Primitive Polynomial	What <code>gftuple</code> Produces
Polynomial format of any degree	Integer $m > 1$	Polynomial format of degree $< m$ , using <i>default</i> primitive polynomial for $\text{GF}(p^m)$ to simplify
Example: <code>tp = gftuple([0 0 0 0 0 0 1], 2, 3);</code>		
Polynomial format of any degree	Vector of coefficients of primitive polynomial	Polynomial format of degree $< m$ , using the <i>given</i> primitive polynomial for $\text{GF}(p^m)$ to simplify
Example: <code>polynomial = gfprmdf(2, 3); tp = gftuple([0 0 0 0 0 0 1], polynomial, 3);</code>		

The four examples that appear in the table above all produce the same vector `tp = [2, 1]`, but their different inputs to `gftuple` correspond to the lines of the table. Each example expresses the fact that

$$\alpha^6 = 2 + \alpha$$

where  $\alpha$  is a root of the (default) primitive polynomial  $2 + x + x^2$  for  $\text{GF}(3^2)$ .

**Example.** This example shows how `gfconv` and `gftuple` combine to multiply two polynomial-format elements of  $\text{GF}(3^4)$ . Initially, `gfconv` multiplies the two polynomials, treating the primitive element as if it were a variable. This produces a high-order polynomial, which `gftuple` simplifies using the polynomial equation that the primitive element satisfies. The final result is the simplest polynomial format of the product.

```
p = 3; m = 4;
a = [1 2 0 1]; b = [2 2 1 2];
notsimple = gfconv(a, b, p) % a times b, using high powers of alpha

notsimple =

      2      0      2      0      0      1      2

simple = gftuple(notsimple, m, p) % Highest exponent of alpha is m-1
```

```
simple =  
  
2      1      0      1
```

### Example: Generating a List of Galois Field Elements

This example applies the conversion functionality to the task of generating a matrix that lists all elements of a Galois field. A matrix that lists all field elements is an input argument in functions such as `gfadd` and `gfmul`. The variables `field1` and `field2` below have the format that such functions expect.

```
p = 5; % Or any prime number  
m = 4; % Or any positive integer  
field1 = gftuple([-1:p^m-2]', m, p);  
  
primpoly = gfprimdf(m, p); % Or any primitive polynomial  
% for GF(p^m)  
field2 = gftuple([-1:p^m-2]', primpoly, p);
```

### Converting to Simplest Exponential Format

The same function `gftuple` also produces the simplest exponential representation of an element of  $GF(p^m)$ , given either an exponential representation or a polynomial representation of that element. To retrieve this output, use the syntax

```
[polyformat, expformat] = gftuple(...)
```

The input format and the output `polyformat` are as in Table 2-12, Behavior of `gftuple` Depending on Format of Inputs. In addition, the variable `expformat` contains the simplest exponential format of the element represented in `polyformat`. It is *simplest* in the sense that the exponent is either  $-\infty$  or a number between 0 and  $p^m-2$ .

To recover the exponential format of the element  $2 + \alpha$  that the previous section considered, use the commands below. In this case, `polyformat` contains redundant information, while `expformat` contains the desired result.

```
[polyformat, expformat] = gftuple([2 1], 2, 3)  
  
polyformat =  
  
2      1
```

```
expformat =
```

```
6
```

This output appears at first to contradict the information in Table 2-11, Elements of  $\text{GF}(9)$ , but in fact it does not. The table uses a different primitive element; two plus that primitive element has the polynomial and exponential formats shown below. The output below reflects the information in the bottom line of the table.

```
primpoly = [2 2 1];
[polyformat, expformat] = gftuple([2 1], primpoly, 3)
```

```
polyformat =
```

```
2      1
```

```
expformat =
```

```
7
```

## Arithmetic in Galois Fields

You can add, subtract, multiply, and divide elements of Galois fields using the functions `gfadd`, `gfsub`, `gfmul`, and `gfdi v`, respectively. Each of these functions has a mode for prime fields and a mode for extension fields.

### Arithmetic in Prime Fields

Arithmetic in  $\text{GF}(p)$  is the same as arithmetic modulo  $p$ . The functions `gfadd`, `gfmul`, `gfsub`, and `gfdi v` accept two arguments that represent elements of  $\text{GF}(p)$  as integers between 0 and  $p-1$ . An optional third argument specifies  $p$ ; if it does not appear, then the computations are performed in  $\text{GF}(2)$ .

**Example: Addition Table for  $\text{GF}(5)$ .** The code below constructs an addition table for  $\text{GF}(5)$ . If  $a$  and  $b$  are between 0 and 4, then the element `gfp_add(a+1, b+1)` represents the sum  $a+b$  in  $\text{GF}(5)$ . For example, `gfp_add(3, 5) = 1` because  $2+4$  is 1 modulo 5.

```
p = 5;
row = 0: p-1;
```

```
table = ones(p, 1)*row;
gfp_add = gfadd(table, table', p)
```

```
gfp_add =
```

0	1	2	3	4
1	2	3	4	0
2	3	4	0	1
3	4	0	1	2
4	0	1	2	3

Other values of  $p$  produce tables for different prime fields  $GF(p)$ . Replacing `gfadd` by `gfmul`, `gfsub`, or `gfdiv` produces a table for the corresponding arithmetic operation in  $GF(p)$ .

### Arithmetic in Extension Fields

The same arithmetic functions can add elements of  $GF(p^m)$  when  $m > 1$ , but the format of the arguments is more complicated than in the case above. In general, arithmetic in extension fields is more complicated than arithmetic in prime fields; see the works listed in “Selected Bibliography for Galois Fields” on page 2-103 for details about how the arithmetic operations work.

When working in extension fields, the functions `gfadd`, `gfmul`, `gfsub`, and `gfdiv` use the first two arguments to represent elements of  $GF(p^m)$  in exponential format. The third argument, which is required, lists all elements of  $GF(p^m)$  as described in the section, “List of All Elements of a Galois Field” on page 2-91. The result is in exponential format.

**Example: Addition Table for  $GF(9)$ .** The code below constructs an addition table for  $GF(3^2)$ , using exponential formats relative to a root of the default primitive polynomial for  $GF(9)$ . If  $a$  and  $b$  are between -1 and 7, then the element `gfp_add(a+2, b+2)` represents the sum of  $\alpha^a$  and  $\alpha^b$  in  $GF(9)$ . For example, `gfp_add(4, 6) = 5` because

$$\alpha^2 + \alpha^4 = \alpha^5$$

Using the fourth and sixth rows of the matrix `field`, you can verify that

$$\alpha^2 + \alpha^4 = (1 + 2\alpha) + (2 + 0\alpha) = 3 + 2\alpha = 0 + 2\alpha = \alpha^5 \text{ modulo } 3.$$

```
p = 3; m = 2; % Work in GF(3^2).
field = gftuple([-1:p^m-2]', m, p); % Construct list of elements.
```

```
row = -1:p^m-2;
table = ones(p^m, 1)*row;
gfpm_add = gfadd(table, table', field)
```

```
gfpm_add =
```

- Inf	0	1	2	3	4	5	6	7
0	4	7	3	5	- Inf	2	1	6
1	7	5	0	4	6	- Inf	3	2
2	3	0	6	1	5	7	- Inf	4
3	5	4	1	7	2	6	0	- Inf
4	- Inf	6	5	2	0	3	7	1
5	2	- Inf	7	6	3	1	4	0
6	1	3	- Inf	0	7	4	2	5
7	6	2	4	- Inf	1	0	5	3

---

**Note** If you used a different primitive polynomial, then the tables would look different. This makes sense because the ordering of the rows and columns of the tables was based on that particular choice of primitive polynomial and not on any natural ordering of the elements of  $GF(9)$ .

---

Other values of  $p$  and  $m$  produce tables for different prime fields  $GF(p^m)$ . Replacing `gfadd` by `gfmul`, `gfsub`, or `gfdi v` produces a table for the corresponding arithmetic operation in  $GF(p^m)$ .

## Polynomials over Prime Fields

A polynomial over  $GF(p)$  is a polynomial whose coefficients are elements of  $GF(p)$ . The Communications Toolbox provides functions for:

- Changing polynomials in cosmetic ways
- Performing polynomial arithmetic
- Characterizing polynomials as primitive or irreducible
- Finding roots of polynomials in a Galois field

---

**Note** The Galois field functions in this toolbox represent a polynomial as a vector that lists the coefficients in order of *ascending* powers of the variable. This is the opposite of the order that other MATLAB functions use.

---

### Cosmetic Changes of Polynomials

To display the traditionally formatted polynomial that corresponds to a row vector containing coefficients, use `gfpretty`. To truncate a polynomial by removing all zero-coefficient terms that have exponents *higher* than the degree of the polynomial, use `gftrunc`. For example,

```
polynom = gftrunc([1 20 394 10 0 0 29 3 0 0])

polynom =

    1    20   394    10     0     0    29     3

gfpretty(polynom)
```

$$1 + 20 X + 394 X^2 + 10 X^3 + 29 X^6 + 3 X^7$$


---

**Note** If you do not use a fixed-width font, then the spacing in the display might not look correct.

---

### Polynomial Arithmetic

The functions `gfadd` and `gfsub` add and subtract, respectively, polynomials over  $\text{GF}(p)$ . The `gfconv` function multiplies polynomials over  $\text{GF}(p)$ . The `gfdeconv` function divides polynomials in  $\text{GF}(p)$ , producing a quotient polynomial and a remainder polynomial. For example, the commands below show that  $2 + x + x^2$  times  $1 + x$  over the field  $\text{GF}(3)$  is  $2 + 2x^2 + x^3$ .

```
a = gfconv([2 1 1], [1 1], 3)
```

```

a =

      2      0      2      1

[quot, remd] = gfdeconv(a, [2 1 1], 3)

quot =

      1      1

remd =

      0

```

The previously discussed functions `gfadd` and `gfsub` add and subtract, respectively, polynomials. Because it uses a vector of coefficients to represent a polynomial, MATLAB does not distinguish between adding two polynomials and adding two row vectors elementwise.

### Characterization of Polynomials

Given a polynomial over  $\text{GF}(p)$ , the `gfprimck` function determines whether it is irreducible and/or primitive. By definition, if it is primitive then it is irreducible; however, the reverse is not necessarily true.

Given an element of  $\text{GF}(p^m)$ , the `gfminpol` function computes its minimal polynomial over  $\text{GF}(p)$ .

For example, the code below reflects the irreducibility of all minimal polynomials. However, the minimal polynomial of a nonprimitive element is not a primitive polynomial.

```

p = 2; m = 4;
% Use default primitive polynomial here.

primpoly = gfminpol(1, m, p);
ckprim = gfprimck(primpoly, p);
% ckprim = 1, since primpoly represents a primitive polynomial.

notprimpoly = gfminpol(3, m, p);
cknotprim = gfprimck(notprimpoly, p);
% cknotprim = 0 (irreducible but not primitive)

```

```
% since alpha^3 is not a primitive element when p = 2.

ckreducible = gfprimck([0 1 1], p);
% ckreducible = -1 since the polynomial is reducible.
```

Roots of Polynomials

Given a polynomial over  $GF(p)$ , the `gfroots` function finds the roots of the polynomial in a suitable extension field  $GF(p^m)$ . If  $p$  is not specified, then the default is 2. If  $m$  is not specified, then the default is the degree of the polynomial. There are two ways to tell MATLAB the degree  $m$  of the extension field  $GF(p^m)$ , as shown in the table below.

Table 2-13: Formats for Second Argument of `gfroots`

Second Argument	Represents
A positive integer	$m$ as in $GF(p^m)$ . MATLAB uses the default primitive polynomial in its computations.
A row vector	a primitive polynomial for $GF(p^m)$ . Here $m$ is the degree of this primitive polynomial.

**Example: Roots of a Polynomial in  $GF(9)$ .** The code below finds roots of the polynomial  $1 + x^2 + x^3$  in  $GF(9)$  and then checks that they are indeed roots. The exponential format of elements of  $GF(9)$  is used throughout.

```
p = 3; m = 2;
field = gftuple([-1;p^m-2]', m, p); % List of all elements of GF(9)
% Use default primitive polynomial here.
polynomial = [1 0 1 1]; % 1 + x^2 + x^3
roots = gfroots(polynomial, m, p) % Find roots in exponential format
% Check that each one is actually a root.
for ii = 1:3
    root = roots(ii);
    rootsquared = gfmul(root, root, field);
    rootcubed = gfmul(root, rootsquared, field);
    answer(ii) = ...
        gfadd(gfadd(0, rootsquared, field), rootcubed, field);
    % Recall that 1 is really alpha to the zero power.
    % If answer = -Inf, then the variable root represents
    % a root of the polynomial.
```



```
end
answer
```

The output shows that  $\alpha^0$  (which equals 1),  $\alpha^5$ , and  $\alpha^7$  are roots.

```
roots =
```

```
    0
    5
    7
```

```
answer =
```

```
-Inf -Inf -Inf
```

See the reference page for `gfroots` to see how `gfroots` can also provide you with the polynomial formats of the roots and the list of all elements of the field.

## Other Galois Field Functions

See the reference pages for information about these other Galois field functions in the Communications Toolbox:

- `gfcosets`, which produces cyclotomic cosets
- `gffilter`, which filters data using  $\text{GF}(p)$  polynomials
- `gflineq`, which solves a linear matrix equation over  $\text{GF}(p)$
- `gfpriofd`, which finds primitive polynomials
- `gfrank`, which computes the rank of a matrix over  $\text{GF}(p)$
- `gfepcov`, which converts one  $\text{GF}(2)$  polynomial representation to another

## Selected Bibliography for Galois Fields

- [1] Blahut, Richard E. *Theory and Practice of Error Control Codes*. Reading, Mass.: Addison-Wesley, 1983, p.105.
- [2] Lang, Serge. *Algebra*. Third Edition. Reading, Mass.: Addison-Wesley, 1993.
- [3] Lin, Shu and Daniel J. Costello, Jr. *Error Control Coding: Fundamentals and Applications*. Englewood Cliffs, N.J.: Prentice-Hall, 1983.

[4] van Lint, J. H. *Introduction to Coding Theory*. New York: Springer-Verlag, 1982.

## Reference

---

This chapter contains detailed descriptions of all Communications Toolbox functions. To access the descriptions, use the links in the second column of the table below.

Organization of Functions	Section
By category	“Functions by Category”
Alphabetical	“Alphabetical List of Functions”

## Functions by Category

**Table 3-1: Signal Sources**

Function	Purpose
randerr	Generate bit error patterns
randint	Generate matrix of uniformly distributed random integers
randsrc	Generate random matrix using prescribed alphabet
wgn	Generate white Gaussian noise

**Table 3-2: Signal Analysis Functions**

Function	Purpose
biterr	Compute number of bit errors and bit error rate
eyediagram	Generate an eye diagram
scatterplot	Generate a scatter plot
symerr	Compute number of symbol errors and symbol error rate

**Table 3-3: Source Coding**

Function	Purpose
compand	Source code mu-law or A-law compressor or expander
dpcmdeco	Decode using differential pulse code modulation
dpcmenco	Encode using differential pulse code modulation
dpcmopt	Optimize differential pulse code modulation parameters

Table 3-3: Source Coding (Continued)

Function	Purpose
lloyd s	Optimize quantization parameters using the Lloyd algorithm
quantiz	Produce a quantization index and a quantized output value

Table 3-4: Error-Control Coding

Function	Purpose
bchpol y	Produce parameters or generator polynomial for binary BCH code
convenc	Convolutionally encode binary data
cycl gen	Produce parity-check and generator matrices for cyclic code
cycl pol y	Produce generator polynomials for a cyclic code
decode	Block decoder
encode	Block encoder
gen2par	Convert between parity-check and generator matrices
gfwei ght	Calculate the minimum distance of a linear block code
hammgen	Produce parity-check and generator matrices for Hamming code
rsdec of	Decode an ASCII file that was encoded using Reed-Solomon code
rsenc of	Encode an ASCII file using Reed-Solomon code
rspol y	Produce Reed-Solomon code generator polynomial

**Table 3-4: Error-Control Coding (Continued)**

Function	Purpose
syndtbl e	Produce syndrome decoding table
vitdec	Convolutionally decode binary data using the Viterbi algorithm

**Table 3-5: Lower-Level Functions for Error-Control Coding**

Function	Purpose
bchdeco	BCH decoder
bchenco	BCH encoder
rsdeco	Reed-Solomon decoder
rsdecode	Reed-Solomon decoding using the exponential format
rsenco	Reed-Solomon encoder
rsencode	Reed-Solomon encoding using the exponential format

**Table 3-6: Modulation and Demodulation**

Function	Purpose
ademod	Analog passband demodulator
ademodce	Analog baseband demodulator
amod	Analog passband modulator
amodce	Analog baseband modulator
apkconst	Plot a combined circular ASK-PSK signal constellation
ddemod	Digital passband demodulator

Table 3-6: Modulation and Demodulation (Continued)

Function	Purpose
ddemodce	Digital baseband demodulator
demodmap	Demap a digital message from a demodulated signal
dmod	Digital passband modulator
dmodce	Digital baseband modulator
modmap	Map a digital signal to an analog signal
qaskdeco	Demap a message from a QASK square signal constellation
qaskenco	Map a message to a QASK square signal constellation

Table 3-7: Special Filters

Function	Purpose
hank2sys	Convert a Hankel matrix to a linear system model
hilbiir	Design a Hilbert transform IIR filter
rcosflt	Filter the input signal using a raised cosine filter
rcosine	Design a raised cosine filter

Table 3-8: Lower-Level Functions for Special Filters

Function	Purpose
rcosfir	Design a raised cosine FIR filter
rcosiir	Design a raised cosine IIR filter



**Table 3-9: Channel Functions**

Function	Purpose
awgn	Add white Gaussian noise to a signal

**Table 3-10: Galois Field Computation**

Function	Purpose
gfadd	Add polynomials over a Galois field
gfconv	Multiply polynomials over a Galois field
gfcosets	Produce cyclotomic cosets for a Galois field
gfdeconv	Divide polynomials over a Galois field
gfdiv	Divide elements of a Galois field
gffilter	Filter data using polynomials over a prime Galois field
gflin	Solve the linear equation $Ax = b$ over a prime Galois field
gfminpol	Find the minimal polynomial of an element of a Galois field
gfmul	Multiply elements of a Galois field
gfplus	Add elements of a Galois field of characteristic two
gfpretty	Display a polynomial in traditional format
gfpri mck	Check whether a polynomial over a Galois field is primitive
gfpri mdf	Provide default primitive polynomials for a Galois field
gfpri mfd	Find primitive polynomials for a Galois field
gfrank	Compute the rank of a matrix over a Galois field
gfrepconv	Convert one GF(2) polynomial representation to another

**Table 3-10: Galois Field Computation (Continued)**

Function	Purpose
gfroots	Find the roots of a polynomial over a prime Galois field
gfsub	Subtract polynomials over a Galois field
gftrunc	Minimize the length of a polynomial representation
gftuple	Simplify or convert the format of elements of a Galois field

**Table 3-11: Utilities**

Function	Purpose
bi2de	Convert binary vectors to decimal numbers
de2bi	Convert decimal numbers to binary vectors
erf	Error function
erfc	Complementary error function
istrellis	Check if the input is a valid trellis structure
marcumq	Generalized Marcum Q function
oct2dec	Convert octal numbers to decimal numbers
poly2trellis	Convert convolutional code polynomials to trellis description
vec2mat	Convert a vector into a matrix

# Alphabetical List of Functions

ademod	3-12
ademodce	3-16
amod	3-20
amodce	3-25
apkconst	3-28
awgn	3-32
bchdeco	3-34
bchenco	3-36
bchpoly	3-37
bi2de	3-41
biterr	3-43
compand	3-49
convenc	3-51
cyclgen	3-53
cyclpoly	3-55
ddemod	3-57
ddemodce	3-62
de2bi	3-67
decode	3-69
demodmap	3-73
dmod	3-78
dmodce	3-82
dpcmdeco	3-86
dpcmenco	3-87
dpcmopt	3-88
encode	3-89
eyediagram	3-95
gen2par	3-97
gfadd	3-99
gfconv	3-101
gfcosets	3-103
gfdeconv	3-105
gfdiv	3-108
gffilter	3-110
gflineq	3-112
gfminpol	3-114
gfmul	3-116

gfplus .....	3-117
gfpretty .....	3-118
gfprimck .....	3-120
gfprimdf .....	3-121
gfprimfd .....	3-122
gfrank .....	3-124
gfrepcov .....	3-125
gfroots .....	3-126
gfsb .....	3-128
gftrunc .....	3-130
gftuple .....	3-131
gfweight .....	3-134
hamngen .....	3-135
hank2sys .....	3-138
hilbiir .....	3-140
istrellis .....	3-143
lloyds .....	3-145
marcumq .....	3-148
modmap .....	3-149
oct2dec .....	3-154
poly2trellis .....	3-155
qaskdeco .....	3-158
qaskenco .....	3-160
quantiz .....	3-163
randerr .....	3-165
randint .....	3-167
randsrc .....	3-168
rcosfir .....	3-170
rcosflt .....	3-172
rcosiir .....	3-175
rcosine .....	3-177
rsdeco .....	3-179
rsdecode .....	3-182
rsdecof .....	3-184
rsenco .....	3-185
rsencode .....	3-188
rsencof .....	3-190
rspoly .....	3-191
scatterplot .....	3-193

<b>symerr</b>	<b>3-195</b>
<b>syndtable</b>	<b>3-198</b>
<b>vec2mat</b>	<b>3-199</b>
<b>vitdec</b>	<b>3-201</b>
<b>wgn</b>	<b>3-205</b>

# ademod

**Purpose** Analog passband demodulator

**Syntax**

```
z = ademod(y, Fc, Fs, ' amlsb- tc' , offset, num, den) ;  
z = ademod(y, Fc, Fs, ' amlsb- tc/costas' , offset, num, den) ;  
z = ademod(y, Fc, Fs, ' amlsb- sc' , num, den) ;  
z = ademod(y, Fc, Fs, ' amlsb- sc/costas' , num, den) ;  
z = ademod(y, Fc, Fs, ' amssb' , num, den) ;  
z = ademod(y, Fc, Fs, ' qam' , num, den) ;  
z = ademod(y, Fc, Fs, ' fm' , num, den, vcoconst) ;  
z = ademod(y, Fc, Fs, ' pm' , num, den, vcoconst) ;  
z = ademod(y, Fc, [Fs phase], ... ) ;
```

Optional Inputs	Input	Default Value
	offset	Appropriate value so that each output signal has zero mean
	num, den	[ num, den] = butter(5, Fc*2/Fs) ;
	vcoconst	1

**Description** The function ademod performs analog passband demodulation. The corresponding modulation function is amod. The table below lists the demodulation schemes that ademod supports.

Demodulation Scheme	Fourth Input Argument
Amplitude demodulation	' <b>amlsb- tc</b> ' or ' <b>amlsb- tc/costas</b> '
Amplitude demodulation, double sideband suppressed carrier	' <b>amlsb- sc</b> ' or ' <b>amlsb- sc/costas</b> '
Amplitude demodulation, single sideband suppressed carrier	' <b>amssb</b> '
Quadrature amplitude demodulation	' <b>qam</b> '
Frequency demodulation	' <b>fm</b> '
Phase demodulation	' <b>pm</b> '

### For All Syntaxes

The generic syntax `z = ademod(y, Fc, Fs, ...)` demodulates the received signal that `y` represents. `Fc` is the carrier frequency in Hertz, and `Fs` is the sampling rate in Hertz. The initial phase of the carrier signal is zero.

`y` and `z` are real matrices whose sizes depend on the demodulation method:

- **(QAM method)** If `y` is a length- $n$  vector, then `z` is an  $n$ -by-2 matrix. Otherwise, if `y` is  $n$ -by- $m$ , then `z` is  $n$ -by- $2m$  and each column of `y` is processed separately. The odd-numbered columns in `z` represent in-phase components and the even-numbered columns represent quadrature components.
- **(Other methods)** `y` and `z` have the same dimensions. If `y` is a two-dimensional matrix, then each column of `y` is processed separately.

The generic syntax `z = ademod(y, Fc, [Fs phase], ...)` is the same, except that the third input argument is a two-element vector instead of a scalar. The first entry, `Fs`, is the sampling rate. The second entry, `phase`, is the initial phase of the carrier signal, measured in radians.

`ademod` uses a lowpass filter with sample time  $1/F_s$  while demodulating, in order to filter out the carrier signal. To specify the lowpass filter, include `num` and `den` in the list of input arguments. `num` and `den` are row vectors that give the coefficients, in *descending* order, of the numerator and denominator of the filter's transfer function. If `num` is empty, zero, or absent, then the default filter is a Butterworth filter whose parameters come from the command below.

`butter` is in the Signal Processing Toolbox.

```
[num, den] = butter(5, Fc*2/Fs);
```

### For Specific Syntaxes

`z = ademod(y, Fc, Fs, 'andsb-tc', offset, num, den)` implements double-sideband amplitude demodulation. `offset` is a vector whose  $k$ th entry is subtracted from the  $k$ th signal after the demodulation. If `offset` is empty, then by default `z` will be adjusted so that each column has mean zero (or, so that `z` has mean zero in case `z` is a vector).

`z = ademod(y, Fc, Fs, 'andsb-tc/costas', offset, num, den)` is the same as the syntax above, except that the algorithm includes a Costas phase-locked loop.

`z = ademod(y, Fc, Fs, 'amdsb-sc', num, den)` implements double-sideband suppressed-carrier amplitude demodulation.

`z = ademod(y, Fc, Fs, 'amdsb-sc/costas', num, den)` is the same as the syntax above, except that the algorithm includes a Costas phase-locked loop.

`z = ademod(y, Fc, Fs, 'amssb', num, den)` implements single-sideband suppressed-carrier amplitude demodulation.

`z = ademod(y, Fc, Fs, 'qam', num, den)` implements quadrature amplitude demodulation.

`z = ademod(y, Fc, Fs, 'fm', num, den, vcoconst)` implements frequency demodulation. The spectrum of the demodulated signal is between  $\min(y) + F_c$  and  $\max(y) + F_c$ . The demodulation process uses a phase-locked loop composed of a multiplier (as a phase detector), a lowpass filter, and a voltage-controlled oscillator (VCO). If  $F_s$  is a two-element vector, then its second element is the initial phase of the VCO, in radians. The optional argument `vcoconst` is a scalar that represents the VCO constant in Hz/V.

`z = ademod(y, Fc, Fs, 'pm', num, den, vcoconst)` implements phase demodulation. The demodulation process uses a phase-locked loop (which acts as an FM demodulator) cascaded with an integrator. The phase-locked loop consists of a multiplier (as a phase detector), a lowpass filter, and a voltage-controlled oscillator (VCO). If  $F_s$  is a two-element vector, then its second element is the initial phase of the VCO, in radians. The optional argument `vcoconst` is a scalar that represents the input signal's sensitivity.

## Examples

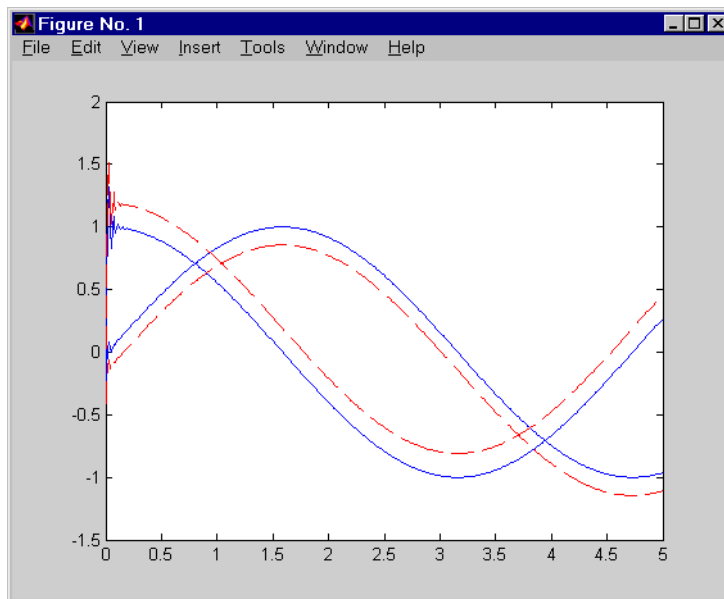
This example illustrates the use of the `offset` argument. Since the first `ademod` command uses the same `offset` value of `.3` that the `amod` command used, `z1` is similar to the original message signal. Since the second `ademod` command omits `offset`, `z2` has mean close to zero (not exactly zero because of roundoff error).

```
Fc = 25; % Carrier signal frequency
Fs = 100; % Sampling rate of signal
t = [0:1/Fs:5]'; % Times to sample the signals
x = [cos(t), sin(t)]; % Cosine signal and sine signal
y = amod(x, Fc, Fs, 'amdsb-tc', .3); % Modulate
% and shift the values up by .3.
z1 = ademod(y, Fc, Fs, 'amdsb-tc', .3); % Demodulate.
```



```
z2 = ademod(y, Fc, Fs, 'amdsb-tc'); % Demodulate.
plot(t, z1, 'b', t, z2, 'r--') % Plot recovered signal.
```

The plot shows z1 as a solid line and z2 as a dashed line.



Other examples using ademod are the Hilbert Filter Example on the reference page for amod, and in the section “Example: Varying the Filter’s Cutoff Frequency” on page 2-63.

## See Also

amod, dmod, ddemod, amodce, ademodce

# ademodce

**Purpose** Analog baseband demodulator

**Syntax**

```
z = ademodce(y, Fs, ' amdsb-tc' , offset, num, den) ;
z = ademodce(y, Fs, ' amdsb-tc/costas' , offset, num, den) ;
z = ademodce(y, Fs, ' amdsb-sc' , num, den) ;
z = ademodce(y, Fs, ' amdsb-sc/costas' , num, den) ;
z = ademodce(y, Fs, ' amssb' , num, den) ;
z = ademodce(y, Fs, ' qam' , num, den) ;
z = ademodce(y, Fs, ' fm' , num, den, vcoconst) ;
z = ademodce(y, Fs, ' pm' , num, den, vcoconst) ;
z = ademodce(y, [Fs phase], ... ) ;
```

Optional Inputs	Input	Default Value, or Default Behavior if Input is Omitted
	offset	Appropriate value so that each output signal has zero mean
	num, den	Omitting these arguments prevents ademodce from using a filter.
	vcoconst	1

**Description** The function ademodce performs analog baseband demodulation. The corresponding modulation function is amodce. The table below lists the demodulation schemes that ademodce supports.

Demodulation Scheme	Third Input Argument
Amplitude demodulation	' <b>amdsb-tc</b> '
Amplitude demodulation, double sideband suppressed carrier	' <b>amdsb-sc</b> ' or ' <b>amdsb-sc/costas</b> '
Amplitude demodulation, single sideband suppressed carrier	' <b>amssb</b> '
Quadrature amplitude demodulation	' <b>qam</b> '
Frequency demodulation	' <b>fm</b> '
Phase demodulation	' <b>pm</b> '

### For All Syntaxes

The generic syntax  $z = \text{ademodce}(y, F_s, \dots)$  demodulates the received signal that  $y$  represents.  $F_s$  is the sampling rate in Hertz. The initial phase of the carrier signal is zero.  $y$  is a complex matrix and  $z$  is a real matrix. Their sizes depend on the demodulation method:

- **(QAM method)** If  $y$  is a vector of length  $n$ , then  $z$  is an  $n$ -by-2 matrix. Otherwise, if  $y$  is  $n$ -by- $m$ , then  $z$  is  $n$ -by- $2m$  and each column of  $y$  is processed separately. The odd-numbered columns in  $z$  represent in-phase components and the even-numbered columns represent quadrature components.
- **(Other methods)**  $y$  and  $z$  have the same dimensions. If  $y$  is a two-dimensional matrix, then each column of  $y$  is processed separately.

The generic syntax  $z = \text{ademodce}(y, [F_s \text{ phase}], \dots)$  is the same, except that the second input argument is a two-element vector instead of a scalar. The first entry,  $F_s$ , is the sampling rate as described in the paragraph above. The second entry,  $\text{phase}$ , is the initial phase of the carrier signal, measured in radians.

To use a lowpass filter in the demodulation, include  $\text{num}$  and  $\text{den}$  in the list of input arguments.  $\text{num}$  and  $\text{den}$  are row vectors that give the coefficients, in *descending* order, of the numerator and denominator of the filter's transfer function. If  $\text{num}$  is empty, zero, or absent, then  $\text{ademodce}$  does not use a filter.

### For Specific Syntaxes

$z = \text{ademodce}(y, F_s, \text{'amlsb-tc'}, \text{offset}, \text{num}, \text{den})$  implements double-sideband amplitude demodulation.  $\text{offset}$  is a vector whose  $k$ th entry is subtracted from the  $k$ th column of demodulated data. If  $\text{offset}$  is empty, then by default  $z$  will be adjusted so that each column has mean zero (or, so that  $z$  has mean zero in case  $z$  is a vector).

$z = \text{ademodce}(y, F_s, \text{'amlsb-tc/costas'}, \text{offset}, \text{num}, \text{den})$  is the same as the syntax above, except that the algorithm includes a Costas phase-locked loop.

$z = \text{ademodce}(y, F_s, \text{'amlsb-sc'}, \text{num}, \text{den})$  implements double-sideband suppressed-carrier amplitude demodulation.

$z = \text{ademodce}(y, F_s, \text{'amlsb-sc/costas'}, \text{num}, \text{den})$  is the same as the syntax above, except that the algorithm includes a Costas phase-locked loop.

`z = ademodce(y, Fs, 'amssb', num, den)` implements single-sideband suppressed-carrier amplitude demodulation.

`z = ademodce(y, Fs, 'qam', num, den)` implements quadrature amplitude demodulation.

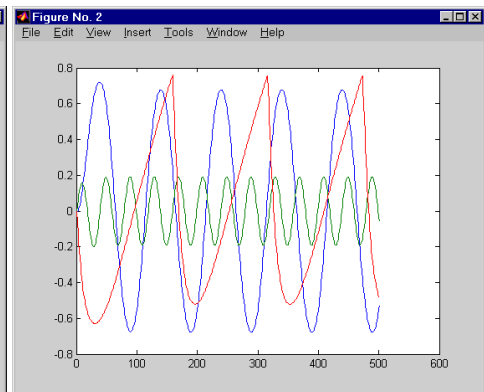
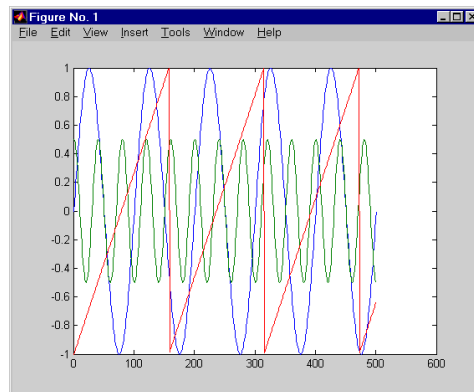
`z = ademodce(y, Fs, 'fm', num, den, vcoconst)` implements frequency demodulation. The optional argument `vcoconst` is a scalar that represents the VCO constant in the demodulation.

`z = ademodce(y, Fs, 'pm', num, den, vcoconst)` implements phase demodulation. The optional argument `vcoconst` specifies the VCO constant in the demodulation.

## Examples

The example below processes sine, cosine, and sawtooth signals simultaneously. All three signals have the same sampling rate and the same number of samples. The code also plots the original and demodulated signals.

```
Fs = 100; % Sampling rate of signal
t = [0:1/Fs:5]'; % Times to sample the signals
% Combine three signals into a three-column matrix.
% Each signal occupies one column.
x = [sin(2*pi*t), .5*cos(5*pi*t), sawtooth(4*t)];
y = amodce(x, Fs, 'fm'); % Modulate.
z = ademodce(y, Fs, 'fm'); % Demodulate.
plot(x); figure; plot(z); % Original and demodulated signals
```



Other examples using `ademodce` are in the sections “Simple Analog Modulation Example” on page 2-61 and “Example: Time Lag From Filtering” on page 2-64.

**See Also**

`amodce`, `dmodce`, `ddemodce`, `amod`, `ademod`

# amod

**Purpose** Analog passband modulator

**Syntax**

```
y = amod(x, Fc, Fs, 'amdsb-sc');
y = amod(x, Fc, Fs, 'amdsb-tc', offset);
y = amod(x, Fc, Fs, 'amssb/opt');
y = amod(x, Fc, Fs, 'amssb/opt', num, den);
y = amod(x, Fc, Fs, 'amssb/opt', hilbertflag);
y = amod(x, Fc, Fs, 'qam');
y = amod(x, Fc, Fs, 'fmi', deviation);
y = amod(x, Fc, Fs, 'pmi', deviation);
y = amod(x, Fc, [Fs phase], ...);
[y, t] = amod(...);
```

Optional Inputs	Input	Default Value, or Default Behavior if Input is Omitted
	offset	-min(min(x))
	opt	Omitting this argument causes amod to produce the lower sideband instead of the upper sideband.
	deviation	1

**Description** The function amod performs analog passband modulation. The corresponding demodulation function is ademod. The table below lists the modulation schemes that amod supports.

Modulation Scheme	Fourth Input Argument
Amplitude modulation, double sideband with transmission carrier	'amdsb-tc'
Amplitude modulation, double sideband suppressed carrier	'amdsb-sc'
Amplitude modulation, single sideband suppressed carrier	'amssb' or 'amssb/up'
Quadrature amplitude modulation	'qam'

Modulation Scheme	Fourth Input Argument
Frequency modulation	'fmi'
Phase modulation	'pmi'

### For All Syntaxes

The generic syntax  $y = \text{amod}(x, F_c, F_s, \dots)$  modulates the message signal that  $x$  represents.  $F_c$  is the carrier frequency in Hertz, and  $F_s$  is the sampling rate in Hertz. (Thus  $1/F_s$  represents the time interval between two consecutive samples in  $x$ .) The initial phase of the carrier signal is zero. By the Nyquist theorem, the sampling rate must be at least twice as large as the modulation carrier frequency.  $x$  and  $y$  are real matrices whose sizes depend on the demodulation method:

- **(QAM method)**  $x$  must have an even number of columns. The odd-numbered columns in  $x$  represent in-phase components and the even-numbered columns represent quadrature components. If  $x$  is  $n$ -by- $2m$ , then  $y$  is  $n$ -by- $m$  and each *pair* of columns of  $x$  is processed separately.
- **(Other methods)**  $x$  and  $y$  have the same dimensions. If  $x$  is a two-dimensional matrix, then each column of  $x$  is processed separately.

The generic syntax  $y = \text{amod}(x, F_c, [F_s \text{ phase}], \dots)$  is the same, except that the third input argument is a two-element vector instead of a scalar. The first entry,  $F_s$ , is the sampling rate as described in the paragraph above. The second entry, *phase*, is the initial phase of the carrier signal, measured in radians.

### For Specific Syntaxes

$y = \text{amod}(x, F_c, F_s, \text{'amlsb-tc'}, \text{offset})$  implements double-sideband amplitude modulation. *offset* is the value added to  $x$  prior to the modulation. If you omit *offset*, then its default value is  $-\min(\min(x))$ . This default value produces 100% modulation.

$y = \text{amod}(x, F_c, F_s, \text{'amlsb-sc'})$  implements double-sideband suppressed-carrier amplitude modulation.

$y = \text{amod}(x, F_c, F_s, \text{'amssb/opt'})$  implements single-sideband suppressed-carrier amplitude modulation. By default, it produces the lower

sideband; if *opt* is **up**, then the function produces the upper sideband. This syntax does a Hilbert transform in the frequency domain.

`y = amod(x, Fc, Fs, 'amssb/opt', num, den)` is the same as the syntax above, except that it specifies a time-domain Hilbert filter. `num` and `den` are row vectors that give the coefficients, in *descending* order, of the numerator and denominator of the filter's transfer function. You can use the function `hilbiir` to design the Hilbert filter.

`y = amod(x, Fc, Fs, 'amssb/opt', hilbertflag)` is the same as the syntax above, except that it uses a default time-domain Hilbert filter. The filter's transfer function is defined by `[num, den] = hilbiir(1/Fs)`, where `num` and `den` are as in the paragraph above. The input argument `hilbertflag` can have any value.

`y = amod(x, Fc, Fs, 'qam')` implements quadrature amplitude modulation. `x` is a two-column matrix whose first column represents the in-phase signal and whose second column represents the quadrature signal. `y` is a column vector.

`y = amod(x, Fc, Fs, 'fm', deviation)` implements frequency modulation. The spectrum of the modulated signal is between `min(x) + Fc` and `max(x) + Fc`. The optional argument `deviation` is a scalar that represents the frequency deviation constant of the modulation. The command `y = amod(x, Fc, Fs, 'fm', deviation)` is equivalent to the command `y = amod(x*deviation, Fc, Fs, 'fm')`.

`y = amod(x, Fc, Fs, 'pm', deviation)` implements phase modulation. The optional argument `deviation` is a scalar that represents the phase deviation constant of the modulation. The command `y = amod(x, Fc, Fs, 'pm', deviation)` is equivalent to the command `y = amod(x*deviation, Fc, Fs, 'pm')`.

`[y, t] = amod(...)` returns the computation time in `t`.

## Examples

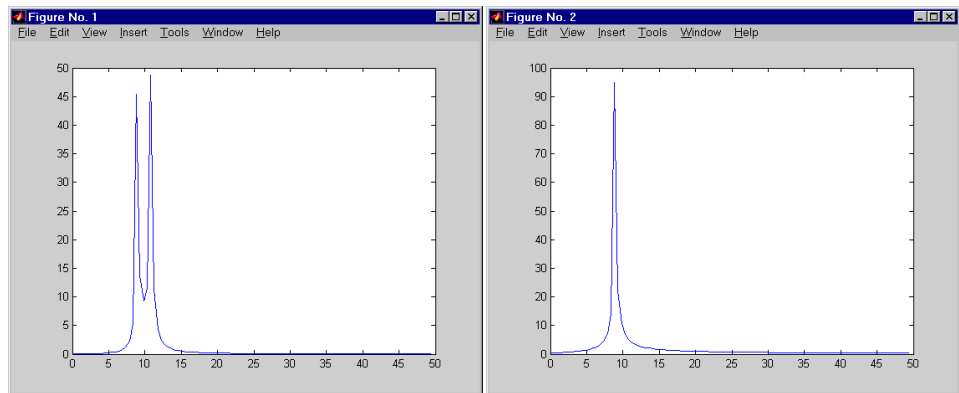
### Double- and Single-Sideband Comparison Example

The first example compares the spectra of signals after modulation using the double-sideband and single-sideband techniques. The message signal is a frequency-one sine wave and the carrier signal is a 10 Hz sine wave. The script below uses the '**amdsb-sc**' and '**amssb**' arguments in the `amod` function to



produce modulated signals `ydouble` and `ysingle`, respectively. It then plots the spectra of both modulated signals.

```
% Sample the signal 100 times per second, for 2 seconds.
Fs = 100;
t = [0:2*Fs+1]' /Fs;
Fc = 10; % Carrier frequency
x = sin(2*pi*t); % Sinusoidal signal
% Modulate x using single- and double-sideband AM
ydouble = amod(x, Fc, Fs, 'amdsb-sc');
ysingle = amod(x, Fc, Fs, 'amssb');
% Plot spectra of both modulated signals.
zdouble = fft(ydouble);
zdouble = abs(zdouble(1:length(zdouble)/2+1));
frqdouble = [0:length(zdouble)-1]*Fs/length(zdouble)/2;
plot(frqdouble, zdouble); % The plot on the left-hand side below
figure;
zsingle = fft(ysingle);
zsingle = abs(zsingle(1:length(zsingle)/2+1));
frqsingle = [0:length(zsingle)-1]*Fs/length(zsingle)/2;
plot(frqsingle, zsingle); % The plot on the right-hand side below
```



Notice that the spectrum in the left plot has two peaks; these are the lower and the upper sidebands of the modulated signal. The two sidebands are symmetrical with respect to the 10 Hz carrier frequency,  $F_c$ . The spectrum of a DSB-SC AM modulated signal is twice as wide as the input signal bandwidth.

In the right plot, there is one peak because the SSB AM technique requires `amod` to transmit only one sideband.

## Hilbert Filter Example

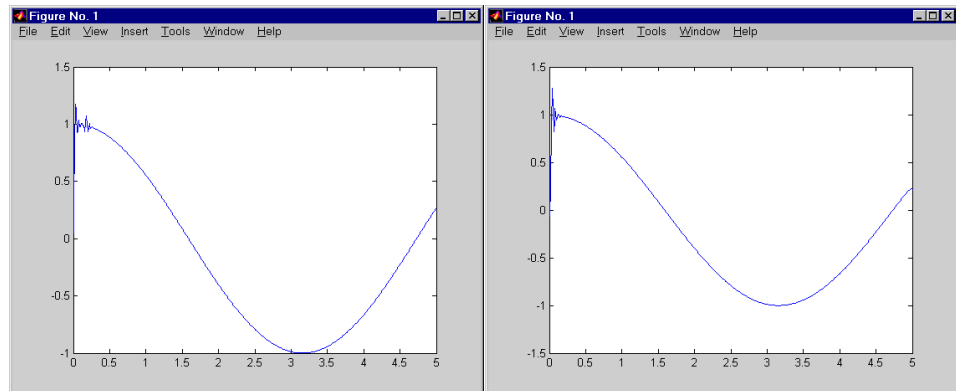
The next example uses a Hilbert filter in the time domain.

```
Fc = 25; % Carrier signal frequency
Fs = 100; % Sampling rate of signal
[numh, denh] = hilbiir(1/Fs, 15/Fs, 15); % Design Hilbert filter.
t = [0:1/Fs:5]'; % Times to sample the signal
x = cos(t); % Signal is a cosine wave.
y = amod(x, Fc, [Fs pi/4], 'amssb', numh, denh); % Modulate,
% using a Hilbert filter in the time domain.
z = ademod(y, Fc, [Fs pi/4], 'amssb'); % Demodulate.
plot(t, z) % Plot recovered signal.
```

The resulting plot is on the left below. If you replace the sixth line above with

```
y = amod(x, Fc, [Fs pi/4], 'amssb'); % Modulate,
```

then modulation uses a Hilbert transform in the frequency domain. The result is the plot on the right below. The two plots differ slightly in their initial errors.



## See Also

`ademod`, `dmod`, `ddemod`, `amodce`, `ademodce`

**Purpose**                      Analog baseband modulator

**Syntax**

```
y = amodce(x, Fs, 'amdsb-tc', offset);
y = amodce(x, Fs, 'amdsb-sc');
y = amodce(x, Fs, 'amssb');
y = amodce(x, Fs, 'amssb/time', num, den);
y = amodce(x, Fs, 'amssb/time');
y = amodce(x, Fs, 'qam');
y = amodce(x, Fs, 'fm', deviation);
y = amodce(x, Fs, 'pm', deviation);
y = amodce(x, [Fs phase], ...);
```

Optional Inputs	Input	Default Value, or Default Behavior if Input is Omitted
	offset	-min(min(x))
	deviation	1

**Description**                      The function amodce performs analog baseband modulation. The corresponding demodulation function is ademodce. The table below lists the modulation schemes that amodce supports.

Modulation Scheme	Third Input Argument
Amplitude modulation, double sideband	'amdsb-tc'
Amplitude modulation, double sideband suppressed carrier	'amdsb-sc'
Amplitude modulation, single sideband suppressed carrier	'amssb' or 'amssb/time'
Quadrature amplitude modulation	'qam'
Frequency modulation	'fm'
Phase modulation	'pm'

## For All Syntaxes

The generic syntax  $y = \text{amodce}(x, F_s, \dots)$  modulates the message signal that  $x$  represents, and returns the modulated signal's complex envelope. The input and output signals share the same sampling rate  $F_s$ , measured in Hertz. (Thus  $1/F_s$  represents the time interval between two consecutive samples in  $x$ .) The initial phase of the carrier signal is zero.  $x$  is a real matrix and  $y$  is a complex matrix. Their sizes depend on the modulation method:

- **(QAM method)**  $x$  must have an even number of columns. The odd-numbered columns in  $x$  represent in-phase components and the even-numbered columns represent quadrature components. If  $x$  is  $n$ -by- $2m$ , then  $y$  is  $n$ -by- $m$  and each *pair* of columns of  $x$  is processed separately.
- **(Other methods)**  $x$  and  $y$  have the same dimensions. If  $x$  is a two-dimensional matrix, then each column of  $x$  is processed separately.

The generic syntax  $y = \text{amodce}(x, [F_s \text{ phase}], \dots)$  is the same, except that the second input argument is a two-element vector instead of a scalar. The first entry,  $F_s$ , is the sampling rate as described in the paragraph above. The second entry, *phase*, is the initial phase of the carrier signal, measured in radians.

## For Specific Syntaxes

$y = \text{amodce}(x, F_s, ' \text{amdsb- tc}' , \text{offset})$  implements double-sideband amplitude modulation. *offset* is the value added to  $x$  prior to the modulation. If you omit *offset*, then its default value is  $-\min(\min(x))$ . This default value produces 100% modulation.

$y = \text{amodce}(x, F_s, ' \text{amdsb- sc}' )$  implements double-sideband suppressed-carrier amplitude modulation.

$y = \text{amodce}(x, F_s, ' \text{amssb}' )$  implements single-sideband suppressed-carrier amplitude modulation. By default, it produces the lower sideband. It does a Hilbert transform in the frequency domain.

$y = \text{amodce}(x, F_s, ' \text{amssb/time}' , \text{num}, \text{den})$  is the same as the syntax above, except that it specifies a time-domain Hilbert filter. *num* and *den* are row vectors that give the coefficients, in *descending* order, of the numerator and denominator of the filter's transfer function. You can use the function `hilb1r` to design the Hilbert filter.

`y = amodce(x, Fs, 'amssb/time')` is the same as the syntax above, except that it uses a default time-domain Hilbert filter. The filter's transfer function is defined by `[num, den] = hilbiir(1/Fs)`, where `num` and `den` are as in the paragraph above.

`y = amodce(x, Fs, 'qam')` implements quadrature amplitude modulation. `x` is a two-column matrix whose first column represents the in-phase signal and whose second column represents the quadrature signal. `y` is a column vector.

`y = amodce(x, Fs, 'fm', deviation)` implements frequency modulation. The bandwidth of the modulated signal is  $\max(x) - \min(x)$ . The optional argument `deviation` is a scalar that represents the frequency deviation constant of the modulation.

`y = amodce(x, Fs, 'pm', deviation)` implements phase modulation. The optional argument `deviation` is a scalar that represents the phase deviation constant of the modulation.

## Examples

This example is similar to the one under the heading “Hilbert Filter Example” on the `amod` reference page, except that it uses baseband simulation. The plots in the passband (`amod`) example show far more obvious errors in the recovered signal. The output from this example shows that the average difference between the original and recovered signals is smaller than  $10^{-16}$ .

```
Fs = 100; % Sampling rate of signal
[numh, denh] = hilbiir(1/Fs, 15/Fs, 15); % Design Hilbert filter.
t = [0:1/Fs:5]'; % Times to sample the signal
x = cos(t); % Signal is a cosine wave.
y = amodce(x, [Fs pi/4], 'amssb/time', numh, denh); % Modulate,
% using a Hilbert filter in the time domain.
z = ademodce(y, [Fs pi/4], 'amssb'); % Demodulate.
d = ceil(log10(sum(abs(x-z))/length(x)))

d =
```

- 16

Other examples using `amodce` are in the sections “Representing Analog Signals” on page 2-59 and “Simple Analog Modulation Example” on page 2-61.

## See Also

`ademodce`, `dmodce`, `ddemodce`, `amod`, `ademod`

# apkconst

---

**Purpose** Plot a combined circular ASK-PSK signal constellation

**Syntax**

```
apkconst (numsi g);  
apkconst (numsi g, amp);  
apkconst (numsi g, amp, phs);  
apkconst (numsi g, amp, 'n');  
apkconst (numsi g, amp, phs, pl ot spec);  
y = apkconst (. . .);
```

**Description** APK refers to a hybrid of amplitude- and phase-keying modulation. See the reference listed below for more details.

`apkconst (numsi g)` plots a circular signal constellation. `numsi g` is a vector of positive integers. The plot contains `length(numsi g)` circles. The  $k$ th circle has radius  $k$  and contains `numsi g(k)` evenly spaced constellation points. One point on each circle has zero phase.

`apkconst (numsi g, amp)` is the same as the previous syntax, except that `amp(k)` is the radius of the  $k$ th circle. `amp` is a vector of positive real numbers. The lengths of `amp` and `numsi g` must be the same.

`apkconst (numsi g, amp, phs)` is the same as the previous syntax, except that it is not necessarily true that one point on each circle has zero phase. However, one point on the  $k$ th circle has phase `phs(k)`. The lengths of `phs`, `amp` and `numsi g` must all be the same.

`apkconst (numsi g, amp, phs, 'n')` is the same as the previous syntax, except that the plot includes a number next to each constellation point. The number indicates how symbols would be mapped to constellation points if you were using `numsi g`, `amp`, and `phs` in modulation and demodulation functions such as `dmodce/ddemodce` or `modmap/demodmap`.

`apkconst (numsi g, amp, phs, pl ot spec)` is the same as `apkconst (numsi g, amp, phs)`, except that `pl ot spec` influences the appearance of the constellation points via MATLAB's `pl ot` function. `pl ot spec` is a

two-character string made up of one character from each odd-numbered column in the table below.

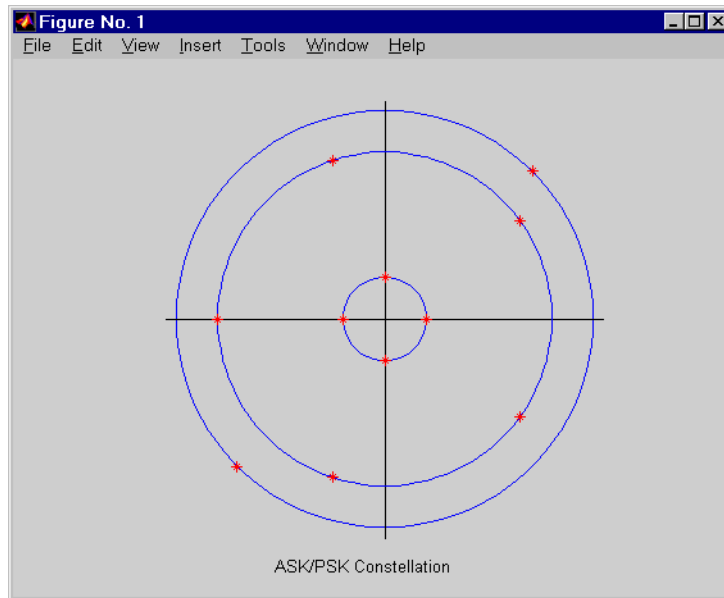
Color Character	Meaning	Marker-Type Character	Meaning
y	yellow	.	point
m	magenta	o	circle
c	cyan	x	cross
r	red	+	plus sign
g	green	*	asterisk
b	blue	s	square
w	white	d	diamond
k	black	v	triangle (down)
		^	triangle (up)
		<	triangle (left)
		>	triangle (right)
		p	five-pointed star
		h	six-pointed star

`y = apkconst(...)` does *not* produce a plot, but instead returns a complex vector `y` that represents the coordinates of the points in the constellation. The real part of `y` gives the in-phase component of each point and the imaginary part of `y` gives the quadrature component of each point.

## Examples

The command below produces a plot having three circles. One circle has radius 1 and four points, one of which has zero phase. Another circle has radius 4 and five points, one of which has phase  $\pi$ . The outermost circle has radius 5 and two points, one of which has phase  $\pi/4$ . The plot follows.

```
apkconst([4 5 2], [1 4 5], [0 pi pi/4])
```



The command below produces a vector containing the coordinates in the complex plane of the points in the figure above.

```
y = apkconst([4 5 2],[1 4 5],[0 pi pi/4])
```

```
y =
```

```
Columns 1 through 4
```

```
1.0000      0.0000 + 1.0000i  -1.0000 + 0.0000i  -0.0000 - 1.0000i
```

```
Columns 5 through 8
```

```
-4.0000 + 0.0000i  -1.2361 - 3.8042i   3.2361 - 2.3511i   3.2361 + 2.3511i
```

```
Columns 9 through 11
```

```
-1.2361 + 3.8042i   3.5355 + 3.5355i  -3.5355 - 3.5355i
```

## See Also

dmod, modmap, ddemod, demodmap



**References**

Thomas, C. Melvil, Michael Y. Weidner, and S. H. Durrani. "Digital Amplitude-Phase Keying with M-ary Alphabets." *IEEE Transactions on Communications*. Vol Com-22, No. 2, Feb. 1974, 168-180.

**Purpose** Add white Gaussian noise to a signal

**Syntax**

```
y = awgn(x, snr);  
y = awgn(x, snr, si gpower);  
y = awgn(x, snr, 'measured');  
y = awgn(x, snr, si gpower, state);  
y = awgn(x, snr, 'measured', state);  
y = awgn(..., powertype);
```

**Description** `y = awgn(x, snr)` adds white Gaussian noise to the vector signal `x`. The scalar `snr` specifies the signal-to-noise ratio in decibels. If `x` is complex, then `awgn` adds complex noise. This syntax assumes that the power of `x` is 0 dB.

`y = awgn(x, snr, si gpower)` is the same as the syntax above, except that `si gpower` is the power of `x` in dB.

`y = awgn(x, snr, 'measured')` is the same as `y = awgn(x, snr)`, except that `awgn` measures the power of `x` before adding noise.

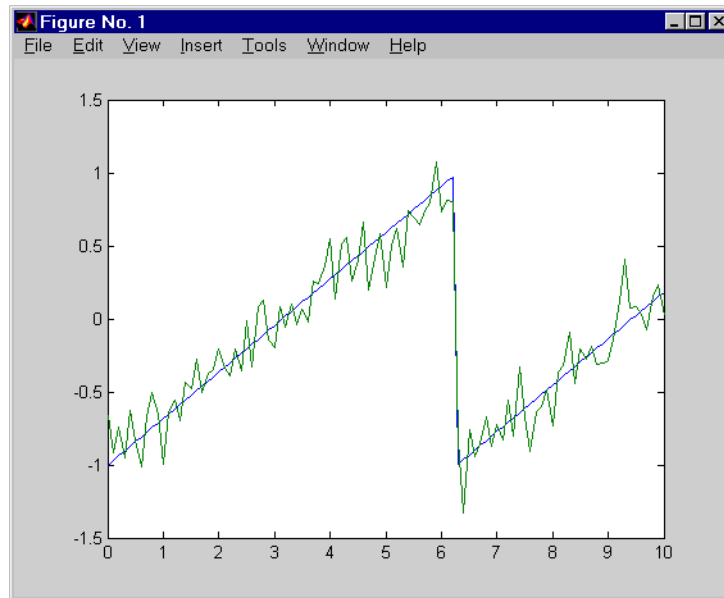
`y = awgn(x, snr, si gpower, state)` is the same as `y = awgn(x, snr, si gpower)`, except that `awgn` first resets the state of MATLAB's normal random number generator `randn` to the integer `state`.

`y = awgn(x, snr, 'measured', state)` is the same as `y = awgn(x, snr, 'measured')`, except that `awgn` first resets the state of MATLAB's normal random number generator `randn` to the integer `state`.

`y = awgn(..., powertype)` is the same as the previous syntaxes, except that the string `powertype` specifies the units of `snr` and `si gpower`. Choices for `powertype` are `'db'` and `'linear'`. Linear power is measured in Watts.

**Examples** The commands below add white Gaussian noise to a sawtooth signal. It then plots the original and noisy signals.

```
t = 0: .1: 10;  
x = sawtooth(t); % Create sawtooth signal.  
y = awgn(x, 10, 'measured'); % Add white Gaussian noise.  
plot(t, x, t, y) % Plot both signals.
```



**See Also**

wgn, randn

**Purpose** BCH decoder

**Syntax**

```
msg = bchdeco(code, k, t);  
msg = bchdeco(code, k, t, pri mpoly);  
[msg, err] = bchdeco(...);  
[msg, err, ccode] = bchdeco(...);
```

**Description** `msg = bchdeco(code, k, t)` decodes `code` using the BCH method. `k` is the message length. The codeword length  $n$  must have the form  $2^m - 1$  for some integer  $m$  greater than or equal to 3. `code` is a binary matrix with  $n$  columns, each row of which represents one codeword. `msg` is a binary matrix with `k` columns, each row of which represents one message. `t` is the error-correction capability. BCH decoding requires a primitive polynomial for  $GF(2^m)$ ; this syntax uses MATLAB's default primitive polynomial, `gfprimdf(m)`.

`msg = bchdeco(code, k, t, pri mpoly)` is the same as the first syntax, except that `pri mpoly` is a row vector that gives the coefficients, in order of ascending powers, of the primitive polynomial for  $GF(2^m)$  that will be used during processing.

`[msg, err] = bchdeco(...)` returns a column vector `err` that gives information about error correction. A nonnegative integer in `err(r)` indicates the number of errors corrected in the  $r$ th codeword; a negative integer indicates that there are more errors in the  $r$ th codeword than can be corrected.

`[msg, err, ccode] = bchdeco(...)` returns the corrected code in `ccode`.

**Examples** The script below encodes a (random) message, simulates the addition of noise to the code, and then decodes the message.

```
m = 4; n = 2^m - 1; % Codeword length  
params = bchpoly(n);  
% Arbitrarily focus on 3rd row of params.  
k = params(3, 2); % Codeword length  
t = params(3, 3); % Error-correction capability  
msg = randint(100, k);  
code = bchenco(msg, n, k); % Encode the message.  
% Corrupt up to t bits in each codeword.  
noisycode = rem(code + randerr(100, n, 1:t), 2);
```

```

% Decode the noisy code.
[newmsg, err, ccode] = bchdeco(noisycode, k, t);
if ccode==code
    disp('All errors were corrected. ')
end
if newmsg==msg
    disp('The message was recovered perfectly. ')
end

```

In this case, all errors are corrected and the message is recovered perfectly. However, if the ninth line is changed to

```
noisycode = rem(code + randerr(100, n, 1: (t+1)), 2);
```

then some codewords will contain more than  $t$  errors. This is too many errors, and some will go uncorrected.

## See Also

bchenco, encode, decode

# bchenco

---

**Purpose** BCH encoder

**Syntax** `code = bchenco(msg, n, k);`  
`code = bchenco(msg, n, k, genpoly);`

**Description** `code = bchenco(msg, n, k)` encodes `msg` using the BCH technique and the generator polynomial `genpoly = bchpoly(n, k)`. `n` is the codeword length and `k` is the message length. `msg` is a binary matrix with `k` columns. Each row of `msg` represents a message. `code` is a binary matrix with `n` columns. Each row of `code` represents a codeword.

`code = bchenco(msg, n, k, genpoly)` is the same as the first syntax, except that `genpoly` is a row vector that gives the coefficients of the generator polynomial in order of ascending powers.

**Examples** See the example on the reference page for the function `bchdeco`.

**See Also** `bchdeco`, `encode`, `decode`, `bchpoly`, `cyclgen`

**Purpose** Produce parameters or generator polynomial for binary BCH code

**Syntax**

```
bchpoly  
params = bchpoly  
params = bchpoly(n);  
genpoly = bchpoly(n, k);  
genpoly = bchpoly(primpoly, k);  
[genpoly, factors] = bchpoly(..., k);  
[genpoly, factors, cosets] = bchpoly(..., k);  
[genpoly, factors, cosets, parmat] = bchpoly(..., k);  
[genpoly, factors, cosets, parmat, errorcorr] = bchpoly(..., k);
```

**Description** bchpoly produces a figure window containing a table that lists valid codeword and message lengths of binary BCH codes, as well as the corresponding error-correction capabilities. The codeword lengths listed are 7, 15, 31, 63, 127, 255, and 511. The codeword lengths, message length, and error-correction capabilities are denoted by N, K, and T, respectively.

params = bchpoly produces a three-column matrix containing the same information that is in the table mentioned in the syntax above. The first column of params gives the codeword length, the second column gives the message length, and the third column gives the error-correction capability.

params = bchpoly(n) produces a matrix params containing valid codeword and message lengths of binary BCH codes in its first and second columns, respectively. If  $n < 1024$ , then params has a third column that lists the corresponding error-correction capabilities. The codeword lengths listed in params are all equal to the smallest number of the form  $2^m - 1$  that is at least as big as n, where  $m$  is an integer greater than or equal to 3.

genpoly = bchpoly(n, k) produces a generator polynomial for a binary BCH code having codeword length n and message length k. genpoly is a row vector that gives the coefficients, in order of ascending powers, of the generator polynomial. n must have the form  $2^m - 1$  for some integer  $m$  greater than or equal to 3. k must be a valid message length, as reported in the second column of the output of the command genpoly = bchpoly(n). The primitive polynomial used for the  $GF(2^m)$  calculations is MATLAB's default primitive polynomial, gfprimdf(m).

`genpoly = bchpoly(primpoly, k)` produces a generator polynomial for a binary BCH code having codeword length  $n$  and message length  $k$ . `primpoly` represents a degree- $m$  primitive polynomial for the field  $\text{GF}(2^m)$ . Both `primpoly` and `genpoly` are row vectors that represent polynomials by giving the coefficients in order of ascending powers. Given the degree  $m$  of the primitive polynomial, the message length  $n$  is  $2^m-1$ .  $k$  must be a valid message length, as reported in the second column of the output of the command `genpoly = bchpoly(n)`.

The remaining syntaxes, of the form

`[genpoly, ...] = bchpoly(..., k)`

return some or all of the output variables listed in the table below.

**Table 3-12: Additional Output Variables for `bchpoly(...,k)`**

Output Variable	Significance	Format
<code>factors</code>	Irreducible factors of the generator polynomial	Binary matrix, each row of which gives the coefficients of a factor polynomial in order of ascending powers
<code>cosets</code>	Cyclotomic cosets of the field $\text{GF}(2^m)$	Same as <code>gfcosets(m)</code>
<code>parmat</code>	Parity-check matrix of the code	$(n-k)$ -by- $n$ binary matrix
<code>errorcorr</code>	Error-correction capability of the code	Positive integer

## Examples

The script below uses `bchpoly` to find out what message lengths are valid for a BCH code with codeword length  $2^4-1$ . It then chooses one of the possible message lengths and uses `bchpoly` to find the generator polynomial and parity-check matrix for such a code.

```
m = 4;
n = 2^m-1; % Codeword length is 15.
% Want to find out possible valid message lengths.
```



```

params = bchpoly(n);
disp(['Possible message lengths are ', num2str(params(:, 2)')])
disp(' ')

ii = 1; % Arbitrarily choose first row.
k = params(ii, 2); % Message lengths are in 2nd column.
% Get generator polynomial and other facts.
[genpoly, factors, cosets, parmat, errorcorr] = bchpoly(n, k);
disp(['For k = ', num2str(k), ' the generator polynomial is'])
gfpretty(genpoly)
disp('and the parity-check matrix is')
parmat

```

The full output is below.

Possible message lengths are 11 7 5

For k = 11 the generator polynomial is

$$1 + X + X^4$$

and the parity-check matrix is

parmat =

Columns 1 through 12

1	0	0	0	1	0	0	1	1	0	1	0
0	1	0	0	1	1	0	1	0	1	1	1
0	0	1	0	0	1	1	0	1	0	1	1
0	0	0	1	0	0	1	1	0	1	0	1

Columns 13 through 15

1	1	1
1	0	0
1	1	0
1	1	1

## See Also

cyclpoly, encode, decode

## References

Peterson, W. Wesley and E. J. Weldon, Jr. *Error-correcting Codes*, 2nd ed. Cambridge, Mass.: MIT Press, 1972.

<b>Purpose</b>	Convert binary vectors to decimal numbers
<b>Syntax</b>	<pre> d = bi2de(b); d = bi2de(b, flag) d = bi2de(b, p); d = bi2de(b, p, flag); </pre>
<b>Description</b>	<p><code>d = bi2de(b)</code> converts a binary row vector <code>b</code> to a nonnegative decimal integer. If <code>b</code> is a matrix, then each row is interpreted separately as a binary number. In this case, the output <code>d</code> is a column vector, each element of which is the decimal representation of the corresponding row of <code>b</code>.</p> <hr/> <p><b>Note</b> By default, <code>bi2de</code> interprets the first column of <code>b</code> as the <i>lowest-order</i> digit.</p> <hr/> <p><code>d = bi2de(b, flag)</code> is the same as the syntax above, except that <code>flag</code> is a string that determines whether the first column of <code>b</code> contains the lowest-order or highest-order digits. Possible values for <code>flag</code> are '<b>right-msb</b>' and '<b>left-msb</b>'. The value '<b>right-msb</b>' produces the default behavior.</p> <p><code>d = bi2de(b, p)</code> converts a base-<code>p</code> row vector <code>b</code> to a nonnegative decimal integer, where <code>p</code> is an integer greater than or equal to two. The first column of <code>b</code> is the <i>lowest</i> base-<code>p</code> digit. If <code>b</code> is a matrix, then the output <code>d</code> is a nonnegative decimal vector, each row of which is the decimal form of the corresponding row of <code>b</code>.</p> <p><code>d = bi2de(b, p, flag)</code> is the same as the syntax above, except that <code>flag</code> is a string that determines whether the first column of <code>b</code> contains the lowest-order or highest-order digits. Possible values for <code>flag</code> are '<b>right-msb</b>' and '<b>left-msb</b>'. The value '<b>right-msb</b>' produces the default behavior.</p>
<b>Examples</b>	<p>The code below generates a matrix that contains binary representations of five random numbers between 0 and 15. It then converts all five numbers to decimal integers.</p> <pre> b = randi nt(5, 4);    % Generate a 5-by-4 random binary matrix. de = bi2de(b); </pre>

```
disp('      Dec          Bi nary' )
disp(' -----' )
disp([ de, b])
```

Sample output is below. Your results may vary since the numbers are random.

Dec	Bi nary			
-----	-----	-----	-----	-----
13	1	0	1	1
7	1	1	1	0
15	1	1	1	1
4	0	0	1	0
9	1	0	0	1

The command below converts a base-five number into its decimal counterpart, using the leftmost base-five digit (4 in this case) as the most significant digit. The example reflects the fact that  $4(5^3) + 2(5^2) + 5^0 = 551$ .

```
d = bi2de([4 2 0 1], 5, 'left-msb' )

d =

    551
```

See Also

de2bi

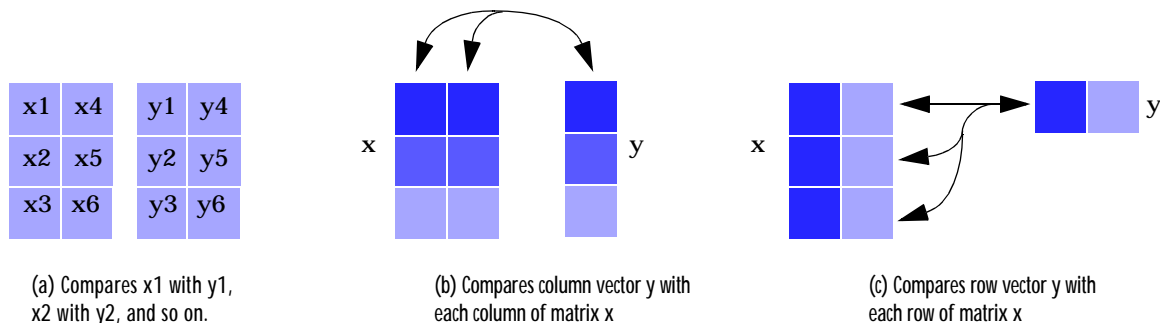
**Purpose** Compute number of bit errors and bit error rate

**Syntax**

```
[number, ratio] = biterr(x, y);
[number, ratio] = biterr(x, y, k);
[number, ratio] = biterr(..., flag);
[number, ratio, individual] = biterr(...)
```

**Description** For All Syntaxes

The `biterr` function compares unsigned binary representations of elements in `x` with those in `y`. The schematics below illustrate how the shapes of `x` and `y` determine which elements `biterr` compares.



Each element of `x` and `y` must be a nonnegative decimal integer; `biterr` converts each element into its natural unsigned binary representation. `number` is a scalar or vector that indicates the number of bits that differ. `ratio` is `number` divided by the *total number of bits*. The total number of bits, the size of `number`, and the elements that `biterr` compares are determined by the dimensions of `x` and `y` and by the optional parameters.

**For Specific Syntaxes**

`[number, ratio] = biterr(x, y)` compares the elements in `x` and `y`. If the largest among all elements of `x` and `y` has exactly  $k$  bits in its simplest binary representation, then the total number of bits is  $k$  times the number of entries in the *smaller* input. The sizes of `x` and `y` determine which elements are compared:

- If  $x$  and  $y$  are matrices of the same dimensions, then `biterr` compares  $x$  and  $y$  element-by-element. `number` is a scalar. See schematic (a) in the figure.
- If one is a row (respectively, column) vector and the other is a two-dimensional matrix, then `biterr` compares the vector element-by-element with *each row (resp., column)* of the matrix. The length of the vector must equal the number of columns (resp., rows) in the matrix. `number` is a column (resp., row) vector whose  $m$ th entry indicates the number of bits that differ when comparing the vector with the  $m$ th row (resp., column) of the matrix. See schematics (b) and (c) in the figure.

`[number, ratio] = biterr(x, y, k)` is the same as the first syntax, except that it considers each entry in  $x$  and  $y$  to have  $k$  bits. The total number of bits is  $k$  times the number of entries of the smaller of  $x$  and  $y$ . An error occurs if the binary representation of an element of  $x$  or  $y$  would require more than  $k$  digits.

`[number, ratio] = biterr(x, y, k, flag)` is similar to the previous syntaxes, except that `flag` can override the defaults that govern which elements `biterr` compares and how `biterr` computes the outputs. The possible values of `flag` are '**row-wise**', '**column-wise**', and '**overall**'. The table below describes the differences that result from various combinations of inputs. As always, `ratio` is `number` divided by the total number of bits. If you do not provide `k` as an input argument, then the function defines it internally as the number of bits in the simplest binary representation of the largest among all elements of  $x$  and  $y$ .

Table 3-13: Comparing a Two-Dimensional Matrix *x* with Another Input *y*

Shape of <i>y</i>	flag	Type of Comparison	number	Total Number of Bits
Two-dimensional matrix	'overall' (default)	Element-by-element	Total number of bit errors	k times number of entries of <i>y</i>
	'row-wise'	<i>m</i> th row of <i>x</i> vs. <i>m</i> th row of <i>y</i>	Column vector whose entries count bit errors in each row	k times number of entries of <i>y</i>
	'column-wise'	<i>m</i> th column of <i>x</i> vs. <i>m</i> th column of <i>y</i>	Row vector whose entries count bit errors in each column	k times number of entries of <i>y</i>
Row vector	'overall'	<i>y</i> vs. each row of <i>x</i>	Total number of bit errors	k times number of entries of <i>x</i>
	'row-wise' (default)	<i>y</i> vs. each row of <i>x</i>	Column vector whose entries count bit errors in each row of <i>x</i>	k times size of <i>y</i>
Column vector	'overall'	<i>y</i> vs. each column of <i>x</i>	Total number of bit errors	k times number of entries of <i>x</i>
	'column-wise' (default)	<i>y</i> vs. each column of <i>x</i>	Row vector whose entries count bit errors in each column of <i>x</i>	k times size of <i>y</i>

[ number, ratio, individual ] = biterr(...) returns a matrix *individual* whose dimensions are those of the larger of *x* and *y*. Each entry of *individual* corresponds to a comparison between a pair of elements of *x* and *y*, and specifies the number of bits by which the elements in the pair differ.

## Examples

### Example 1

The commands below compare the column vector `[0; 0; 0]` to each column of a random binary matrix. The output is the number, proportion, and locations of ones in the matrix. In this case, `i n d i v i d u a l` is the same as the random matrix.

```
format rat;
[number, ratio, i n d i v i d u a l] = biterr([0; 0; 0], randint(3, 5))

number =

     2         0         0         3         1

ratio =

    2/3         0         0         1    1/3

i n d i v i d u a l =

     1         0         0         1         0
     1         0         0         1         0
     0         0         0         1         1
```

### Example 2

The commands below illustrate the use of *flag* to override the default row-by-row comparison. Notice that `number` and `ratio` are scalars, while `i n d i v i d u a l` has the same dimensions as the larger of the first two arguments of `biterr`.

```
format rat;
[number, ratio, i n d i v i d u a l] = biterr([1 2; 3 4], [1 3], 3, 'overall')

number =

     5
```



```
ratio =
```

```
5/12
```

```
individual =
```

```
0      1
1      3
```

### Example 3

The script below adds errors to 10% of the elements in a matrix. Each entry in the matrix is a two-bit number in decimal form. The script computes the bit error rate using `biterr` and the symbol error rate using `symerr`.

```
x = randint(100, 100, 4); % Original signal
% Create errors to add to ten percent of the elements of x.
% Errors can be either 1, 2, or 3 (not zero).
errorplace = (rand(100, 100) > .9); % Where to put errors
errorvalue = randint(100, 100, [1, 3]); % Value of the errors
error = errorplace.*errorvalue;
y = rem(x+error, 4); % Signal with errors added, mod 4
format short
[num_bit, ratio_bit] = biterr(x, y, 2)
[num_sym, ratio_sym] = symerr(x, y)
```

Sample output is below. Notice that `ratio_sym` is close to the target value of 0.10. Your results might vary because the example uses random numbers.

```
num_bit =
```

```
1304
```

```
ratio_bit =
```

```
0.0652
```

# biterr

---

num\_sym =

981

ratio\_sym =

0.0981

## See Also

symerr

**Purpose**

Source code mu-law or A-law compressor or expander

**Syntax**

```
out = compand(i n, mu, maxi m);  
out = compand(i n, mu, maxi m, ' mu/compressor' );  
out = compand(i n, mu, maxi m, ' mu/expander' );  
out = compand(i n, A, maxi m, ' A/compressor' );  
out = compand(i n, A, maxi m, ' A/expander' );
```

**Description**

`out = compand(i n, param, maxi m)` implements a  $\mu$ -law compressor for the input vector `i n`. `mu` specifies  $\mu$  and `maxi m` is the input signal's maximum magnitude. `out` has the same dimensions and maximum magnitude as `i n`.

`out = compand(i n, mu, maxi m, ' mu/compressor' )` is the same as the syntax above.

`out = compand(i n, mu, maxi m, ' mu/expander' )` implements a  $\mu$ -law expander for the input vector `i n`. `mu` specifies  $\mu$  and `maxi m` is the input signal's maximum magnitude. `out` has the same dimensions and maximum magnitude as `i n`.

`out = compand(i n, A, maxi m, ' A/compressor' )` implements an A-law compressor for the input vector `i n`. The scalar `A` is the A-law parameter, and `maxi m` is the input signal's maximum magnitude. `out` is a vector of the same length and maximum magnitude as `i n`.

`out = compand(i n, A, maxi m, ' A/expander' )` implements an A-law expander for the input vector `i n`. The scalar `A` is the A-law parameter, and `maxi m` is the input signal's maximum magnitude. `out` is a vector of the same length and maximum magnitude as `i n`.

---

**Note** The prevailing parameters used in practice are  $\mu = 255$  and  $A = 87.6$ .

---

**Examples**

The examples below illustrate the fact that compressors and expanders perform inverse operations.

```
compressed = compand(1:5, 87.6, 5, ' a/compressor' )
```

# compand

---

compressed =

3. 5296      4. 1629      4. 5333      4. 7961      5. 0000

expanded = compand(compressed, 87.6, 5, 'a/expander')

expanded =

1. 0000      2. 0000      3. 0000      4. 0000      5. 0000

## Algorithm

For a given signal  $x$ , the output of the  $\mu$ -law compressor is

$$y = \frac{V \log(1 + \mu|x|/V)}{\log(1 + \mu)} \operatorname{sgn}(x)$$

where  $V$  is the maximum value of the signal  $x$ ,  $\mu$  is the  $\mu$ -law parameter of the compander,  $\log$  is the natural logarithm and  $\operatorname{sgn}$  is the signum function (`sign` in MATLAB).

The output of the A-law compressor is

$$y = \begin{cases} \frac{A|x|}{1 + \log A} \operatorname{sgn}(x) & \text{for } 0 \leq |x| \leq \frac{V}{A} \\ \frac{V(1 + \log(A|x|/V))}{1 + \log A} \operatorname{sgn}(x) & \text{for } \frac{V}{A} < |x| \leq V \end{cases}$$

where  $A$  is the A-law parameter of the compander and the other elements are as in the  $\mu$ -law case.

## See Also

`quantiz`, `lloyd`s, `dpcmopt`, `dpcmenco`, `dpcmdeco`

## References

Sklar, Bernard. *Digital Communications: Fundamentals and Applications*. Englewood Cliffs, N.J.: Prentice-Hall, 1988.

<b>Purpose</b>	Convolutionally encode binary data
<b>Syntax</b>	<pre>code = convenc(msg, trellis); code = convenc(msg, trellis, initstate); [code, finalstate] = convenc(...);</pre>
<b>Description</b>	<p><code>code = convenc(msg, trellis)</code> encodes the binary vector <code>msg</code> using the convolutional encoder whose MATLAB trellis structure is <code>trellis</code>. For details about MATLAB trellis structures, see “Trellis Description of a Convolutional Encoder” on page 2-46. Each symbol in <code>msg</code> consists of <math>\log_2(\text{trellis.numInputSymbols})</math> bits. The vector <code>msg</code> contains one or more symbols. The output vector <code>code</code> contains one or more symbols, each of which consists of <math>\log_2(\text{trellis.numOutputSymbols})</math> bits.</p> <p><code>code = convenc(msg, trellis, initstate)</code> is the same as the syntax above, except that <code>initstate</code> specifies the starting state of the encoder registers. The scalar <code>initstate</code> is an integer between 0 and <code>trellis.numStates-1</code>. If the encoder schematic has more than one input stream, then the shift register that receives the first input stream provides the least significant bits in <code>initstate</code>, while the shift register that receives the last input stream provides the most significant bits in <code>initstate</code>. To use the default value for <code>initstate</code>, specify <code>initstate</code> as 0 or [].</p> <p><code>[code, finalstate] = convenc(...)</code> encodes the input message and also returns in <code>finalstate</code> the encoder’s state. <code>finalstate</code> has the same format as <code>initstate</code>.</p>
<b>Examples</b>	<p>The command below encodes five two-bit symbols using a rate 2/3 convolutional code. A schematic of this encoder is on the reference page for the <code>poly2trellis</code> function.</p> <pre>code1 = convenc(randint(10, 1, 2, 123), ... poly2trellis([5 4], [27 33 0; 0 5 13]));</pre> <p>The commands below define the encoder’s trellis structure explicitly and then use <code>convenc</code> to encode ten one-bit symbols. A schematic of this encoder is in the section, “Trellis Description of a Convolutional Encoder” on page 2-46.</p> <pre>trell = struct('numInputSymbols', 2, 'numOutputSymbols', 4, ... 'numStates', 4, 'nextStates', [0 2; 0 2; 1 3; 1 3], ...</pre>

```
'outputs', [0 3; 1 2; 3 0; 2 1]);  
code2 = convenc(randi nt(10, 1), trel);
```

The commands below illustrate how to use the final state and initial state arguments when invoking `convenc` repeatedly. Notice that [code3; code4] is the same as the earlier example's output, code1.

```
trel = poly2trellis([5 4], [27 33 0; 0 5 13]);  
msg = randi nt(10, 1, 2, 123);  
% Encode part of msg, recording final state for later use.  
[code3, fstate] = convenc(msg(1:6), trel);  
% Encode the rest of msg, using state as an input argument.  
code4 = convenc(msg(7:10), trel, fstate);
```

### See Also

`vitdec`, `poly2trellis`, `istrellis`

### References

Gitlin, Richard D., Jeremiah F. Hayes, and Stephen B. Weinstein. *Data Communications Principles*. New York: Plenum, 1992.

<b>Purpose</b>	Produce parity-check and generator matrices for cyclic code
<b>Syntax</b>	<pre> parmat = cycl gen(n, pol); parmat = cycl gen(n, pol, opt); [parmat, genmat] = cycl gen(...); [parmat, genmat, k] = cycl gen(...); </pre>
<b>Description</b>	<p>For all syntaxes, the codeword length is <math>n</math> and the message length is <math>k</math>. A polynomial can generate a cyclic code with codeword length <math>n</math> and message length <math>k</math> if and only if the polynomial is a degree-<math>(n-k)</math> divisor of <math>x^n-1</math>. (Over the binary field <math>GF(2)</math>, <math>x^n-1</math> is the same as <math>x^n+1</math>.) This implies that <math>k</math> equals <math>n</math> minus the degree of the generator polynomial.</p> <p><code>parmat = cycl gen(n, pol)</code> produces an <math>(n-k)</math>-by-<math>n</math> parity-check matrix for a systematic binary cyclic code having codeword length <math>n</math>. The row vector <code>pol</code> gives the binary coefficients, in order of ascending powers, of the degree-<math>(n-k)</math> generator polynomial.</p> <p><code>parmat = cycl gen(n, pol, opt)</code> is the same as the syntax above, except that the argument <code>opt</code> determines whether the matrix should be associated with a systematic or nonsystematic code. The values for <code>opt</code> are '<b>system</b>' and '<b>nonsys</b>'.</p> <p><code>[parmat, genmat] = cycl gen(...)</code> is the same as <code>parmat = cycl gen(...)</code> except that it also produces the <math>k</math>-by-<math>n</math> generator matrix <code>genmat</code> that corresponds to the parity-check matrix <code>parmat</code>.</p> <p><code>[parmat, genmat, k] = cycl gen(...)</code> is the same as <code>[parmat, genmat] = cycl gen(...)</code> except that it also returns the message length <math>k</math>.</p>
<b>Examples</b>	<p>The code below produces parity-check and generator matrices for a binary cyclic code with codeword length 7 and message length 4.</p> <pre> pol = cycl poly(7, 4); [parmat, genmat, k] = cycl gen(7, pol) </pre>

```
parmat =
```

1	0	0	1	1	1	0
0	1	0	0	1	1	1
0	0	1	1	1	0	1

```
genmat =
```

1	0	1	1	0	0	0
1	1	1	0	1	0	0
1	1	0	0	0	1	0
0	1	1	0	0	0	1

```
k =
```

4

In the output below, notice that the parity-check matrix is different from `parmat` above, since it corresponds to a nonsystematic cyclic code. In particular, `parmatn` does not have a 3-by-3 identity matrix in its leftmost three columns, as `parmat` does.

```
parmatn = cyclgen(7, cyclpoly(7, 4), 'nonsys')
```

```
parmatn =
```

1	1	1	0	1	0	0
0	1	1	1	0	1	0
0	0	1	1	1	0	1

## See Also

`cyclpoly`, `encode`



**Purpose** Produce generator polynomials for a cyclic code

**Syntax** `pol = cyclpoly(n, k);`  
`pol = cyclpoly(n, k, opt);`

**Description** For all syntaxes, a polynomial is represented as a row containing the coefficients in order of ascending powers.

`pol = cyclpoly(n, k)` returns the row vector representing one nontrivial generator polynomial for a cyclic code having codeword length  $n$  and message length  $k$ .

`pol = cyclpoly(n, k, opt)` searches for one or more nontrivial generator polynomials for cyclic codes having codeword length  $n$  and message length  $k$ . The output `pol` depends on the argument `opt` as shown in the table below.

opt	Significance of pol	Format of pol
'min'	One generator polynomial having the smallest possible weight	The row vector representing the polynomial
'max'	One generator polynomial having the greatest possible weight	The row vector representing the polynomial
'all'	All generator polynomials	A matrix, each row of which represents one such polynomial
a positive integer	All generator polynomials having weight <i>opt</i>	A matrix, each row of which represents one such polynomial

The weight of a binary polynomial is the number of nonzero terms it has. If no generator polynomial satisfies the given conditions, then the output `pol` is empty and an error message is displayed.

**Examples** The first command below produces representations of three generator polynomials for a [15,4] cyclic code. The second command shows that  $1 + x + x^2 + x^3 + x^5 + x^7 + x^8 + x^{11}$  is one such polynomial having the largest number of nonzero terms. The third command shows that no generator polynomial for a [15,4] cyclic code has exactly three nonzero terms.

```
c1 = cyclpoly(15, 4, 'all')

c1 =
    1    1    0    0    0    1    1    0    0    0    1    1
    1    0    0    1    1    0    1    0    1    1    1    1
    1    1    1    1    0    1    0    1    1    0    0    1

c2 = cyclpoly(15, 4, 'max')

c2 =
    1    1    1    1    0    1    0    1    1    0    0    1

c3 = cyclpoly(15, 4, 3)

No generator polynomial satisfies the given constraints.

c3 =

[]
```

**Algorithm** If *opt* is '**min**', '**max**', or omitted, then polynomials are constructed by converting decimal integers to base *p*. Based on the decimal ordering, `gfprimfd` returns the first polynomial it finds that satisfies the appropriate conditions. This algorithm is similar to the one used in `gfprimfd`.

**See Also** `cyclgen`, `encode`

**Purpose** Digital passband demodulator

**Syntax**

```

z = ddemod(y, Fc, Fd, Fs, 'ask/opt', M, num, den);
z = ddemod(y, Fc, Fd, Fs, 'fsk/opt', M);
z = ddemod(y, Fc, Fd, Fs, 'msk');
z = ddemod(y, Fc, Fd, Fs, 'psk/opt', M, num, den);
z = ddemod(y, Fc, Fd, Fs, 'qask/opt', M, num, den);
z = ddemod(y, Fc, Fd, Fs, 'qask/arb/opt', i nphase, quadr, num, den);
z = ddemod(y, Fc, Fd, Fs, 'qask/ci r/opt', numsi g, amp, phs, num, den);
z = ddemod(y, Fc, Fd, [Fs phase], ...);

```

<b>Optional Inputs</b>	<b>Input</b>	<b>Default Value, or Default Behavior if Input is Omitted</b>
	<i>opt</i>	ddemod demaps after demodulating. If the method is ASK, then the algorithm does not use a Costas loop. If the method is FSK, then demodulation is coherent.
	num, den	Omitting these arguments prevents ddemod from using a filter.
	amp	[ 1: length(numsi g) ]
	phs	numsi g*0

**Description** The function ddemod performs digital passband demodulation. The corresponding modulation function is dmod. The table below lists the demodulation schemes that ddemod supports.

Demodulation Scheme	Fifth Input Argument	Where /opt can contain
M-ary amplitude shift keying	'ask/opt'	/nomap; /costas
M-ary frequency shift keying	'fsk/opt'	/noncoherence
M-ary phase shift keying	'psk/opt'	/nomap
Quadrature amplitude shift keying	'qask/opt', 'qask/ci r/opt', or 'qask/arb/opt'	/nomap

The second column of the table indicates in bold type the required portion of the fifth input argument for ddemod. The third column indicates optional flags

that you can append to the fifth argument. The order of optional flags does not matter.

### To Demodulate Without Demapping (ASK, PSK, QASK only)

Ordinarily, the `ddemod` function first demodulates the analog signal it receives and then demaps the demodulated signal in order to recover the digital message signal. The optional `/nomap` flag, appended to the fifth input argument, prevents `ddemod` from demapping. The output is then an analog signal  $x$  whose sampling rate is  $F_s$ . You can use the `demodmap` function to perform the demapping step. The `/nomap` option is not available for FSK or MSK demodulation.

### To Demodulate a Digital Signal (General Information)

The generic syntax `z = ddemod(y, Fc, Fd, Fs, ...)` demodulates the digital message signal  $z$  from a received analog signal  $y$ . After measuring the distance from the received signal to all possible digits in the coding scheme, `ddemod` returns the nearest digit.

$y$  and  $z$  are real matrices whose sizes depend on the demodulation method:

- **(ASK, FSK, MSK methods)** If  $y$  is a vector of length  $n \cdot F_s / F_d$ , then  $z$  is a column vector of length  $n$ . Otherwise, if  $y$  is  $(n \cdot F_s / F_d)$ -by- $m$ , then  $z$  is  $n$ -by- $m$  and each column of  $y$  is processed separately.
- **(PSK, QASK methods)** If  $y$  is  $(n \cdot F_s / F_d)$ -by- $m$ , then  $z$  is  $n$ -by- $2m$ . The odd-numbered columns in  $z$  represent in-phase components and the even-numbered columns represent quadrature components. Each column of  $y$  is processed separately.

The carrier frequency in Hertz is  $F_c$ . The sampling rates in Hertz of  $y$  and  $z$ , respectively, are  $F_s$  and  $F_d$ . (Thus  $1/F_s$  represents the time interval between two consecutive samples in  $y$ , and similarly for  $z$ .) The ratio  $F_s/F_d$  must be a positive integer. The time interval between two decision points is  $1/F_d$ .

The generic syntax `z = ddemod(y, Fc, Fd, [Fs phase], ...)` is the same, except that the fourth input argument is a two-element vector instead of a scalar. The first entry,  $F_s$ , is the sampling rate as described in the paragraph above. The second entry, `phase`, is the initial phase of the carrier signal, measured in radians.

ddemod can use a lowpass filter with sample time  $1/F_s$  while demodulating, in order to filter out the carrier signal. To specify the lowpass filter, include num and den in the list of input arguments. num and den are row vectors that give the coefficients, in *descending* order, of the numerator and denominator of the filter's transfer function. If num is empty, zero, or absent, then the function does not use a filter.

### To Demodulate a Digital Signal (Specific Syntax Information)

`z = ddemod(y, Fc, Fd, Fs, 'ask', M)` implements M-ary amplitude shift keying demodulation. Each entry of `z` is in the range  $[0, M-1]$ .

`z = ddemod(y, Fc, Fd, Fs, 'ask/costas', M)` is the same as the syntax above, except that the algorithm includes a Costas loop

`z = ddemod(y, Fc, Fd, Fs, 'fsk', M, tone)` implements coherent M-ary frequency shift keying demodulation. The optional argument `tone` is the separation between successive frequencies in the modulated signal `z`. The default value of `tone` is `Fd`. Each entry of `z` is in the range  $[0, M-1]$ .

`z = ddemod(y, Fc, Fd, Fs, 'fsk/noncoherence', M, tone)` is the same as the syntax above, except that it uses noncoherent demodulation.

`z = ddemod(y, Fc, Fd, Fs, 'msk')` implements minimum shift keying demodulation. Each entry of `z` is either 0 or 1. The separation between the two frequencies is  $F_d/2$ .

`z = ddemod(y, Fc, Fd, Fs, 'psk', M)` implements M-ary correlation phase shift keying demodulation. Each entry of `z` is in the range  $[0, M-1]$ .

`z = ddemod(y, Fc, Fd, Fs, 'qask', M)` implements M-ary quadrature amplitude shift keying demodulation with a square signal constellation. The table below

shows the maximum among in-phase and quadrature coordinates of constellation points, for several small values of M.

M	Maximum of Coordinates of Constellation Points	M	Maximum of Coordinates of Constellation Points
2	1	32	5
4	1	64	7
8	3 (quadrature maximum is 1)	128	11
16	3	256	15

**Note** To see how symbols are mapped to the constellation points, generate a square constellation plot using `qaskenco(M)`.

`z = ddemod(y, Fc, Fd, Fs, 'qask/arb', inphase, quadr)` implements quadrature amplitude shift keying demodulation, with a signal constellation that you define using the vectors `inphase` and `quadr`. The signal constellation point for the  $k$ th message has in-phase component `inphase(k+1)` and quadrature component `quadr(k+1)`.

`z = ddemod(y, Fc, Fd, Fs, 'qask/cir', numsig, amp, phs)` implements quadrature amplitude shift keying demodulation with a circular signal constellation. `numsig`, `amp`, and `phs` are vectors of the same length. The entries in `numsig` and `amp` must be positive. If  $k$  is an integer in the range  $[1, \text{length}(\text{numsig})]$ , then `amp(k)` is the radius of the  $k$ th circle, `numsig(k)` is the number of constellation points on the  $k$ th circle, and `phs(k)` is the phase of the first constellation point plotted on the  $k$ th circle. All points on the  $k$ th circle are evenly spaced. If you omit `phs`, then its default value is `numsig*0`. If you omit `amp`, then its default value is `[1: length(numsig)]`.

---

**Note** To see how symbols are mapped to the constellation points, generate a labeled circle constellation plot using `apkconst (numsi g, amp, phs, 'n')`.

---

## Examples

This example mimics the one in the section “Simple Digital Modulation Example” on page 2-74 but uses passband simulation. It generates a random digital signal, modulates it using `dmod`, and adds noise. Then it demodulates the noisy signal and computes the symbol error rate. The `ddemod` function demodulates the analog signal `y` and then demaps to produce the digital signal `z`.

Important differences between this example and the original baseband example are the explicit reference to the carrier signal frequency `Fc` and the fact that `y` and `ynoi sy` are real, not complex. For variety, this example uses ASK instead of PSK, as well as a different sampling rate `Fd`.

```
M = 16; % Use 16-ary modulation.
Fc = 10; % Carrier signal frequency is 10 Hz.
Fd = 1; % Sampling rates of original and modulated signals
Fs = 50; % are 1 and 50, respectively (samples per second).
x = randint(100, 1, M); % Random digital message
% Use M-ary PSK modulation to produce y.
y = dmod(x, Fc, Fd, Fs, 'ask', M);
% Add some Gaussian noise.
ynoi sy = y + .01*randn(Fs/Fd*100, 1);
% Demodulate y to recover the message.
z = ddemod(ynoi sy, Fc, Fd, Fs, 'ask', M);
s = symerr(x, z) % Check symbol error rate.
```

`s =`

`0`

## See Also

`dmod`, `amod`, `ademod`, `dmodce`, `ddemodce`, `demodmap`, `modmap`, `eyediagram`, `scatterplot`

# ddemodce

**Purpose** Digital baseband demodulator

**Syntax**

```
z = ddemodce(y, Fd, Fs, 'ask/opt', M, num, den);
z = ddemodce(y, Fd, Fs, 'fsk/opt', M);
z = ddemodce(y, Fd, Fs, 'msk');
z = ddemodce(y, Fd, Fs, 'psk/opt', M, num, den);
z = ddemodce(y, Fd, Fs, 'qask/opt', M, num, den);
z = ddemodce(y, Fd, Fs, 'qask/arb/opt', inphase, quadr, num, den);
z = ddemodce(y, Fd, Fs, 'qask/cir/opt', numsig, amp, phs, num, den);
z = ddemodce(y, Fd, [Fs phase], ...);
```

Optional Inputs	Input	Default Value, or Default Behavior if Input is Omitted
	opt	ddemodce demaps after demodulating. If the method is ASK, then the algorithm does not use a Costas loop. If the method is FSK, then demodulation is coherent.
	num, den	Omitting these arguments prevents ddemodce from using a filter.
	amp	[ 1: length(numsig) ]
	phs	numsig*0

**Description** The function ddemodce performs digital baseband demodulation. The corresponding modulation function is dmodce. The table below lists the demodulation schemes that ddemodce supports.

Demodulation Scheme	Fourth Input Argument	Where /opt can contain
M-ary amplitude shift keying	'ask/opt'	/nomap; /costas
M-ary frequency shift keying	'fsk/opt'	/noncoherence
M-ary phase shift keying	'psk/opt'	/nomap
Quadrature amplitude shift keying	'qask/opt', 'qask/cir/opt', or 'qask/arb/opt'	/nomap

The second column of the table indicates in bold type the required portion of the fourth input argument for ddemodce. The third column indicates optional



flags that you can append to the fourth argument. The order of optional flags does not matter.

### To Demodulate Without Demapping (ASK, PSK, QASK only)

Ordinarily, the `ddemodce` function first demodulates the analog signal it receives and then demaps the demodulated signal in order to recover the digital message signal. The optional `/nomap` flag, appended to the fourth input argument, prevents `ddemodce` from demapping. The output is then an analog signal  $z$  whose sampling rate is  $F_s$ . The size of  $z$  depends on the size of  $y$  and the demodulation method:

- **(ASK method)**  $z$  has the same size as  $y$ .
- **(PSK and QASK methods)** If  $y$  is a vector of length  $n$ , then  $z$  is an  $n$ -by-2 matrix. Otherwise, if  $y$  is  $n$ -by- $m$ , then  $z$  is  $n$ -by- $2m$  and each column of  $y$  is processed separately. In either case, the odd-numbered columns in  $z$  represent in-phase components and the even-numbered columns represent quadrature components.

You can use the `demodmap` function to perform the demapping step. The `/nomap` option is not available for FSK or MSK demodulation.

### To Demodulate a Digital Signal (General Information)

The generic syntax `z = ddemodce(y, Fd, Fs, ...)` demodulates the digital message signal  $z$  from a received analog signal  $y$ . After measuring the distance from the received signal to all possible digits in the coding scheme, `ddemodce` returns the nearest digit.

$y$  is a complex matrix and  $z$  is a real matrix. The sizes of  $y$  and  $z$  depend on the demodulation method:

- **(ASK, FSK, MSK methods)** If  $y$  is a vector of length  $n \cdot F_s / F_d$ , then  $z$  is a column vector of length  $n$ . Otherwise, if  $y$  is  $(n \cdot F_s / F_d)$ -by- $m$ , then  $z$  is  $n$ -by- $m$  and each column of  $y$  is processed separately.
- **(PSK, QASK methods)** If  $y$  is  $(n \cdot F_s / F_d)$ -by- $m$ , then  $z$  is  $n$ -by- $2m$ . The odd-numbered columns in  $z$  represent in-phase components and the even-numbered columns represent quadrature components. Each column of  $y$  is processed separately.

The sampling rates in Hertz of  $y$  and  $z$ , respectively, are  $F_s$  and  $F_d$ . (Thus  $1/F_s$  represents the time interval between two consecutive samples in  $y$ , and

similarly for  $z$ .) The ratio  $F_s/F_d$  must be a positive integer. The time interval between two decision points is  $1/F_d$ .

The generic syntax `z = ddemodce(y, Fd, [Fs phase], ...)` is the same, except that the third input argument is a two-element vector instead of a scalar. The first entry,  $F_s$ , is the sampling rate as described in the paragraph above. The second entry, *phase*, is the initial phase of the carrier signal, measured in radians.

To use a lowpass filter in conjunction with ASK, PSK, or QASK demodulation, include *num* and *den* in the list of input arguments. *num* and *den* are row vectors that give the coefficients, in *descending* order, of the numerator and denominator of the filter's transfer function. If *num* is empty, zero, or absent, then `ddemodce` does not use a filter.

### To Demodulate a Digital Signal (Specific Syntax Information)

`z = ddemodce(y, Fd, Fs, 'ask', M)` implements  $M$ -ary amplitude shift keying demodulation. Each entry of  $z$  is in the range  $[0, M-1]$ .

`z = ddemodce(y, Fd, Fs, 'ask/costas', M)` is the same as the syntax above, except that the algorithm includes a Costas loop

`z = ddemodce(y, Fd, Fs, 'fsk', M, tone)` implements coherent  $M$ -ary frequency shift keying demodulation. The optional argument *tone* is the separation between successive frequencies in the modulated signal  $z$ . The default value of *tone* is  $F_d$ . Each entry of  $z$  is in the range  $[0, M-1]$ .

`z = ddemodce(y, Fd, Fs, 'fsk/noncoherence', M, tone)` is the same as the syntax above, except that it uses noncoherent demodulation.

`z = ddemodce(y, Fd, Fs, 'msk')` implements minimum shift keying demodulation. Each entry of  $z$  is either 0 or 1. The separation between the two frequencies is  $F_d/2$ .

`z = ddemodce(y, Fd, Fs, 'psk', M)` implements  $M$ -ary correlation phase shift keying demodulation. Each entry of  $z$  is in the range  $[0, M-1]$ .

`z = ddemodce(y, Fd, Fs, 'qask', M)` implements  $M$ -ary quadrature amplitude shift keying demodulation with a square signal constellation. The table below

shows the maximum among in-phase and quadrature coordinates of constellation points, for several small values of  $M$ .

$M$	Maximum of Coordinates of Constellation Points	$M$	Maximum of Coordinates of Constellation Points
2	1	32	5
4	1	64	7
8	3 (quadrature maximum is 1)	128	11
16	3	256	15

**Note** To see how symbols are mapped to the constellation points, generate a square constellation plot using `qaskenco(M)`.

`z = ddemodce(y, Fd, Fs, 'qask/arb', inphase, quadr)` implements quadrature amplitude shift keying demodulation, with a signal constellation that you define using the vectors `inphase` and `quadr`. The signal constellation point for the  $k$ th message has in-phase component `inphase(k+1)` and quadrature component `quadr(k+1)`.

`z = ddemodce(y, Fd, Fs, 'qask/cir', numsig, amp, phs)` implements quadrature amplitude shift keying demodulation with a circular signal constellation. `numsig`, `amp`, and `phs` are vectors of the same length. The entries in `numsig` and `amp` must be positive. If  $k$  is an integer in the range  $[1, \text{length}(\text{numsig})]$ , then `amp(k)` is the radius of the  $k$ th circle, `numsig(k)` is the number of constellation points on the  $k$ th circle, and `phs(k)` is the phase of the first constellation point plotted on the  $k$ th circle. All points on the  $k$ th circle are evenly spaced. If you omit `phs`, then its default value is `numsig*0`. If you omit `amp`, then its default value is `[1: length(numsig)]`.

## ddemodce

---

---

**Note** To see how symbols are mapped to the constellation points, generate a labeled circle constellation plot using `apkconst (numsig, amp, phs, 'n')`.

---

### See Also

`dmodce`, `amodce`, `ademodce`, `dmod`, `ddemod`, `demodmap`, `eyediagram`, `scatterplot`

**Purpose** Convert decimal numbers to binary vectors

**Syntax**

```
b = de2bi(d);
b = de2bi(d, n);
b = de2bi(d, n, p);
b = de2bi(d, [], p);
b = de2bi(d, ..., flag)
```

**Description** `b = de2bi(d)` converts a nonnegative decimal integer `d` to a binary row vector. If `d` is a vector, then the output `b` is a matrix, each row of which is the binary form of the corresponding element in `d`. If `d` is a matrix, then `de2bi` treats it like the vector `d(:)`.

---

**Note** By default, `de2bi` uses the first column of `b` as the *lowest-order* digit.

---

`b = de2bi(d, n)` is the same as `b = de2bi(d)`, except that its output has `n` columns, where `n` is a positive integer. An error occurs if the binary representations would require more than `n` digits. If necessary, the binary representation of `d` is padded with extra zeros.

`b = de2bi(d, n, p)` converts a nonnegative decimal integer `d` to a base-`p` row vector, where `p` is an integer greater than or equal to two. The first column of `b` is the *lowest* base-`p` digit. `b` is padded with extra zeros if necessary, so that it has `n` columns, where `n` is a positive integer. An error occurs if the base-`p` representations would require more than `n` digits. If `d` is a nonnegative decimal vector, then the output `b` is a matrix, each row of which is the (possibly zero-padded) base-`p` form of the corresponding element in `d`. If `d` is a matrix, then `de2bi` treats it like the vector `d(:)`.

`b = de2bi(d, [], p)` specifies the base `p` but not the number of columns.

`b = de2bi(d, ..., flag)` uses the string `flag` to determine whether the first column of `b` contains the lowest-order or highest-order digits. Values for `flag` are **'right-msb'** and **'left-msb'**. The value **'right-msb'** produces the default behavior.

Examples

The code below counts to ten in decimal and binary.

```
d = (1:10)';
b = de2bi(d);
disp('      Dec          Bi nary      ')
disp('      -----  -')
disp([d, b])
```

The output is below.

Dec	Bi nary				
	-----				
1	1	0	0	0	0
2	0	1	0	0	0
3	1	1	0	0	0
4	0	0	1	0	0
5	1	0	1	0	0
6	0	1	1	0	0
7	1	1	1	0	0
8	0	0	0	1	0
9	1	0	0	1	0
10	0	1	0	1	0

The command below shows how de2bi pads its output with zeros.

```
bb = de2bi([3 9],5) % Zero-padding the output

bb =

     1     1     0     0     0
     1     0     0     1     0
```

The command below shows how to convert a decimal integer to base three without specifying the number of columns in the output matrix.

```
t = de2bi(12,[],3) % Convert 12 to base 3.

t =

     0     1     1
```

See Also

bi2de

Purpose	Block decoder	
Syntax	<pre>msg = decode(code, n, k, 'hamming/format', pri mpoly); msg = decode(code, n, k, 'linear/format', genmat, trt); msg = decode(code, n, k, 'cyclic/format', genpoly, trt); msg = decode(code, n, k, 'bch/format', errorcorr, pri mpoly); msg = decode(code, n, k, 'rs/format', field); msg = decode(code, n, k); [msg, err] = decode(...); [msg, err, ccode] = decode(...); [msg, err, ccode, cerr] = decode(...);</pre>	
Optional Inputs	Input	Default Value
	<i>format</i>	<b>binary</b>
	<i>pri mpoly</i>	<i>gfprimdf</i> ( <i>m</i> ) where $n = 2^m - 1$
	<i>genpoly</i>	<i>cycl poly</i> ( <i>n</i> , <i>k</i> )
	<i>trt</i>	Uses <i>syndtable</i> to create the syndrome decoding table associated with the method's parity-check matrix.
Description	For All Syntaxes	
	The decode function aims to recover messages that were encoded using an error-correction coding technique. The technique and the defining parameters must match those that were used to encode the original signal.	
	The “For All Syntaxes” section on the reference page for the encode function explains the meanings of <i>n</i> and <i>k</i> , the possible values of <i>format</i> , and the possible formats for <i>code</i> and <i>msg</i> . You should be familiar with the conventions described there before reading the rest of this section. Using the decode function with an input argument <i>code</i> that was <i>not</i> created by the encode function may cause errors.	
	For Specific Syntaxes	
	<i>msg</i> = <i>decode</i> ( <i>code</i> , <i>n</i> , <i>k</i> , 'hamming/format', <i>pri mpoly</i> ) decodes <i>code</i> using the Hamming method. For this syntax, <i>n</i> must have the form $2^m - 1$ for some integer <i>m</i> greater than or equal to 3, and <i>k</i> must equal <i>n-m</i> . <i>pri mpoly</i> is a row	

vector that gives the binary coefficients, in order of ascending powers, of the primitive polynomial for  $GF(2^m)$  that is used in the encoding process. The default value of `primpoly` is `gfprimdf(m)`. The decoding table that the function uses to correct a single error in each codeword is `syndtable(hammgen(m))`.

`msg = decode(code, n, k, 'linear/format', genmat, trt)` decodes `code`, which is a linear block code determined by the  $k$ -by- $n$  generator matrix `genmat`. `genmat`, a  $k$ -by- $n$  matrix, is required as input. `decode` tries to correct errors using the decoding table `trt`, where `trt` is a  $2^{n-k}$ -by- $n$  matrix.

`msg = decode(code, n, k, 'cyclic/format', genpoly, trt)` decodes the cyclic code `code` and tries to correct errors using the decoding table `trt`, where `trt` is a  $2^{n-k}$ -by- $n$  matrix. `genpoly` is a row vector that gives the coefficients, in order of ascending powers, of the binary generator polynomial of the code. The default value of `genpoly` is `cyclpoly(n, k)`. By definition, the generator polynomial for an  $[n, k]$  cyclic code must have degree  $n-k$  and must divide  $x^n-1$ .

`msg = decode(code, n, k, 'bch/format', errorcorr, primpoly)` decodes `code` using the BCH method. `primpoly` is a row vector that gives the coefficients, in order of ascending powers, of the primitive polynomial for  $GF(2^m)$  that will be used during processing. The default value of `primpoly` is `gfprimdf(m)`. For this syntax,  $n$  must have the form  $2^m-1$  for some integer  $m$  greater than or equal to 3.  $k$  and `errorcorr` must be a valid message length and error-correction capability, respectively, as reported in the second and third columns of a row of `params` in the command

```
params = bchpoly(n)
```

`msg = decode(code, n, k, 'rs/format', field)` decodes `code` using the Reed-Solomon method.  $n$  must have the form  $2^m-1$  for some integer  $m$  greater than or equal to 3. `field` is a matrix that lists all elements of  $GF(2^m)$  in the format described in "List of All Elements of a Galois Field" on page 2-91. The default value of `field` is `gftuple([-1:2^m-2], m)`.

`msg = decode(code, n, k)` is the same as  
`msg = decode(code, n, k, 'hamming/binary')`.

`[msg, err] = decode(...)` returns a column vector `err` that gives information about error correction. If the code is a convolutional code, then `err` contains the



metric calculations used in the decoding decision process. For other types of codes, a nonnegative integer in the  $r$ th row of `err` (or the  $r$ th row of `vec2mat(err, k)` if code is a column vector) indicates the number of errors corrected in the  $r$ th *message* word; a negative integer indicates that there are more errors in the  $r$ th word than can be corrected.

`[msg, err, ccode] = decode(...)` returns the corrected code in `ccode`.

`[msg, err, ccode, cerr] = decode(...)` returns a column vector `cerr` whose meaning depends on the format of code:

- If code is a binary vector, then a nonnegative integer in the  $r$ th row of `vec2mat(cerr, n)` indicates the number of errors corrected in the  $r$ th *codeword*; a negative integer indicates that there are more errors in the  $r$ th codeword than can be corrected.
- If code is not a binary vector, then `cerr = err`.

## Examples

On the reference page for `encode`, some of the example code illustrates the use of the `decode` function.

The example below illustrates the use of `err` and `cerr` when the coding method is not convolutional code and the code is a binary vector. The script encodes two five-bit messages using BCH code. Each codeword has fifteen bits. Errors are added to the first two bits of the first codeword and the first bit of the second codeword. Then `decode` is used to recover the original message. As a result, the errors are corrected. `err` is the same size as `msg` and `cerr` is the same size as `code`. `err` reflects the fact that the first *message* was recovered after correcting two errors, while the second message was recovered after correcting one error. `cerr` reflects the fact that the first *codeword* was decoded after correcting two errors, while the second codeword was decoded after correcting one error.

```
m = 4; n = 2^m-1; % Codeword length is 15.
k = 5; % Valid message length for BCH code when n = 15
t = 3; % Corresponding error-correction capability
msg = ones(10, 1); % Two messages, five bits each
code = encode(msg, n, k, 'bch'); % Encode the message.
% Now place two errors in first word and one error
% in the second word. Create errors by reversing bits.
noisycode = code;
noisycode(1:2) = bitxor(noisycode(1:2), [1 1]');
```

## decode

---

```
noisycode(16) = bitxor(noisycode(16), 1);  
% Decode and try to correct the errors.  
[newmsg, err, ccode, cerr] = decode(noisycode, n, k, 'bch', t);  
disp('Transpose of err is'); disp(err')  
disp('Transpose of cerr is'); disp(cerr')
```

The output is below.

Transpose of err is

2 2 2 2 2 1 1 1 1 1

Transpose of cerr is

Columns 1 through 12

2 2 2 2 2 2 2 2 2 2 2 2

Columns 13 through 24

2 2 2 1 1 1 1 1 1 1 1 1

Columns 25 through 30

1 1 1 1 1 1

### Algorithm

Depending on the decoding method, decode relies on such lower-level functions as `hammgen`, `syndtable`, `cyclgen`, `bchdeco`, and `rsdeco`.

### See Also

`encode`, `hammgen`, `syndtable`, `cyclpoly`, `cyclgen`, `bchpoly`, `bchdeco`, `rspoly`, `rsdeco`, `rsdecode`, `vitdec`

**Purpose** Demap a digital message from a demodulated signal

**Syntax**

```
z = demodmap(x, Fd, Fs, 'ask', M);  
z = demodmap(x, Fd, Fs, 'fsk', M, tone);  
z = demodmap(x, Fd, Fs, 'msk');  
z = demodmap(x, Fd, Fs, 'psk', M);  
z = demodmap(x, Fd, Fs, 'qask', M);  
z = demodmap(x, Fd, Fs, 'qask/arb', inphase, quadr);  
z = demodmap(x, Fd, Fs, 'qask/cir', numsig, amp, phs);  
z = demodmap(x, [Fd offset], Fs, ...)
```

Optional Inputs	Input	Default Value
	tone	Fd
	amp	[ 1: length(numsig) ]
	phs	numsig*0

**Description** The digital demodulation process consists of two steps: demodulating an analog signal and demapping the demodulated signal to a digital signal. You can perform the first step using `ademod`, `ademodce`, or your own custom demodulator. The function `demodmap` performs the second step. The table below lists the demodulation schemes that `demodmap` supports.

Demodulation Scheme	Fourth Input Argument
M-ary amplitude shift keying	'ask'
M-ary frequency shift keying	'fsk'
Minimum shift keying	'msk'
M-ary phase shift keying	'psk'
Quadrature amplitude shift keying	'qask', 'qask/cir', or 'qask/arb'

### To Demap a Digital Signal (General Information)

The generic syntax `z = demodmap(x, Fd, Fs, ...)` demaps the digital message signal `z` from a received analog signal `x`. After measuring the distance from the

received signal to all possible digits in the coding scheme, the demapper returns the nearest digit.

$x$  is a matrix. The sizes of  $x$  and  $z$  depend on the demodulation method:

- **(ASK, FSK, MSK methods)** If  $x$  is a vector of length  $n \cdot F_s / F_d$ , then  $z$  is a column vector of length  $n$ . Otherwise, if  $x$  is  $(n \cdot F_s / F_d)$ -by- $m$ , then  $z$  is  $n$ -by- $m$  and each column of  $x$  is processed separately.
- **(PSK, QASK methods)**  $x$  must have an even number of columns. The odd-numbered columns in  $x$  represent in-phase components and the even-numbered columns represent quadrature components. Each *pair* of columns of  $x$  is processed separately. If  $x$  is  $(n \cdot F_s / F_d)$ -by- $2m$ , then  $z$  is  $n$ -by- $m$ .

The sampling rates in Hertz of  $x$  and  $z$ , respectively, are  $F_s$  and  $F_d$ . (Thus  $1/F_s$  represents the time interval between two consecutive samples in  $x$ , and similarly for  $z$ .) The ratio  $F_s/F_d$  must be a positive integer. The time interval between two decision points is  $1/F_d$ .

To shift the decision times ahead by the integer offset, use the alternative syntax

```
z = demodmap(x, [Fd offset], ...)
```

instead of the demapping syntaxes listed in this section and the next. The default decision offset is 0.

## To Demap a Digital Signal (Specific Syntax Information)

`z = demodmap(x, Fd, Fs, 'ask', M)` demaps from an  $M$ -ary amplitude shift keying signal constellation. Each entry of  $z$  is in the range  $[0, M-1]$ .

`z = demodmap(x, Fd, Fs, 'fsk', M, tone)` demaps using the coherent  $M$ -ary frequency shift keying method. The optional argument `tone` is the separation between successive frequencies in the modulated signal  $x$ . The default value of `tone` is  $F_d$ . Each entry of  $z$  is in the range  $[0, M-1]$ .

`z = demodmap(x, Fd, Fs, 'msk')` demaps using the minimum shift keying method. Each entry of  $z$  is either 0 or 1. The separation between the two frequencies is  $F_d/2$ .

$z = \text{demodmap}(x, F_d, F_s, ' \text{psk}' , M)$  demaps from an  $M$ -ary phase shift keying signal constellation. Each entry of  $z$  is in the range  $[0, M-1]$ .

$z = \text{demodmap}(x, F_d, F_s, ' \text{qask}' , M)$  demaps from an  $M$ -ary quadrature amplitude shift keying square signal constellation. The table below shows the maximum among in-phase and quadrature coordinates of constellation points, for several small values of  $M$ .

M	Maximum of Coordinates of Constellation Points	M	Maximum of Coordinates of Constellation Points
2	1	32	5
4	1	64	7
8	3 (quadrature maximum = 1)	128	11
16	3	256	15

**Note** To see how symbols are mapped to the constellation points, generate a square constellation plot using `qaskenco(M)`.

$z = \text{demodmap}(x, F_d, F_s, ' \text{qask/arb}' , i_{\text{nphase}}, \text{quadr})$  demaps from a quadrature amplitude shift keying signal constellation that you define using the vectors  $i_{\text{nphase}}$  and  $\text{quadr}$ . The signal constellation point for the  $k$ th message has in-phase component  $i_{\text{nphase}}(k+1)$  and quadrature component  $\text{quadr}(k+1)$ .

$z = \text{demodmap}(x, F_d, F_s, ' \text{qask/cir}' , \text{numsi\_g}, \text{amp}, \text{phs})$  demaps from a quadrature amplitude shift keying circular signal constellation.  $\text{numsi\_g}$ ,  $\text{amp}$ , and  $\text{phs}$  are vectors of the same length. The entries in  $\text{numsi\_g}$  and  $\text{amp}$  must be positive. If  $k$  is an integer in the range  $[1, \text{length}(\text{numsi\_g})]$ , then  $\text{amp}(k)$  is the radius of the  $k$ th circle,  $\text{numsi\_g}(k)$  is the number of constellation points on the  $k$ th circle, and  $\text{phs}(k)$  is the phase of the first constellation point plotted on the  $k$ th circle. All points on the  $k$ th circle are evenly spaced. If you omit  $\text{phs}$ , then

its default value is `numsig*0`. If you omit `amp`, then its default value is `[1:length(numsig)]`.

---

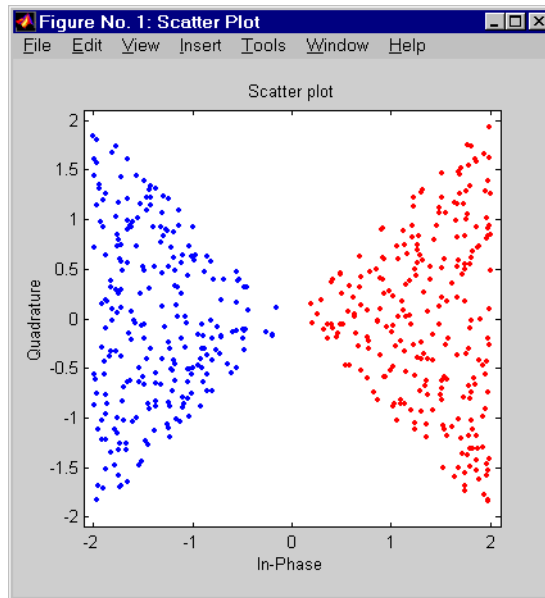
**Note** To see how symbols are mapped to the constellation points, generate a labeled circle constellation plot using `apkconst(numsig, amp, phs, 'n')`.

---

## Examples

The script below suggests which regions in the in-phase/quadrature plane are associated with different digits. It demaps random points, looks for points that were demapped to the digits 0 and 2, and plots those points in red and blue, respectively. The horizontal axis shows in-phase components and the vertical axis shows quadrature components.

```
% Construct [in-phase, quadrature] for random points.
x = 4*(rand(1000, 2) - 1/2);
% Demap to a digital signal, using 4-PSK method.
y = demodmap(x, 1, 1, 'psk', 4);
red = find(y==0); % Indices of points that mapped to the digit 0
h = scatterplot(x(red,:), 1, 0, 'r. '); hold on % Plot in red.
blue = find(y==2); % Indices of points that mapped to the digit 2
scatterplot(x(blue,:), 1, 0, 'b. ', h); hold off % Plot in blue.
```

**See Also**

modmap, ddemod, ddemodce, ademod, ademodce, eyediagram, scatterplot

# dmod

**Purpose** Digital passband modulator

**Syntax**

```
y = dmod(x, Fc, Fd, Fs, 'method/nomap' ... );
y = dmod(x, Fc, Fd, Fs, 'ask', M);
y = dmod(x, Fc, Fd, Fs, 'fsk', M, tone);
y = dmod(x, Fc, Fd, Fs, 'msk' );
y = dmod(x, Fc, Fd, Fs, 'psk', M);
y = dmod(x, Fc, Fd, Fs, 'qask', M);
y = dmod(x, Fc, Fd, Fs, 'qask/arb', inphase, quadr);
y = dmod(x, Fc, Fd, Fs, 'qask/cir', numsig, amp, phs);
y = dmod(x, Fc, Fd, [Fs phase], ... );
[y, t] = dmod(... );
```

Optional Inputs	Input	Default Value
	tone	Fd
	amp	[ 1:length(numsig) ]
	phs	numsig*0

**Description** The function dmod performs digital passband modulation and some related tasks. The corresponding demodulation function is ddemod. The table below lists the modulation schemes that dmod supports.

Modulation Scheme	Fifth Input Argument
M-ary amplitude shift keying	'ask'
M-ary frequency shift keying	'fsk'
Minimum shift keying	'msk'
M-ary phase shift keying	'psk'
Quadrature amplitude shift keying	'qask', 'qask/cir', or 'qask/arb'

**To Avoid the Mapping Process**  
Ordinarily, the dmod function first maps the digital message signal to an analog signal and then modulates the analog signal. The generic syntax



```
y = dmod(x, Fc, Fd, Fs, 'method/nomap' . . .)
```

uses the **nomap** flag to tell dmod that the digital message has already been mapped to an analog signal  $x$  whose sampling rate is  $F_s$ . As a result, dmod skips its usual mapping step. You can use the `modmap` function to perform the mapping step. In this generic syntax, *method* is one of the seven values listed in the table above and the other variables are as in the next section.

### To Modulate a Digital Signal (General Information)

The generic syntax  $y = \text{dmod}(x, F_c, F_d, F_s, \dots)$  modulates the digital message signal that  $x$  represents.  $x$  is a matrix of nonnegative integers. If  $x$  is a vector of length  $n$ , then  $y$  is a vector of length  $n \cdot F_s / F_d$ . Otherwise, if  $x$  is  $n$ -by- $m$ , then  $y$  is  $(n \cdot F_s / F_d)$ -by- $m$  and each column of  $x$  is processed separately.

$F_c$  is the carrier frequency in Hertz. The sampling rates in Hertz of  $x$  and  $y$ , respectively, are  $F_d$  and  $F_s$ . (Thus  $1/F_d$  represents the time interval between two consecutive samples in  $x$ , and similarly for  $y$ .) The ratio  $F_s/F_d$  must be a positive integer. For best results, use values such that  $F_s > F_c > F_d$ . The initial phase of the carrier signal is zero.

The generic syntax  $y = \text{dmod}(x, F_c, F_d, [F_s \text{ phase}], \dots)$  is the same, except that the fourth input argument is a two-element vector instead of a scalar. The first entry,  $F_s$ , is the sampling rate as described in the paragraph above. The second entry, *phase*, is the initial phase of the carrier signal, measured in radians.

### To Modulate a Digital Signal (Specific Syntax Information)

$y = \text{dmod}(x, F_c, F_d, F_s, \text{'ask'}, M)$  performs  $M$ -ary amplitude shift keying modulation. Each entry of  $x$  must be in the range  $[0, M-1]$ . The maximum value of the modulated signal is 1.

$y = \text{dmod}(x, F_c, F_d, F_s, \text{'fsk'}, M, \text{tone})$  performs  $M$ -ary frequency shift keying modulation. Each entry of  $x$  must be in the range  $[0, M-1]$ . The optional argument *tone* is the separation between successive frequencies in the modulated signal  $y$ . The default value of *tone* is  $F_d$ . The maximum value of  $y$  is 1.

$y = \text{dmod}(x, F_c, F_d, F_s, \text{'msk'})$  performs minimum shift keying modulation. Each entry of  $x$  is either 0 or 1. The maximum value of  $y$  is 1.

`y = dmod(x, Fc, Fd, Fs, 'psk', M)` performs M-ary phase shift keying modulation. Each entry of `x` must be in the range `[0, M-1]`. The maximum value of `y` is 1.

`y = dmod(x, Fc, Fd, Fs, 'qask', M)` performs M-ary quadrature amplitude shift keying modulation with a square signal constellation. The table below shows the maximum value of `y`, for several small values of `M`.

M	Maximum Value of y	M	Maximum Value of y
2	1	32	5
4	1	64	7
8	3	128	11
16	3	256	15

**Note** To see how symbols are mapped to the constellation points, generate a square constellation plot using `qaskenco(M)`.

`y = dmod(x, Fc, Fd, Fs, 'qask/arb', inphase, quadr)` performs quadrature amplitude shift keying modulation, with a signal constellation that you define using the vectors `inphase` and `quadr`. The constellation point for the  $k$ th message has in-phase component `inphase(k+1)` and quadrature component `quadr(k+1)`.

`y = dmod(x, Fc, Fd, Fs, 'qask/cir', numsig, amp, phs)` performs quadrature amplitude shift keying modulation with a circular signal constellation. `numsig`, `amp`, and `phs` are vectors of the same length. The entries in `numsig` and `amp` must be positive. If  $k$  is an integer in the range `[1, length(numsig)]`, then `amp(k)` is the radius of the  $k$ th circle, `numsig(k)` is the number of constellation points on the  $k$ th circle, and `phs(k)` is the phase of the first constellation point plotted on the  $k$ th circle. All points on the  $k$ th circle are evenly spaced. If you omit `phs`, then its default value is `numsig*0`. If you omit `amp`, then its default value is `[1: length(numsig)]`.

---

**Note** To see how symbols are mapped to the constellation points, generate a labeled circle constellation plot using `apkconst (numsi g, amp, phs, 'n')`.

---

`[y, t] = dmod(...)` returns the computation time in `t`. `t` is a vector whose length is the number of rows of `y`.

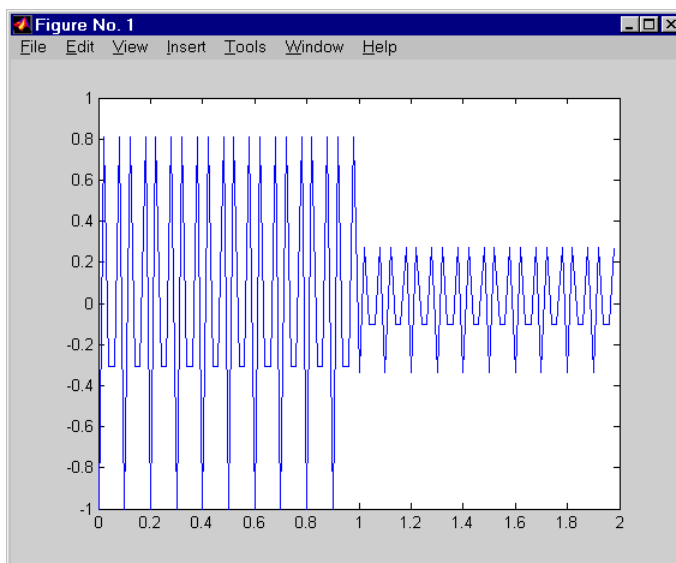
## Examples

An example on the reference page for `ddemod` uses `dmod`. Also, the code below shows the waveforms used to communicate the digits 0 and 1 using 4-ASK modulation. Notice that the `dmod` command has two output arguments. The second output, `t`, is used to scale the horizontal axis in the plot.

```

Fc = 20; Fd = 10; Fs = 50;
M = 4; % Use 4-ASK modulation.
x = ones(Fd, 1)*[0 1]; x=x(:);
% Modulate, keeping track of time.
[y, t] = dmod(x, Fc, Fd, Fs, 'ask', M);
plot(t, y) % Plot signal versus time.

```



## See Also

`ddemod`, `dmodce`, `ddemodce`, `amod`, `amodce`

# dmodce

**Purpose** Digital baseband modulator

**Syntax**

```
y = dmodce(x, Fd, Fs, 'method/nomap' . . . );
y = dmodce(x, Fd, Fs, 'ask', M);
y = dmodce(x, Fd, Fs, 'fsk', M, tone);
y = dmodce(x, Fd, Fs, 'msk' );
y = dmodce(x, Fd, Fs, 'psk', M);
y = dmodce(x, Fd, Fs, 'qask', M);
y = dmodce(x, Fd, Fs, 'qask/arb', inphase, quadr);
y = dmodce(x, Fd, Fs, 'qask/cir', numsig, amp, phs);
y = dmodce(x, Fd, [Fs phase], . . . );
```

Optional Inputs	Input	Default Value
	tone	Fd
	amp	[ 1:length(numsig) ]
	phs	numsig*0

**Description** The function dmodce performs digital baseband modulation and some related tasks. The corresponding demodulation function is ddemodce. The table below lists the modulation schemes that dmodce supports.

Modulation Scheme	Fourth Input Argument
M-ary amplitude shift keying	'ask'
M-ary frequency shift keying	'fsk'
Minimum shift keying	'msk'
M-ary phase shift keying	'psk'
Quadrature amplitude shift keying	'qask', 'qask/cir', or 'qask/arb'

**To Modulate Without Mapping**  
Ordinarily, the dmodce function first maps the digital message signal to an analog signal and then modulates the analog signal. The generic syntax

```
y = dmodce(x, Fd, Fs, 'method/nomap' . . . )
```

uses the `/nomap` flag to tell `dmodce` that the digital message has already been mapped to an analog signal `x` whose sampling rate is `Fs`. As a result, `dmodce` skips its usual mapping step. You can use the `modmap` function to perform the mapping step. In this generic syntax, *method* is one of the seven values listed in the table above, and the other variables are as in the next section.

### To Modulate a Digital Signal (General Information)

The generic syntax `y = dmodce(x, Fd, Fs, ...)` modulates the digital message signal that `x` represents. `x` is a matrix of nonnegative integers. If `x` is a vector of length `n`, then `y` is a vector of length  $n \cdot F_s / F_d$ . Otherwise, if `x` is `n`-by-`m`, then `y` is  $(n \cdot F_s / F_d)$ -by-`m` and each column of `x` is processed separately. Since `dmodce` implements baseband simulation, the entries of `y` are *complex*.

The sampling rates in Hertz of `x` and `y`, respectively, are `Fd` and `Fs`. (Thus  $1/F_d$  represents the time interval between two consecutive samples in `x`, and similarly for `y`.) The ratio  $F_s/F_d$  must be a positive integer. The initial phase in the modulation is zero.

The generic syntax `y = dmodce(x, Fd, [Fs phase], ...)` is the same, except that the third input argument is a two-element vector instead of a scalar. The first entry, `Fs`, is the sampling rate as described in the paragraph above. The second entry, `phase`, is the initial phase in the modulation, measured in radians.

### To Modulate a Digital Signal (Specific Syntax Information)

`y = dmodce(x, Fd, Fs, 'ask', M)` performs M-ary amplitude shift keying modulation. Each entry of `x` must be in the range  $[0, M-1]$ . The maximum value of the modulated signal is 1.

`y = dmodce(x, Fd, Fs, 'fsk', M, tone)` performs M-ary frequency shift keying modulation. Each entry of `x` must be in the range  $[0, M-1]$ . The optional argument `tone` is the separation between successive frequencies in the modulated signal `y`. The default value of `tone` is `Fd`. The maximum value of `y` is 1.

`y = dmodce(x, Fd, Fs, 'msk')` performs minimum shift keying modulation. Each entry of `x` is either 0 or 1. The maximum value of `y` is 1. The separation between the two frequencies is  $F_d/2$ .

`y = dmodce(x, Fd, Fs, 'psk', M)` performs M-ary phase shift keying modulation. Each entry of `x` must be in the range `[0, M-1]`. The maximum value of `y` is 1.

`y = dmodce(x, Fd, Fs, 'qask', M)` performs M-ary quadrature amplitude shift keying modulation with a square signal constellation. The table below shows the maximum value of `y`, for several small values of `M`.

M	Maximum Value of y	M	Maximum Value of y
2	1	32	5
4	1	64	7
8	3	128	11
16	3	256	15

**Note** To see how symbols are mapped to the constellation points, generate a square constellation plot using `qaskenco(M)`.

`y = dmodce(x, Fd, Fs, 'qask/arb', inphase, quadr)` performs quadrature amplitude shift keying modulation, with a signal constellation that you define using the vectors `inphase` and `quadr`. The constellation point for the  $k$ th message has in-phase component `inphase(k+1)` and quadrature component `quadr(k+1)`.

`y = dmodce(x, Fd, Fs, 'qask/cir', numsig, amp, phs)` performs quadrature amplitude shift keying modulation with a circular signal constellation. `numsig`, `amp`, and `phs` are vectors of the same length. The entries in `numsig` and `amp` must be positive. If  $k$  is an integer in the range `[1, length(numsig)]`, then `amp(k)` is the radius of the  $k$ th circle, `numsig(k)` is the number of constellation points on the  $k$ th circle, and `phs(k)` is the phase of the first constellation point plotted on the  $k$ th circle. All points on the  $k$ th circle are evenly spaced. If you omit `phs`, then its default value is `numsig*0`. If you omit `amp`, then its default value is `[1: length(numsig)]`.

---

**Note** To see how symbols are mapped to the constellation points, generate a labeled circle constellation plot using `apkconst (numsi g, amp, phs, 'n')`.

---

## Examples

This example uses FSK modulation and demodulation with different frequency separations `tone`. The output indicates that the symbol error rate varies depending on the value of `tone`. Your results might be different from those shown below, because the example uses random numbers.

```
M = 4; Fd = 1; Fs = 32;
SNRperBit = 5;
adjSNR = SNRperBit - 10*log10(Fs/Fd) + 10*log10(log2(M));
x = randint(5000, 1, M); % Original signal
% Modulate using FSK with orthogonal tone spacing.
tone = .5;
randn('state', 1945724); % Seed the Gaussian generator.
w1 = dmodce(x, Fd, Fs, 'fsk', M, tone);
y1 = awgn(w1, adjSNR, 'measured', [], 'dB');
z1 = ddemodce(y1, Fd, Fs, 'fsk', M, tone);
ser1 = symerr(x, z1)

ser1 =

    67

% Modulate using FSK with nonorthogonal tone spacing.
tone = .25;
randn('state', 1945724); % Reseed the Gaussian generator.
w2 = dmodce(x, Fd, Fs, 'fsk', M, tone);
y2 = awgn(w2, adjSNR, 'measured', [], 'dB');
z2 = ddemodce(y2, Fd, Fs, 'fsk', M, tone);
ser2 = symerr(x, z2)

ser2 =

   258
```

## See Also

`ddemodce`, `dmod`, `ddemod`, `amod`, `amodce`, `modmap`, `apkconst`

# dpcmdeco

---

**Purpose** Decode using differential pulse code modulation

**Syntax** `sig = dpcmdeco(indx, codebook, predictor);`  
`[sig, quanterror] = dpcmdeco(indx, codebook, predictor);`

**Description** `sig = dpcmdeco(indx, codebook, predictor)` implements differential pulse code demodulation to decode the vector `indx`. The vector `codebook` represents the predictive-error quantization codebook. The vector `predictor` specifies the predictive transfer function. If the transfer function has predictive order  $M$ , then `predictor` has length  $M+1$  and an initial entry of 0. To decode correctly, use the same codebook and predictor in `dpcmenco` and `dpcmdeco`.

See either “Representing Quantization Parameters” on page 2-14 or the reference page for `quantiz` in this chapter, for a description of the formats of `partition` and `codebook`.

`[sig, quanterror] = dpcmdeco(indx, codebook, predictor)` is the same as the syntax above, except that the vector `quanterror` is the quantization of the predictive error based on the quantization parameters. `quanterror` is the same size as `sig`.

---

**Note** You can estimate the input parameters `codebook`, `partition`, and `predictor` using the function `dpcmopt`.

---

**Examples** See the sections “Example: DPCM Encoding and Decoding” on page 2-20 and “Example: Comparing Optimized and Nonoptimized DPCM Parameters” on page 2-21 for examples that use `dpcmdeco`.

**See Also** `dpcmenco`, `dpcmopt`, `quantiz`, `compand`

**References** Kondo, A. M. *Digital Speech*. Chichester, England: John Wiley & Sons, 1994.



<b>Purpose</b>	Encode using differential pulse code modulation
<b>Syntax</b>	<pre>indx = dpcmenco(sig, codebook, partition, predictor) [indx, quants] = dpcmenco(sig, codebook, partition, predictor)</pre>
<b>Description</b>	<p><code>indx = dpcmenco(sig, codebook, partition, predictor)</code> implements differential pulse code modulation to encode the vector <code>sig</code>. <code>partition</code> is a vector whose entries give the endpoints of the partition intervals. <code>codebook</code>, a vector whose length exceeds the length of <code>partition</code> by one, prescribes a value for each partition in the quantization. <code>predictor</code> specifies the predictive transfer function. If the transfer function has predictive order <math>M</math>, then <code>predictor</code> has length <math>M+1</math> and an initial entry of 0. The output vector <code>indx</code> is the quantization index.</p> <p>See “Implementing Differential Pulse Code Modulation” on page 2-19 for more about the format of <code>predictor</code>. See either “Representing Quantization Parameters” on page 2-14 or the reference page for <code>quantiz</code> in this chapter, for a description of the formats of <code>partition</code> and <code>codebook</code>.</p> <p><code>[indx, quants] = dpcmenco(sig, codebook, partition, predictor)</code> is the same as the syntax above, except that <code>quants</code> contains the quantization of <code>sig</code> based on the quantization parameters. <code>quants</code> is a vector the same size as <code>sig</code>.</p> <hr/> <p><b>Note</b> If <code>predictor</code> is an order-one transfer function, then the modulation is called a delta-modulation.</p> <hr/>
<b>Examples</b>	See the sections “Example: DPCM Encoding and Decoding” on page 2-20 and “Example: Comparing Optimized and Nonoptimized DPCM Parameters” on page 2-21 for examples that use <code>dpcmenco</code> .
<b>See Also</b>	<code>dpcmdeco</code> , <code>dpcmopt</code> , <code>quantiz</code> , <code>compand</code>
<b>References</b>	Kondoz, A. M. <i>Digital Speech</i> . Chichester, England: John Wiley & Sons, 1994.

# dpcmopt

---

<b>Purpose</b>	Optimize differential pulse code modulation parameters
<b>Syntax</b>	<pre>predictor = dpcmopt(trainingset, ord); [predictor, codebook, partition] = dpcmopt(trainingset, ord, length); [predictor, codebook, partition] =     dpcmopt(trainingset, ord, initcodebook);</pre>
<b>Description</b>	<p><code>predictor = dpcmopt(trainingset, ord)</code> returns a vector representing a predictive transfer function of order <code>ord</code> that is appropriate for the training data in the vector <code>trainingset</code>. <code>predictor</code> is a row vector of length <code>ord+1</code>. See “Representing Quantization Parameters” on page 2-14 for more about its format.</p> <hr/> <p><b>Note</b> <code>dpcmopt</code> optimizes for the data in <code>trainingset</code>. For best results, <code>trainingset</code> should be similar to the data that you plan to quantize.</p> <hr/> <p><code>[predictor, codebook, partition] = dpcmopt(trainingset, ord, length)</code> is the same as the syntax above, except that it also returns corresponding optimized codebook and partition vectors <code>codebook</code> and <code>partition</code>. <code>length</code> is an integer that prescribes the length of codebook. <code>partition</code> is a vector of length <code>length-1</code>. See either “Representing Quantization Parameters” on page 2-14 or the reference page for <code>quantiz</code> in this chapter, for a description of the formats of <code>partition</code> and <code>codebook</code>.</p> <p><code>[predictor, codebook, partition] = dpcmopt(trainingset, ord, initcodebook)</code> is the same as the first syntax, except that it also returns corresponding optimized codebook and partition vectors <code>codebook</code> and <code>partition</code>. <code>initcodebook</code>, a vector of length at least 2, is the initial guess of the codebook values. The output <code>codebook</code> is a vector of the same length as <code>initcodebook</code>. The output <code>partition</code> is a vector whose length is one less than the length of <code>codebook</code>.</p>
<b>Examples</b>	See the section “Example: Comparing Optimized and Nonoptimized DPCM Parameters” on page 2-21 for an example that uses <code>dpcmopt</code> .
<b>See Also</b>	<code>dpcmenco</code> , <code>dpcmdeco</code> , <code>quantiz</code> , <code>lloyd</code> s

Purpose	Block encoder	
Syntax	<pre>code = encode(msg, n, k, 'linear/format', genmat); code = encode(msg, n, k, 'cyclic/format', genpoly); code = encode(msg, n, k, 'bch/format', genpoly); code = encode(msg, n, k, 'hamming/format', primpoly); code = encode(msg, n, k, 'rs/format', genpoly); code = encode(msg, field, k, 'rs/format', genpoly); code = encode(msg, n, k); [ code, added] = encode(...);</pre>	
Optional Inputs	Input	Default Value
	<i>format</i>	<b>binary</b>
	<i>genpoly</i>	cyclpoly(n, k) for cyclic codes; bchpoly(n, k) for BCH codes; rspoly(n, k) or rspoly(n, k, field) for Reed-Solomon codes
	<i>primpoly</i>	gfpri mdf (n- k)
Description	<p><b>For All Syntaxes</b></p> <p>The encode function encodes messages using one of the following error-correction coding methods:</p> <ul style="list-style-type: none"><li>• Linear block</li><li>• Cyclic</li><li>• BCH (Bose, Ray-Chaudhuri, Hocquenghem)</li><li>• Hamming</li><li>• Reed-Solomon</li></ul> <p>For all of these methods, the codeword length is n and the message length is k.</p> <p>msg, which represents the messages, can have one of several formats. Table 3-14, Information Formats for Encoding Methods Other than Reed-Solomon, below, which applies to all coding methods supported by encode <i>except</i> the Reed-Solomon method, shows which formats are allowed for msg, how the argument <i>format</i> should reflect the format of msg, and how the format of the output code depends on these choices. Table 3-15, Information Formats</p>	

for the Reed-Solomon Encoding Method, gives the corresponding information for the Reed-Solomon method. The examples in the tables are for  $k = 4$  and, in Table 3-15, Information Formats for the Reed-Solomon Encoding Method,  $m = 3$ . If *format* is not specified as input, then its default value is **binary**.

**Note** If  $2^n$  or  $2^k$  is large, then you should use the default **binary** format instead of the **decimal** format. This is because the function uses a binary format internally, while the round-off error associated with converting many bits to large decimal numbers and back might be substantial.

Table 3-14: Information Formats for Encoding Methods Other than Reed-Solomon

Format of msg	Value of "format" Argument	Format of code
Binary column vector	<b>binary</b>	Binary column vector
Example: msg = [0 1 1 0, 0 1 0 1, 1 0 0 1]'		
Binary matrix with k columns	<b>binary</b>	Binary matrix with n columns
Example: msg = [0 1 1 0; 0 1 0 1; 1 0 0 1]		
Column vector of integers in the range [0, $2^k-1$ ]	<b>decimal</b>	Column vector of integers in the range [0, $2^n-1$ ]
Example: msg = [6, 10, 9]'		

Table 3-15: Information Formats for the Reed-Solomon Encoding Method

Format of msg (where $n = 2^m-1$ , $m$ = integer greater than or equal to 3)	Value of "format" Argument	Format of code
Binary matrix with $m$ columns	<b>binary</b>	Binary matrix with $m$ columns
Example: msg = [1 1 0; 1 0 1; 1 0 0; 0 1 1; 1 1 0; 1 0 1; 1 0 0; 0 1 1]		

Table 3-15: Information Formats for the Reed-Solomon Encoding Method (Continued)

Format of msg (where $n = 2^m - 1$ , $m = \text{integer}$ greater than or equal to 3)	Value of "format" Argument	Format of code
Binary column vector	<b>binary</b>	Binary column vector
Example: msg = [ 1 1 0, 1 0 1, 1 0 0, 0 1 1, 1 1 0, 1 0 1, 1 0 0, 0 1 1 ]'		
Matrix of integers in the range [0, $2^m - 1$ ], with k columns	<b>decimal</b>	Matrix of integers in the range [0, $2^m - 1$ ], with n columns
Example: msg = [ 3, 5, 1, 6; 3, 5, 1, 6 ]		
Matrix of integers in the range [-1, $2^m - 2$ ], with k columns	<b>power</b>	Matrix of integers in the range [-1, $2^m - 2$ ], with n columns
Example: msg = [ 2, 4, 0, 5; 2, 4, 0, 5 ]		

### For Specific Syntaxes

`code = encode(msg, n, k, 'linear/format', genmat)` encodes msg using genmat as the generator matrix for the linear block encoding method. genmat, a k-by-n matrix, is required as input.

`code = encode(msg, n, k, 'cyclic/format', genpoly)` encodes msg and creates a systematic cyclic encode. genpoly is a row vector that gives the coefficients, in order of ascending powers, of the binary generator polynomial. The default value of genpoly is `cyclpoly(n, k)`. By definition, the generator polynomial for an  $[n, k]$  cyclic code must have degree  $n - k$  and must divide  $x^n - 1$ .

`code = encode(msg, n, k, 'bch/format', genpoly)` encodes msg using the BCH encoding method. genpoly is a row vector that gives the coefficients, in order of ascending powers, of the degree-(n - k) binary BCH generator polynomial. The default value of genpoly is `bchpoly(n, k)`. For this syntax, n must have the form  $2^m - 1$  for integer m greater than or equal to 3. k must be a valid message length as reported in the second column of params in the command

`params = bchpoly(n)`

`code = encode(msg, n, k, 'hamming/format', pri mpol y)` encodes `msg` using the Hamming encoding method. For this syntax, `n` must have the form  $2^m-1$  for some integer  $m$  greater than or equal to 3, and `k` must equal  $n-m$ . `pri mpol y` is a row vector that gives the binary coefficients, in order of ascending powers, of the primitive polynomial for  $\text{GF}(2^m)$  that is used in the encoding process. The default value of `pri mpol y` is the default primitive polynomial `gfpr mdf(m)`.

`code = encode(msg, n, k, 'rs/format', genpol y)` encodes `msg` using the Reed-Solomon encoding method. `n` must have the form  $2^m-1$  for some integer  $m$  greater than or equal to 3. `genpol y` is a row vector that gives the coefficients, in order of ascending powers, of the generator polynomial for the code. Each coefficient is an element of  $\text{GF}(2^m)$  expressed in exponential format. For a description of exponential format, see “Exponential Format” on page 2-90. The default value of `genpol y` is the output of the function `rs pol y`.

`code = encode(msg, fi el d, k, 'rs/format', genpol y)` is the same as the syntax above, except that `fi el d` is a matrix that lists all elements of  $\text{GF}(2^m)$  in the format described in “List of All Elements of a Galois Field” on page 2-91. The size of `fi el d` determines `n`. This syntax is faster than the one above.

`code = encode(msg, n, k)` is the same as `code = encode(msg, n, k, 'hamming/bi nary')`.

`[code, added] = encode(...)` returns the additional variable `added`. `added` is the number of zeros that were placed at the end of the message matrix before encoding, in order for the matrix to have the appropriate shape. “Appropriate” depends on `n`, `k`, the shape of `msg`, and the encoding method.

## Examples

The example below illustrates the three different information formats (binary vector, binary matrix, and decimal vector) for Hamming code. The three messages have identical content in different formats; as a result, the three codes that encode creates have identical content in correspondingly different formats.

```
m = 4; n = 2^m-1; % Codeword length = 15
k = 11; % Message length

% Create 100 messages, k bits each.
msg1 = randint(100*k, 1, [0, 1]); % As a column vector
msg2 = vec2mat(msg1, k); % As a k-column matrix
```

```
msg3 = bi2de(msg2); % As a column of decimal integers

% Create 100 codewords, n bits each.
code1 = encode(msg1, n, k, 'hamming/binary');
code2 = encode(msg2, n, k, 'hamming/binary');
code3 = encode(msg3, n, k, 'hamming/decimal');
if ( vec2mat(code1, n) == code2 & de2bi(code3, n) == code2 )
    disp('All three formats produced the same content.')
end
```

The next example creates a cyclic code, adds noise, and then decodes the noisy code. It uses the decode function. Your error rate results might vary because the noise is random.

```
n = 3; k = 2; % A (3, 2) cyclic code
msg = randint(100, k, [0, 1]); % 100 messages, k bits each
code = encode(msg, n, k, 'cyclic/binary');
% Add noise.
noisycode = rem(code + randerr(100, n, [0 1; .7 .3]), 2);
newmsg = decode(noisycode, n, k, 'cyclic'); % Try to decode.
% Compute error rate for decoding the noisy code.
[number, ratio] = biterr(newmsg, msg);
disp(['The bit error rate is ', num2str(ratio)])
```

The bit error rate is 0.08

The next example encodes the same message using Hamming, BCH, and cyclic methods. Before creating BCH code, it uses the bchpoly command to find out what codeword and message lengths are valid. This example also creates Hamming code with the '**linear**' option of the encode command. It then decodes each code and recovers the original message.

```
n = 6; % Try codeword length = 6.
% Find any valid message length for BCH code.
params = bchpoly(n);
n = params(1, 1); % Redefine codeword length in case earlier one
% was invalid.
k = params(1, 2); % Message length
m = log2(n+1); % Express n as 2^m-1.
msg = randint(100, 1, [0, 2^k-1]); % Column of decimal integers
```

```
% Create various codes.
codehamming = encode(msg, n, k, 'hamming/decimal');
[parmat, genmat] = hamngen(m);
codehamming2 = encode(msg, n, k, 'linear/decimal', genmat);
if codehamming==codehamming2
    disp('The ''linear'' method can create Hamming code.')
end
codebch = encode(msg, n, k, 'bch/decimal');
codecyclic = encode(msg, n, k, 'cyclic/decimal');

% Decode to recover the original message.
decodedhamming = decode(codehamming, n, k, 'hamming/decimal');
decodedbch = decode(codebch, n, k, 'bch/decimal');
decodedcyclic = decode(codecyclic, n, k, 'cyclic/decimal');
if (decodedhamming==msg & decodedbch==msg & decodedcyclic==msg)
    disp('All decoding worked flawlessly in this noiseless world.')
end
```

## Algorithm

Depending on the encoding method, encode relies on such lower-level functions as `hamngen`, `cyclgen`, `bchenco`, and `rsenco`.

## See Also

`decode`, `hamngen`, `cyclpoly`, `cyclgen`, `bchpoly`, `bchenco`, `rspoly`, `rsenco`, `rsencode`, `convenc`



**Purpose** Generate an eye diagram

**Syntax**

```
eyediagram(x, n);
eyediagram(x, n, period);
eyediagram(x, n, period, offset);
eyediagram(x, n, period, offset, plotstring);
eyediagram(x, n, period, offset, plotstring, h);
h = eyediagram(...);
```

**Description** `eyediagram(x, n)` creates an eye diagram for the signal `x`, plotting `n` samples in each trace. `n` must be an integer greater than 1. The labels on the horizontal axis of the diagram range between  $-1/2$  and  $1/2$ . The function assumes that the first value of the signal and every `n`th value thereafter, occur at integer times. The interpretation of `x` and the number of plots depend on the shape and complexity of `x`:

- If `x` is a real two-column matrix, then `eyediagram` interprets the first column as in-phase components and the second column as quadrature components. The two components appear in different subplots of a single figure window.
- If `x` is a complex vector, then `eyediagram` interprets the real part as in-phase components and the imaginary part as quadrature components. The two components appear in different subplots of a single figure window.
- If `x` is a real vector, then `eyediagram` interprets it as a real signal. The figure window contains a single plot.

`eyediagram(x, n, period)` is the same as the syntax above, except that the labels on the horizontal axis range between  $-\text{period}/2$  and  $\text{period}/2$ .

`eyediagram(x, n, period, offset)` is the same as the syntax above, except that the function assumes that the  $(\text{offset}+1)$ st value of the signal, and every `n`th value thereafter, occur at times that are integer multiples of `period`. The variable `offset` must be a nonnegative integer between 0 and `n-1`.

`eyediagram(x, n, period, offset, plotstring)` is the same as the syntax above, except that `plotstring` determines the plotting symbol, line type, and color for the plot. `plotstring` is a string whose format and meaning are the same as in the `plot` function.

# eyediagram

---

`eyediagram(x, n, period, offset, plotstring, h)` is the same as the syntax above, except that the eye diagram is in the figure whose handle is `h`, rather than a new figure. `h` must be a handle to a figure that `eyediagram` previously generated.

---

**Note** You cannot use `hold on` to plot multiple signals in the same figure.

---

`h = eyediagram(...)` is the same as the earlier syntaxes, except that `h` is the handle to the figure that contains the eye diagram.

## Examples

See “Example: Eye Diagrams” on page 2-9 for an example. For an online demonstration, use `scattereyedemo`.

## See Also

`scatterplot`, `plot`

**Purpose** Convert between parity-check and generator matrices

**Syntax**

```
parmat = gen2par(genmat);
genmat = gen2par(parmat);
```

**Description**

`parmat = gen2par(genmat)` converts the standard-form binary generator matrix `genmat` into the corresponding parity-check matrix `parmat`.

`genmat = gen2par(parmat)` converts the standard-form binary parity-check matrix `parmat` into the corresponding generator matrix `genmat`.

The standard forms of the generator and parity-check matrices for an  $[n,k]$  binary linear block code are shown in the table below.

Type of Matrix	Standard Form	Dimensions
Generator	$[I_k \ P]$ or $[P \ I_k]$	$k$ -by- $n$
Parity-check	$[-P \ I_{n-k}]$ or $[I_{n-k} \ -P]$	$(n-k)$ -by- $n$

where  $I_k$  is the identity matrix of size  $k$  and the ' ' symbol indicates matrix transpose. Two standard forms are listed for each type, since different authors use different conventions. For *binary* codes, the minus signs in the parity-check form listed above are irrelevant; that is,  $-1 = 1$  in the binary field.

**Examples**

The commands below convert the parity-check matrix for a Hamming code into the corresponding generator matrix and back again.

```
parmat = hamngen(3)
```

```
parmat =
```

```

1     0     0     1     0     1     1
0     1     0     1     1     1     0
0     0     1     0     1     1     1
```

```
genmat = gen2par(parmat)
```

## gen2par

---

```
genmat =
```

1	1	0	1	0	0	0
0	1	1	0	1	0	0
1	1	1	0	0	1	0
1	0	1	0	0	0	1

```
parmat2 = gen2par(genmat) % Ans should be the same as parmat above
```

```
parmat2 =
```

1	0	0	1	0	1	1
0	1	0	1	1	1	0
0	0	1	0	1	1	1

### See Also

encode, decode, hammgen, cycl gen

<b>Purpose</b>	Add polynomials over a Galois field
<b>Syntax</b>	<pre> c = gfadd(a, b); c = gfadd(a, b, p); c = gfadd(a, b, p, len); c = gfadd(a, b, field); </pre>
<b>Description</b>	<p><code>c = gfadd(a, b)</code> adds two GF(2) polynomials. The inputs and output are row vectors that give the coefficients of the corresponding polynomials in order of ascending powers. Each coefficient is either 0 or 1, since the field is GF(2). If <code>a</code> and <code>b</code> are matrices of the same size, then the function treats each row independently.</p> <p><code>c = gfadd(a, b, p)</code> adds two GF(p) polynomials, where <code>p</code> is a prime number. <code>a</code>, <code>b</code>, and <code>c</code> are row vectors that give the coefficients of the corresponding polynomials in order of ascending powers. Each coefficient is between 0 and <code>p-1</code>. If <code>a</code> and <code>b</code> are matrices of the same size, then the function treats each row independently.</p> <p><code>c = gfadd(a, b, p, len)</code> adds row vectors <code>a</code> and <code>b</code> as in the previous syntax, except that it returns a row vector of length <code>len</code>. The output <code>c</code> is a truncated or extended representation of the sum. If the row vector corresponding to the sum has fewer than <code>len</code> entries (including zeros), then extra zeros are added at the end; if it has more than <code>len</code> entries, then entries from the end are removed.</p> <p><code>c = gfadd(a, b, field)</code> adds two GF(<math>p^m</math>) elements, where <math>m</math> is a positive integer. <code>a</code> and <code>b</code> are the exponential format of the two elements, relative to some primitive element of GF(<math>p^m</math>). <code>field</code> is the matrix listing all elements of GF(<math>p^m</math>), arranged relative to the same primitive element. <code>c</code> is the exponential format of the sum, relative to the same primitive element. See “Representing Elements of Galois Fields” on page 2-90 for an explanation of these formats. If <code>a</code> and <code>b</code> are matrices of the same size, then the function treats each element independently.</p>
<b>Examples</b>	<p>In the code below, <code>sum</code> is the sum of <math>2 + 3x + x^2</math> and <math>4 + 2x + 3x^2</math> over GF(5), and <code>linpart</code> is the degree-one part of <code>sum</code>.</p> <pre> sum = gfadd([2 3 1], [4 2 3], 5) </pre>

```
sum =
```

```
1      0      4
```

```
linpart = gfadd([2 3 1], [4 2 3], 5, 2)
```

```
linpart =
```

```
1      0
```

The code below shows that  $\alpha^2 + \alpha^4 = \alpha^1$ , where  $\alpha$  is a root of the primitive polynomial  $2 + 2x + x^2$  for GF(9).

```
p = 3; m = 2;
primpoly = [2 2 1];
field = gftuple([-1: p^m-2]', primpoly, p);
g = gfadd(2, 4, field)
```

```
g =
```

```
1
```

Other examples are in the section, “Arithmetic in Galois Fields” on page 2-97.

## See Also

gfsub, gfmul, gfconv, gfdi v, gfdeconv, gftuple, bitxor

**Purpose** Multiply polynomials over a Galois field

**Syntax**

```
c = gfconv(a, b);
c = gfconv(a, b, p);
c = gfconv(a, b, field);
```

**Description** The `gfconv` function multiplies polynomials over a Galois field. (To multiply elements of a Galois field, use `gfmul` instead.) Algebraically, multiplying polynomials over a Galois field is equivalent to convolving vectors containing the polynomials' coefficients, where the convolution operation uses arithmetic over the same Galois field.

`c = gfconv(a, b)` multiplies two GF(2) polynomials. The inputs and output are row vectors that give the coefficients of the corresponding polynomials in order of ascending powers. Each coefficient is either 0 or 1, since the field is GF(2).

`c = gfconv(a, b, p)` multiplies two GF(p) polynomials, where  $p$  is a prime number. `a`, `b`, and `c` are row vectors that give the coefficients of the corresponding polynomials in order of ascending powers. Each coefficient is between 0 and  $p-1$ .

`c = gfconv(a, b, field)` multiplies two GF( $p^m$ ) polynomials, where  $p$  is a prime number and  $m$  is a positive integer. `a`, `b`, and `c` are row vectors that list the exponential formats of the coefficients of the corresponding polynomials, in order of ascending powers. The exponential format is relative to some primitive element of GF( $p^m$ ). `field` is the matrix listing all elements of GF( $p^m$ ), arranged relative to the same primitive element. See "Representing Elements of Galois Fields" on page 2-90 for an explanation of these formats.

**Examples** The command below shows that  $(1 + x + x^4)(x + x^2) = x + x^3 + x^5 + x^6$  over GF(2).

```
gfc = gfconv([1 1 0 0 1], [0 1 1])
```

```
gfc =
```

```
0      1      0      1      0      1      1
```

The code below illustrates the identity

$$(x^r + x^s)^p = x^{rp} + x^{sp} \text{ in GF}(p)$$

for the case in which  $p = 7$ ,  $r = 5$ , and  $s = 3$ . (The identity holds when  $p$  is any prime number, and  $r$  and  $s$  are positive integers.)

```
p = 7; r = 5; s = 3;
a = gfrepconv([r s]); % x^r + x^s

% Compute a^p over GF(p).
c = 1;
for ii = 1:p
    c = gfconv(c, a, p);
end;

% Check whether c = x^(rp) + x^(sp).
powers = [];
for ii = 1:length(c)
    if c(ii)~=0
        powers = [powers, ii];
    end;
end;
if (powers==[r*p+1 s*p+1] | powers==[s*p+1 r*p+1])
    disp('The identity is proved for this case of r, s, and p.')
end
```

### See Also

gfdeconv, gfadd, gfsub, gfmul, gftuple, conv



**Purpose** Produce cyclotomic cosets for a Galois field

**Syntax**

```
cs = gfcosets(m);
cs = gfcosets(m, p);
```

**Description** `cs = gfcosets(m)` produces the cyclotomic cosets for  $GF(2^m)$ , where  $m$  is a positive integer.

`cs = gfcosets(m, p)` produces the cyclotomic cosets for  $GF(p^m)$ , where  $m$  is a positive integer and  $p$  is a prime number.

In both cases, the output matrix `cs` is structured so that each row represents one coset. The row represents the coset by giving the exponential format of the elements of the coset, relative to the default primitive polynomial for the field. For a description of exponential formats, see “Representing Elements of Galois Fields” on page 2-90.

The first column contains the coset leaders. Because the lengths of cosets may vary, entries of NaN are used to fill the extra spaces when necessary to make `cs` rectangular.

A cyclotomic coset is a set of elements that all satisfy the same minimal polynomial. For more details on cyclotomic cosets, see the works listed in “References” below.

**Examples** The command below finds the cyclotomic cosets for  $GF(9)$ .

```
cs = gfcosets(2, 3)
```

```
cs =
```

```

0   NaN
1    3
2    6
4   NaN
5    7
```

The `gfmi npol` function can check that the elements of, for example, the third row of `cs` indeed belong in the same coset.

```
m = [gfminpol(2, 2, 3); gfminpol(6, 2, 3)] % Rows are identical.
```

```
m =
```

```
2     0     1
2     0     1
```

### See Also

gfminpol, gfroots, gfpri MDF

### References

Blahut, Richard E. *Theory and Practice of Error Control Codes*. Reading, Mass.: Addison-Wesley, 1983, p.105.

Lin, Shu and Daniel J. Costello, Jr. *Error Control Coding: Fundamentals and Applications*. Englewood Cliffs, N.J.: Prentice-Hall, 1983.

**Purpose** Divide polynomials over a Galois field

**Syntax**

```
[quot, remd] = gfdeconv(b, a);
[quot, remd] = gfdeconv(b, a, p);
[quot, remd] = gfdeconv(b, a, field);
```

**Description** The `gfdeconv` function divides polynomials over a Galois field. (To divide elements of a Galois field, use `gfdiv` instead.) Algebraically, dividing polynomials over a Galois field is equivalent to deconvolving vectors containing the polynomials' coefficients, where the deconvolution operation uses arithmetic over the same Galois field.

`[quot, remd] = gfdeconv(b, a)` divides the polynomial `b` by the polynomial `a` over  $GF(2)$  and returns the quotient in `quot` and the remainder in `remd`. All inputs and outputs are row vectors that give the coefficients of the corresponding polynomials in order of ascending powers. Each coefficient is either 0 or 1, since the field is  $GF(2)$ .

`[quot, remd] = gfdeconv(b, a, p)` divides the polynomial `b` by the polynomial `a` over  $GF(p)$  and returns the quotient in `quot` and the remainder in `remd`. `p` is a prime number. `b`, `a`, `quot`, and `remd` are row vectors that give the coefficients of the corresponding polynomials in order of ascending powers. Each coefficient is between 0 and `p-1`.

`[quot, remd] = gfdeconv(b, a, field)` divides the polynomial `b` by the polynomial `a` over  $GF(p^m)$  and returns the quotient in `quot` and the remainder in `remd`. Here `p` is a prime number and `m` is a positive integer. `b`, `a`, `quot`, and `remd` are row vectors that list the exponential formats of the coefficients of the corresponding polynomials, in order of ascending powers. The exponential format is relative to some primitive element of  $GF(p^m)$ . `field` is the matrix listing all elements of  $GF(p^m)$ , arranged relative to the same primitive element. See "Representing Elements of Galois Fields" on page 2-90 for an explanation of these formats.

**Examples** The code below shows that

$$(x + x^3 + x^4) \div (1 + x) = 1 + x^3 \text{ Remainder } 1$$

in  $GF(2)$ . It also checks the results of the division.

```
p = 2;
b = [0 1 0 1 1]; a = [1 1];
[quot, remd] = gfdeconv(b, a, p)
% Check the result.
bnew = gfadd(gfconv(quot, a, p), remd, p);
if isequal(bnew, b)
    disp('Correct. ')
end;
```

The output is below.

quot =

1      0      0      1

remd =

1

Correct.

Working over GF(3), the code below outputs those polynomials of the form  $x^k - 1$  ( $k = 2, 3, 4, \dots, 8$ ) that  $1 + x^2$  divides evenly.

```
p = 3; m = 2;
a = [1 0 1]; % 1+x^2
for ii = 2:p^m-1
    b = gfrepconv(ii); % x^ii
    b(1) = p-1; % -1+x^ii
    [quot, remd] = gfdeconv(b, a, p);
    % Display -1+x^ii if a divides it evenly.
    if remd==0
        gfpretty(b)
    end
end
```

The output is below.

$$\begin{matrix} & & & 4 \\ & & & 2 + X \end{matrix}$$

$$x^8 + x^2 + 1$$

In light of the discussion in “Algorithm” on the reference page for `gfprimck` along with the irreducibility of  $1 + x^2$  over GF(3), this output indicates that  $1 + x^2$  is not primitive for GF(9).

### Algorithm

The algorithm of `gfdeconv` is similar to that of the MATLAB function `deconv`.

### See Also

`gfconv`, `gfadd`, `gfsub`, `gfdi v`, `gftuple`, `deconv`

**Purpose** Divide elements of a Galois field

**Syntax**

```
quot = gfdiv(b, a);  
quot = gfdiv(b, a, p);  
quot = gfdiv(b, a, field);
```

**Description** The `gfdiv` function divides elements of a Galois field. (To divide polynomials over a Galois field, use `gfdeconv` instead.)

`quot = gfdiv(b, a)` divides `b` by `a` in  $\text{GF}(2)$  and returns the quotient. If `a` and `b` are matrices of the same size, then the function treats each element independently. All entries of `b`, `a`, and `quot` are either 0 or 1, since the field is  $\text{GF}(2)$ .

`quot = gfdiv(b, a, p)` divides `b` by `a` in  $\text{GF}(p)$  and returns the quotient. `p` is a prime number. If `a` and `b` are matrices of the same size, then the function treats each element independently. All entries of `b`, `a`, and `quot` are between 0 and `p-1`.

`quot = gfdiv(b, a, field)` divides `b` by `a` in  $\text{GF}(p^m)$  and returns the quotient. `p` is a prime number and `m` is a positive integer. If `a` and `b` are matrices of the same size, then the function treats each element independently. All entries of `b`, `a`, and `quot` are the exponential formats of elements of  $\text{GF}(p^m)$  relative to some primitive element of  $\text{GF}(p^m)$ . `field` is the matrix listing all elements of  $\text{GF}(p^m)$ , arranged relative to the same primitive element. See “Representing Elements of Galois Fields” on page 2-90 for an explanation of these formats.

In all cases, an attempt to divide by the zero element of the field results in a “quotient” of NaN.

**Examples** The code below displays lists of multiplicative inverses in  $\text{GF}(5)$  and  $\text{GF}(25)$ . It uses column vectors as inputs to `gfdiv`.

```
% Find inverses of nonzero elements of GF(5).  
p = 5;  
b = ones(p-1, 1);  
a = [1:p-1]';  
quot1 = gfdiv(b, a, p);  
disp('Inverses in GF(5):')  
disp('element inverse')  
disp([a, quot1])
```

```
% Find inverses of nonzero elements of GF(25).
m = 2;
field = gftuple([-1:p^m-2]', m, p);
b = zeros(p^m-1, 1); % Numerator is zero since 1 = alpha^0.
a = [0:p^m-2]';
quot2 = gfddiv(b, a, field);
disp('Inverses in GF(25), expressed in EXPONENTIAL FORMAT with')
disp('respect to a root of the default primitive polynomial:')
disp('element inverse')
disp([a, quot2])
```

## See Also

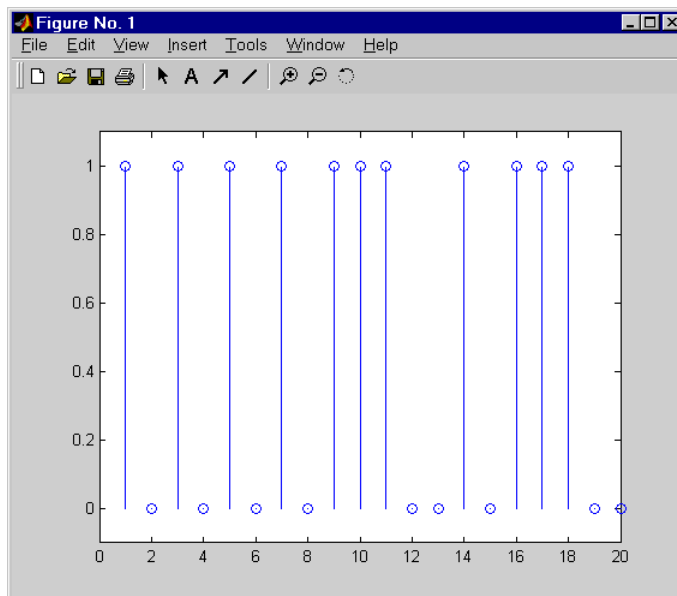
gfmul, gfdeconv, gfconv, gftuple

# gffilter

---

<b>Purpose</b>	Filter data using polynomials over a prime Galois field
<b>Syntax</b>	<pre>y = gffilter(b, a, x); y = gffilter(b, a, x, p);</pre>
<b>Description</b>	<p><code>y = gffilter(b, a, x)</code> filters the data <code>x</code> using the filter described by vectors <code>a</code> and <code>b</code>. <code>y</code> is the filtered data in GF(2).</p> <p><code>y = gffilter(b, a, x, p)</code> filters the data <code>x</code> using the filter described by vectors <code>a</code> and <code>b</code>. <code>y</code> is the filtered data in GF(p). <code>p</code> is a prime number, and all entries of <code>a</code> and <code>b</code> are between 0 and <code>p-1</code>.</p> <p>By definition of the filter, <code>y</code> solves the difference equation below</p> $a(1)y(n) = b(1)x(n) + b(2)x(n-1) + b(3)x(n-2) + \dots + b(B+1)x(n-B) - a(2)y(n-1) - a(3)y(n-2) - \dots - a(A+1)y(n-A)$ <p>where:</p> <ul style="list-style-type: none"><li>• <code>A+1</code> is the length of the vector <code>a</code></li><li>• <code>B+1</code> is the length of the vector <code>b</code></li><li>• <code>n</code> varies between 1 and the length of the vector <code>x</code>.</li></ul> <p>The vector <code>a</code> represents the degree-<math>n_a</math> polynomial</p> $a(1) + a(2)x + a(3)x^2 + \dots + a(A+1)x^A$
<b>Examples</b>	<p>The impulse response of a particular filter is given in the code and diagram below.</p> <pre>b = [1 0 0 1 0 1 0 1]; a = [1 0 1 1]; y = gffilter(b, a, [1, zeros(1, 19)]); stem(y); axis([0 20 -.1 1.1])</pre>





## Algorithm

For filters over  $\text{GF}(2)$  only, `gffilter` uses an algorithm similar to that used by the MATLAB function `filter`. You can use `filter` for filters over  $\text{GF}(2)$  by using the command below.

```
y = abs(rem(filter(b, a, x), 2));
```

However, this may produce an error if  $a$  is not stable in the regular discrete-time system analysis and the vector  $x$  is too long, or for a high order filter. `gffilter` produces an accurate result in all cases.

## See Also

`filter`, `gfconv`, `gfadd`

# gflineq

---

**Purpose** Solve the linear equation  $Ax = b$  over a prime Galois field

**Syntax**

```
x = gflineq(A, b);  
x = gflineq(A, b, p);  
[x, vld] = gflineq(...);
```

**Description** `x = gflineq(A, b)` solves the linear equation  $Ax = b$  over  $GF(2)$ . If  $A$  is a  $k$ -by- $n$  matrix and  $b$  is a vector of length  $k$ , then  $x$  is a vector of length  $n$ . Each entry of  $A$ ,  $x$ , and  $b$  is either 0 or 1. If no solution exists, then  $x$  is empty.

`x = gflineq(A, b, p)` solves the linear equation  $Ax = b$  over  $GF(p)$ , where  $p$  is a prime number. If  $A$  is a  $k$ -by- $n$  matrix and  $b$  is a vector of length  $k$ , then  $x$  is a vector of length  $n$ . Each entry of  $A$ ,  $x$ , and  $b$  is an integer between 0 and  $p-1$ .

`[x, vld] = gflineq(...)` returns a flag `vld` that indicates the existence of a solution. If `vld = 1`, then the solution  $x$  exists and is valid; if `vld = 0`, then no solution exists.

**Examples** The code below produces some valid solutions of a linear equation over  $GF(2)$ .

```
A=[1 0 1;  
   1 1 0;  
   1 1 1];  
% An example in which the solutions are valid  
[x, vld] = gflineq(A, [1; 0; 0])
```

`x =`

```
1  
1  
0
```

`vld =`

```
1
```

By contrast, the command below finds that the linear equation has *no* solutions.

```
[x2, v1 d2] = gflin eq(zeros(3, 3), [1; 0; 0])
This linear equation has no solution.
```

```
x2 =
```

```
[]
```

```
v1 d2 =
```

```
0
```

**Algorithm** `gflin eq` uses Gaussian elimination.

**See Also** `gfrank`, `gfadd`, `gfdiv`, `gfroots`, `gfconv`, `conv`

**Purpose** Find the minimal polynomial of an element of a Galois field

**Syntax**

```
pol = gfminpol (k, m);  
pol = gfminpol (k, primpoly);  
pol = gfminpol (k, m, p);  
pol = gfminpol (k, primpoly, p);
```

**Description** `pol = gfminpol (k, m)` finds the minimal polynomial of  $\alpha^k$  over  $\text{GF}(2)$ , where  $\alpha$  is a root of the default primitive polynomial for  $\text{GF}(2^m)$ .  $m$  is an integer greater than one. The format of the output is listed below:

- If  $k$  is a nonnegative integer, then `pol` is a row vector that gives the coefficients of the minimal polynomial in order of ascending powers.
- If  $k$  is a vector of length `len` all of whose entries are nonnegative integers, then `pol` is a matrix having `len` rows; the  $r$ th row of `pol` gives the coefficients of the minimal polynomial of  $\alpha^{k(r)}$  in order of ascending powers.

`pol = gfminpol (k, primpoly)` is the same as the first syntax listed, except that  $\alpha$  is a root of the primitive polynomial for  $\text{GF}(2^m)$  specified by `primpoly`. `primpoly` is a row vector that gives the coefficients of the degree- $m$  primitive polynomial in order of ascending powers.

`pol = gfminpol (k, m, p)` is the same as the first syntax listed, except that 2 is replaced by a prime number  $p$ .

`pol = gfminpol (k, primpoly, p)` is the same as the first syntax listed, except that 2 is replaced by a prime number  $p$ , and that  $\alpha$  is a root of the primitive polynomial for  $\text{GF}(p^m)$  specified by `primpoly`. `primpoly` is a row vector that gives the coefficients of the degree- $m$  primitive polynomial in order of ascending powers.

**Examples** The syntax `gfminpol (k, m, p)` is used in the sample code in the section “Characterization of Polynomials” on page 2-101.

As another example, the code below determines which elements of  $\text{GF}(2^4)$  are also in  $\text{GF}(2^2)$ , by considering the degrees of their minimal polynomials.

```
p = 2; m = 4; % Consider elements of GF(16).  
primpoly = gfpri mdf (4);
```

```
% Get minimal polys for all elements except 0 and 1.
k = [1:p^m-2];
minpolys = gfminpol(k, primpoly);

% Check which minimal polys have degree 2.
gf4=[];
for ii = 1:p^m-2
    if length(gftrunc(minpolys(ii,:)))==3 % A degree-2 polynomial
        gf4=[gf4, ii];
    end
end

disp(['The elements of GF(4) are 0, 1, alpha^',...
    int2str(gf4(1)), ' and alpha^',int2str(gf4(2))])
disp('where alpha is a root in GF(16) of the polynomial')
gfpretty(primpoly)
```

The output is below.

```
The elements of GF(4) are 0, 1, alpha^5 and alpha^10
where alpha is a root in GF(16) of the polynomial
```

$$1 + X + X^4$$

## See Also

gfprimdf, gfcosets, gfroots

# gfmul

---

**Purpose** Multiply elements of a Galois field

**Syntax**

```
c = gfmul (a, b);  
c = gfmul (a, b, p);  
c = gfmul (a, b, field);
```

**Description** The `gfmul` function multiplies elements of a Galois field. (To multiply polynomials over a Galois field, use `gfconv` instead.)

`c = gfmul (a, b)` multiplies `a` and `b` in  $\text{GF}(2)$ . Each entry of `a` and `b` is either 0 or 1. If `a` and `b` are matrices of the same size, then the function treats each element independently.

`c = gfmul (a, b, p)` multiplies `a` and `b` in  $\text{GF}(p)$ . Each entry of `a` and `b` is between 0 and  $p-1$ .  $p$  is a prime number. If `a` and `b` are matrices of the same size, then the function treats each element independently.

`c = gfmul (a, b, field)` multiplies `a` and `b` in  $\text{GF}(p^m)$ , where  $p$  is a prime number and  $m$  is a positive integer. `a` and `b` represent elements of  $\text{GF}(p^m)$  in exponential format relative to some primitive element of  $\text{GF}(p^m)$ . `field` is the matrix listing all elements of  $\text{GF}(p^m)$ , arranged relative to the same primitive element. `c` is the exponential format of the product, relative to the same primitive element. See “Representing Elements of Galois Fields” on page 2-90 for an explanation of these formats. If `a` and `b` are matrices of the same size, then the function treats each element independently.

**Examples** The section “Arithmetic in Galois Fields” on page 2-97 contains examples. Also, the code below shows that  $\alpha^2 \cdot \alpha^4 = \alpha^6$ , where  $\alpha$  is a root of the primitive polynomial  $2 + 2x + x^2$  for  $\text{GF}(9)$ .

```
p = 3; m = 2;  
primpoly = [2 2 1];  
field = gftuple([-1; p^m-2]', primpoly, p);  
a = gfmul(2, 4, field)
```

```
a =
```

```
6
```

**See Also** `gfdiv`, `gfconv`, `gfdeconv`, `gftuple`

**Purpose** Add elements of a Galois field of characteristic two

**Syntax** `k = gfplus(a, b, fvec, ivec);`

**Description** `k = gfplus(a, b, fvec, ivec)` adds `a` and `b` in  $\text{GF}(2^m)$ , using the exponential format to represent inputs and outputs. If `a` and `b` are matrices, then they must have the same dimensions, and `gfplus` adds them element-by-element. See “Representing Elements of Galois Fields” on page 2-90 for an explanation of the exponential format.

`fvec` and `ivec` are vectors of length  $2^m$ . The entries in both are integers between 0 and  $2^m-1$ . `fvec` contains the same information as the `field` parameter as used by `gfadd`, except that `fvec` has been condensed into a vector. To compute `fvec` and `ivec`, define `m` and then use the commands below.

```
fvec = gftuple([-1 : 2^m-2]', m) * 2.^[0 : m-1]';
ivec(fvec + 1) = 0 : 2^m - 1;
```

Alternatively, define a primitive polynomial vector `pol` and then use the commands below. See `gfpriofd` for information about defining `pol`.

```
fvec = gftuple([-1 : 2^m-2]', pol) * 2.^[0 : m-1]';
ivec(fvec + 1) = 0 : 2^m - 1;
```

**Examples** This example adds two matrices, each of which contains random nonzero elements of  $\text{GF}(2^5)$ .

```
m = 5;
a = randint(3, 6, 2^m-1, 1234); % Create a 3-by-6 matrix in GF(2^5).
b = randint(3, 6, 2^m-1);
fvec = gftuple([-1 : 2^m - 2]', m) * 2.^[0 : m-1]';
ivec(fvec + 1) = 0 : 2^m - 1;
aplsb = gfplus(a, b, fvec, ivec) % Add.

aplsb =

    22    25     4    23     9    29
     9    30    24    23  -Inf    19
    15    17    10    12    25    30
```

**See Also** `gfadd`, `gfsub`

**Purpose** Display a polynomial in traditional format

**Syntax** `gfpretty(a)`  
`gfpretty(a, st)`  
`gfpretty(a, st, n)`

**Description** `gfpretty(a)` displays a polynomial in a traditional format, using `X` as the variable and the entries of the row vector `a` as the coefficients in order of ascending powers. The polynomial is displayed in order of ascending powers. Terms having a zero coefficient are not displayed.

`gfpretty(a, st)` is the same as the first syntax listed, except that the content of the string `st` is used as the variable instead of `X`.

`gfpretty(a, st, n)` is the same as the first syntax listed, except that the content of the string `st` is used as the variable instead of `X`, and each line of the display has width `n` instead of the default value of 79.

---

**Note** For all syntaxes: If you do not use a fixed-width font, then the spacing in the display might not look correct.

---

**Examples** The code below displays statements about the elements of GF(16).

```
p = 2; m = 4;
ii = randint(1, 1, [1, p^m-2]); % Random exponent for prim element
primpolys = gfprimfd(m, 'all');
[rows, cols] = size(primpolys);
jj = randint(1, 1, [1, rows]); % Random primitive polynomial

disp('If A is a root of the primitive polynomial')
gfpretty(primpolys(jj, :)) % Polynomial in X
disp('then the element')
gfpretty([zeros(1, ii), 1], 'A') % The polynomial A^ii
disp('can also be expressed as')
gfpretty(gftuple(ii, m, p), 'A') % Polynomial in A
```

Below is a sample of the output.



If A is a root of the primitive polynomial

$$1 + X^3 + X^4$$

then the element

$$A^5$$

can also be expressed as

$$A + A^2$$

**See Also** `gftuple`, `gfprimdf`

# gfprimck

---

<b>Purpose</b>	Check whether a polynomial over a Galois field is primitive
<b>Syntax</b>	<pre>ck = gfprimck(a); ck = gfprimck(a, p);</pre>
<b>Description</b>	<p><code>ck = gfprimck(a)</code> returns a flag <code>ck</code> that indicates whether a polynomial over GF(2) is irreducible or primitive. <code>a</code> is a row vector that gives the coefficients of the polynomial in order of ascending powers. Each coefficient is either 0 or 1, since the field is GF(2). If <math>m</math> is the degree of the polynomial, then the output <code>ck</code> is:</p> <ul style="list-style-type: none"><li>• -1 if <code>a</code> is not an irreducible polynomial</li><li>• 0 if <code>a</code> is irreducible but not a primitive polynomial for GF(<math>2^m</math>)</li><li>• 1 if <code>a</code> is a primitive polynomial for GF(<math>2^m</math>)</li></ul> <p>This function considers the zero polynomial to be “not irreducible” and considers all polynomials of degree zero or one to be primitive.</p> <p><code>ck = gfprimck(a, p)</code> is the same as the syntax listed above, except that 2 is replaced by a prime number <code>p</code>.</p>
<b>Examples</b>	The section “Characterization of Polynomials” on page 2-101 contains examples.
<b>Algorithm</b>	An irreducible polynomial over GF( $p$ ) of degree at least 2 is primitive if and only if it does not divide $-1 + x^k$ for any positive integer $k$ smaller than $p^m - 1$ .
<b>See Also</b>	<code>gfprimfd</code> , <code>gfprimdf</code> , <code>gftuple</code> , <code>gfminpol</code> , <code>gfadd</code>
<b>References</b>	Clark, George C. Jr. and J. Bibb Cain. <i>Error-Correction Coding for Digital Communications</i> . New York: Plenum Press, 1981.

<b>Purpose</b>	Provide default primitive polynomials for a Galois field
<b>Syntax</b>	<pre>pol = gfprimdf(m); pol = gfprimdf(m, p);</pre>
<b>Description</b>	<p><code>pol = gfprimdf(m)</code> returns the row vector that gives the coefficients, in order of ascending powers, of MATLAB's default primitive polynomial for <math>\text{GF}(2^m)</math>. <code>m</code> is a positive integer.</p> <p><code>pol = gfprimdf(m, p)</code> returns the row vector that gives the coefficients, in order of ascending powers, of MATLAB's default primitive polynomial for <math>\text{GF}(p^m)</math>. <code>m</code> is a positive integer and <code>p</code> is a prime number.</p>
<b>Examples</b>	<p>The command below shows that <math>2 + x + x^2</math> is the default primitive polynomial for <math>\text{GF}(5^2)</math>.</p> <pre>pol = gfprimdf(2, 5)</pre> <pre>pol =</pre> <pre>      2      1      1</pre> <p>The code below displays the default primitive polynomial for each of the fields <math>\text{GF}(2^m)</math>, where <math>m</math> ranges between 3 and 5.</p> <pre>for m = 3:5     gfpretty(gfprimdf(m)) end</pre> <pre>           3       1 + X + X            4       1 + X + X            2      5       1 + X  + X </pre>
<b>See Also</b>	<code>gfprimck</code> , <code>gfprimdf</code> , <code>gftuple</code> , <code>gfminpol</code>

# gfprimfd

**Purpose** Find primitive polynomials for a Galois field

**Syntax**

```
pol = gfprimfd(m);  
pol = gfprimfd(m, opt);  
pol = gfprimfd(m, opt, p);
```

**Description** For all syntaxes:

- If  $m = 1$ , then  $\text{pol} = [1 \ 1]$ .
- A polynomial is represented as a row containing the coefficients in order of ascending powers.

$\text{pol} = \text{gfprimfd}(m)$  returns the row vector representing one primitive polynomial for  $\text{GF}(2^m)$ .  $m$  is a positive integer.

$\text{pol} = \text{gfprimfd}(m, \text{opt})$  searches for one or more primitive polynomials for  $\text{GF}(2^m)$ , where  $m$  is a positive integer. If  $m > 1$ , then the output  $\text{pol}$  depends on the argument  $\text{opt}$  as shown in the table below.

opt	Significance of pol	Format of pol
'min'	One primitive polynomial for $\text{GF}(2^m)$ having the smallest possible number of nonzero terms	The row vector representing the polynomial
'max'	One primitive polynomial for $\text{GF}(2^m)$ having the greatest possible number of nonzero terms	The row vector representing the polynomial
'all'	All primitive polynomials for $\text{GF}(2^m)$	A matrix, each row of which represents one such polynomial
A positive integer	All primitive polynomials for $\text{GF}(2^m)$ that have $\text{opt}$ nonzero terms	A matrix, each row of which represents one such polynomial

$\text{pol} = \text{gfprimfd}(m, \text{opt}, p)$  is the same as  $\text{pol} = \text{gfprimfd}(m, \text{opt})$  except that 2 is replaced by a prime number  $p$ .

## Examples

The code below seeks primitive polynomials for GF(32) having various other properties. Notice that `fourterms` is empty because no primitive polynomial for GF(32) has exactly four nonzero terms. Also notice that `manyterms` represents a *single* polynomial having five terms, while `five terms` represents *all* of the five-term primitive polynomials for GF(32).

```
p = 2; m = 5; % Work in GF(32).
manyterms = gfprimfd(5, 'max')
five terms = gfprimfd(5, 5)
fourterms = gfprimfd(5, 4)
```

The output is below.

```
manyterms =

    1    1    1    1    0    1

five terms =

    1    1    1    1    0    1
    1    1    1    0    1    1
    1    1    0    1    1    1
    1    0    1    1    1    1
```

No primitive polynomial satisfies the given constraints.

```
fourterms =

[]
```

## Algorithm

`gfprimfd` tests for primitivity using `gfprimck`. If `opt` is '**min**', '**max**', or omitted, then polynomials are constructed by converting decimal integers to base `p`. Based on the decimal ordering, `gfprimfd` returns the first polynomial it finds that satisfies the appropriate conditions.

## See Also

`gfprimck`, `gfprimdf`, `gftuple`, `gfminpol`

# gfrank

---

**Purpose** Compute the rank of a matrix over a Galois field

**Syntax** `rk = gfrank(A);`  
`rk = gfrank(A, p);`

**Description** `rk = gfrank(A)` calculates the rank of the matrix A in GF(2).  
`rk = gfrank(A, p)` calculates the rank of the matrix A in GF(p), where p is a prime number.

**Algorithm** gfrank uses an algorithm similar to Gaussian elimination.

**Examples** In the code below, gfrank says that the matrix A has less than full rank. This conclusion makes sense because the determinant of A is zero mod 2.

```
A=[1 0 1;  
   1 1 0;  
   0 1 1];  
det_a = det(A); % Ordinary determinant of A  
detmod2 = rem(det(A), 2); % Determinant mod 2  
rank2 = gfrank(A);  
disp(['determinant = ', num2str(det_a)])  
disp(['determinant mod 2 is ', num2str(detmod2)])  
disp(['rank over GF(2) is ', num2str(rank2)])
```

The output is below.

```
determinant = 2  
determinant mod 2 is 0  
rank over GF(2) is 2
```

Notice that `gfli neq` finds only the trivial solution to the equation  $Ax = 0$ , even though the output above implies that there are infinitely many other solutions.

```
sol = gfli neq(A, [0; 0; 0])'
```

```
sol =
```

```
0      0      0
```

**See Also** `gfli neq`

<b>Purpose</b>	Convert one GF(2) polynomial representation to another
<b>Syntax</b>	<code>polystandard = gfrepcov(poly2)</code>
<b>Description</b>	<p>Two logical ways to represent polynomials over GF(2) are listed below:</p> <p><b>1</b> <code>[A_0 A_1 A_2 ... A_(m-1)]</code> represents the polynomial <math>A_0 + A_1x + A_2x^2 + \dots + A_{(m-1)}x^{m-1}</math>. Each entry <math>A_k</math> is either one or zero.</p> <p><b>2</b> <code>[A_0 A_1 A_2 ... A_(m-1)]</code> represents the polynomial <math>x^{A_0} + x^{A_1} + x^{A_2} + \dots + x^{A_{(m-1)}}</math>. Each entry <math>A_k</math> is a nonnegative integer. All entries must be distinct.</p> <p>Format <b>1</b> is the standard form used by the Galois field functions in this toolbox, but there are some cases in which format <b>2</b> is more convenient.</p> <p><code>polystandard = gfrepcov(poly2)</code> converts from the second format to the first, for polynomials of degree <i>at least</i> 2. <code>poly2</code> and <code>polystandard</code> are row vectors. The entries of <code>poly2</code> are distinct integers, and at least one entry must exceed 1. Each entry of <code>polystandard</code> is either 0 or 1.</p> <hr/> <p><b>Note</b> If <code>poly2</code> is a <i>binary</i> row vector, then <code>gfrepcov</code> assumes that it is already in Format 1 above and returns it unaltered.</p> <hr/>
<b>Examples</b>	<p>The command below converts the representation format of the polynomial <math>1 + x^2 + x^5</math>.</p> <pre>polystandard = gfrepcov([0 2 5])</pre> <pre>polystandard =</pre> <pre> 1      0      1      0      0      1</pre>
<b>See Also</b>	<code>gfpretty</code>

# groots

---

**Purpose** Find the roots of a polynomial over a prime Galois field

**Syntax**

```
rt = groots(f);  
rt = groots(f, m);  
rt = groots(f, primpoly);  
rt = groots(f, m, p);  
rt = groots(f, primpoly, p);  
[rt, rt_tuple] = groots(...);  
[rt, rt_tuple, field] = groots(...);
```

**Description** For all syntaxes,  $f$  is a row vector that gives the coefficients, in order of ascending powers, of a degree- $d$  polynomial.

---

**Note** `groots` lists each root exactly once, ignoring multiplicities of roots.

---

`rt = groots(f)` finds roots in  $\text{GF}(2^d)$  of the polynomial that  $f$  represents.  $rt$  is a column vector each of whose entries is the exponential format of a root. The exponential format is relative to a root of the default primitive polynomial for  $\text{GF}(2^d)$ .

`rt = groots(f, m)` finds roots in  $\text{GF}(2^m)$  of the polynomial that  $f$  represents.  $m$  is an integer greater than or equal to  $d$ .  $rt$  is a column vector each of whose entries is the exponential format of a root. The exponential format is relative to a root of the default primitive polynomial for  $\text{GF}(2^m)$ .

`rt = groots(f, primpoly)` finds roots in  $\text{GF}(2^m)$  of the polynomial that  $f$  represents.  $rt$  is a column vector each of whose entries is the exponential format of a root. The exponential format is relative to a root of the degree- $m$  primitive polynomial for  $\text{GF}(2^m)$  that `primpoly` represents.  $m$  is an integer greater than or equal to  $d$ .

`rt = groots(f, m, p)` is the same as `rt = groots(f, m)` except that 2 is replaced by a prime number  $p$ .

`rt = groots(f, primpoly, p)` is the same as `rt = groots(f, primpoly)` except that 2 is replaced by a prime number  $p$ .



`[rt, rt_tuple] = groots(...)` returns an additional matrix `rt_tuple`, whose  $k$ th row is the polynomial format of the root  $rt(k)$ . The polynomial and exponential formats are both relative to the same primitive element.

`[rt, rt_tuple, field] = groots(...)` returns additional matrices `rt_tuple` and `field`. `rt_tuple` is described in the paragraph above. `field` gives the list of elements of the extension field. The list of elements, the polynomial format, and the exponential format are all relative to the same primitive element.

---

**Note** For a description of the various formats that `groots` uses, see “Representing Elements of Galois Fields” on page 2-90.

---

## Examples

The section, “Roots of Polynomials” on page 2-102, contains a description and example of the use of `groots`.

As another example, the code below finds the polynomial format of the roots of the primitive polynomial  $1 + x^3 + x^4$  for GF(16). It then displays the roots in traditional form as polynomials in `alpha`. Since `primpoly` is both the primitive polynomial and the polynomial whose roots are sought, `alpha` itself is a root.

```
p = 2; m = 4;
primpoly = [1 0 0 1 1]; % A primitive polynomial for GF(16)
f = primpoly; % Find roots of the primitive polynomial.
[rt, rt_tuple] = groots(f, primpoly, p);
% Display roots as polynomials in alpha.
for ii = 1:length(rt_tuple)
    gfpretty(rt_tuple(ii,:), 'alpha')
end
```

## See Also

`gfpri mdf`, `gfl i neq`

**Purpose** Subtract polynomials over a Galois field

**Syntax**

```
c = gfsb(a, b);  
c = gfsb(a, b, p);  
c = gfsb(a, b, p, len);  
c = gfsb(a, b, field);
```

**Description** `c = gfsb(a, b)` calculates  $a$  minus  $b$ , where  $a$  and  $b$  represent polynomials over  $GF(2)$ . The inputs and output are row vectors that give the coefficients of the corresponding polynomials in order of ascending powers. Each coefficient is either 0 or 1, since the field is  $GF(2)$ . If  $a$  and  $b$  are matrices of the same size, then the function treats each row independently.

`c = gfsb(a, b, p)` calculates  $a$  minus  $b$ , where  $a$  and  $b$  represent polynomials over  $GF(p)$  and  $p$  is a prime number.  $a$ ,  $b$ , and  $c$  are row vectors that give the coefficients of the corresponding polynomials in order of ascending powers. Each coefficient is between 0 and  $p-1$ . If  $a$  and  $b$  are matrices of the same size, then the function treats each row independently.

`c = gfsb(a, b, p, len)` subtracts row vectors as in the syntax above, except that it returns a row vector of length `len`. The output  $c$  is a truncated or extended representation of the answer. If the row vector corresponding to the answer has fewer than `len` entries (including zeros), then extra zeros are added at the end; if it has more than `len` entries, then entries from the end are removed.

`c = gfsb(a, b, field)` calculates  $a$  minus  $b$ , where  $a$  and  $b$  are the exponential format of two elements of  $GF(p^m)$ , relative to some primitive element of  $GF(p^m)$ .  $p$  is a prime number and  $m$  is a positive integer. `field` is the matrix listing all elements of  $GF(p^m)$ , arranged relative to the same primitive element.  $c$  is the exponential format of the answer, relative to the same primitive element. See “Representing Elements of Galois Fields” on page 2-90 for an explanation of these formats. If  $a$  and  $b$  are matrices of the same size, then the function treats each element independently.

**Examples** In the code below, `diff` is the difference of  $2 + 3x + x^2$  and  $4 + 2x + 3x^2$  over  $GF(5)$ , and `linpart` is the degree-one part of `diff`.

```
diff = gfsb([2 3 1], [4 2 3], 5)
```

```
diff =
```

```
      3      1      3
```

```
linpart = gfsub([2 3 1], [4 2 3], 5, 2)
```

```
linpart =
```

```
      3      1
```

The code below shows that  $\alpha^2 - \alpha^4 = \alpha^7$ , where  $\alpha$  is a root of the primitive polynomial  $2 + 2x + x^2$  for GF(9).

```
p = 3; m = 2;
primpoly = [2 2 1];
field = gftuple([-1: p^m-2]', primpoly, p);
d = gfsub(2, 4, field)
```

```
d =
```

```
      7
```

## See Also

gfadd, gfmul, gfconv, gfdi v, gfdeconv, gftuple

# gftrunc

---

**Purpose** Minimize the length of a polynomial representation

**Syntax** `c = gftrunc(a);`

**Description** `c = gftrunc(a)` truncates a row vector, `a`, that gives the coefficients of a  $\text{GF}(p)$  polynomial in order of ascending powers. If  $a(k) = 0$  whenever  $k > d + 1$ , then the polynomial has degree  $d$ . The row vector `c` omits these high-order zeros and thus has length  $d + 1$ .

**Examples** In the code below, zeros are removed from the end, but *not* from the beginning or middle, of the row-vector representation of  $x^2 + 2x^3 + 3x^4 + 4x^7 + 5x^8$ .

```
c = gftrunc([0 0 1 2 3 0 0 4 5 0 0])
```

```
c =
```

```
0    0    1    2    3    0    0    4    5
```

**See Also** `gfadd`, `gfconv`, `gfsub`, `gfdeconv`, `gftuple`

**Purpose** Simplify or convert the format of elements of a Galois field

**Syntax**

```
tp = gftuple(a, m);
tp = gftuple(a, primpoly);
tp = gftuple(a, m, p);
tp = gftuple(a, primpoly, p);
tp = gftuple(a, primpoly, p, primck);
[tp, expform] = gftuple(...);
```

**Description** For All Syntaxes

`gftuple` serves to simplify the polynomial or exponential format of Galois field elements, or to convert from one format to another. For an explanation of the formats that `gftuple` uses, see “Representing Elements of Galois Fields” on page 2-90.

In this discussion, the format of an element of  $GF(p^m)$  is called “simplest” if all exponents of the primitive element are:

- Between 0 and  $m-1$  for the polynomial format
- Either  $-Inf$ , or between 0 and  $p^{m-2}$  for the exponential format

For all syntaxes, `a` is a matrix, each row of which represents an element of a Galois field. The format of `a` determines how MATLAB interprets it:

- If `a` is a column of integers, then MATLAB interprets each row as an *exponential* format of an element. Negative integers are equivalent to  $-Inf$  in that they all represent the zero element of the field.
- If `a` has more than one column, then MATLAB interprets each row as a *polynomial* format of an element. (Each entry of `a` must be an integer between 0 and  $p-1$ , where  $p$  is 2 if not specified as an input.)

The exponential or polynomial formats mentioned above are all relative to a primitive element specified by the *second* input argument. The second argument is described below.

**For Specific Syntaxes**

`tp = gftuple(a, m)` returns the simplest polynomial format of the elements that `a` represents, where the  $k$ th row of `tp` corresponds to the  $k$ th row of `a`. The

formats are relative to a root of the default primitive polynomial for  $\text{GF}(2^m)$ .  $m$  is a positive integer. If possible, the default primitive polynomial is used to simplify the polynomial formats.

`tp = gftuple(a, primpoly)` returns the simplest polynomial format of the element that `a` represents, where the  $k$ th row of `tp` corresponds to the  $k$ th row of `a`. The formats are relative to a root of the primitive polynomial whose coefficients are given, in order of ascending powers, by the row vector `primpoly`. If possible, this primitive polynomial is used to simplify the polynomial formats.

`tp = gftuple(a, m, p)` is the same as `tp = gftuple(a, m)` except that 2 is replaced by a prime number  $p$ .

`tp = gftuple(a, primpoly, p)` is the same as `tp = gftuple(a, primpoly)` except that 2 is replaced by a prime number  $p$ .

`tp = gftuple(a, primpoly, p, prim_ck)` is the same as `tp = gftuple(a, primpoly, p)` except that `gftuple` checks whether `primpoly` represents a polynomial that is indeed primitive. If not, then `gftuple` generates an error and `tp` is not returned. The input argument `prim_ck` can be any number or string; only its existence matters.

`[tp, expform] = gftuple(...)` returns the additional matrix `expform`. The  $k$ th row of `expform` is the simplest exponential format of the element that the  $k$ th row of `a` represents. All other features are as described in earlier parts of this “Description” section, depending on the input arguments.

## Examples

Some examples are in these subsections of “Galois Field Computations” on page 2-89:

- “List of All Elements of a Galois Field” on page 2-91 (end of section)
- “Converting to Simplest Polynomial Format” on page 2-94
- “Converting to Simplest Exponential Format” on page 2-96

As another example, the `gftuple` command below generates a list of elements of  $\text{GF}(p^m)$ , arranged relative to a root of the default primitive polynomial. Some functions in this toolbox use such a list as an input argument.

```
p = 5; % Or any prime number
m = 4; % Or any positive integer
field = gftuple([-1:p^m-2]', m, p);
```

Finally, the two commands below illustrate the influence of the *shape* of the input matrix. In the first command, a column vector is treated as a sequence of elements expressed in exponential format. In the second command, a row vector is treated as a single element expressed in polynomial format.

```
tp1 = gftuple([0; 1], 3)
```

```
tp1 =
```

```
1    0    0
0    1    0
```

```
tp2 = gftuple([0, 0, 0, 1], 3)
```

```
tp2 =
```

```
1    1    0
```

The outputs reflect that, according to the default primitive polynomial for GF(8), the relations below are true.

$$\alpha^0 = 1 + 0\alpha + 0\alpha^2$$

$$\alpha^1 = 0 + 1\alpha + 0\alpha^2$$

$$0 + 0\alpha + 0\alpha^2 + \alpha^3 = 1 + \alpha + 0\alpha^2$$

### Algorithm

`gftuple` uses recursive callbacks to determine the exponential format.

### See Also

`gfadd`, `gfmul`, `gfconv`, `gfdiv`, `gfdeconv`, `gfpri mdf`

# gfweight

---

<b>Purpose</b>	Calculate the minimum distance of a linear block code
<b>Syntax</b>	<pre>wt = gfweight(genmat); wt = gfweight(genmat, 'gen'); wt = gfweight(parmat, 'par'); wt = gfweight(genpoly, n);</pre>
<b>Description</b>	<p>The minimum distance, or minimum weight, of a linear block code is defined as the smallest positive number of nonzero entries in any <math>n</math>-tuple that is a codeword.</p> <p><code>wt = gfweight(genmat)</code> returns the minimum distance of the linear block code whose generator matrix is <code>genmat</code>.</p> <p><code>wt = gfweight(genmat, 'gen')</code> returns the minimum distance of the linear block code whose generator matrix is <code>genmat</code>.</p> <p><code>wt = gfweight(parmat, 'par')</code> returns the minimum distance of the linear block code whose parity-check matrix is <code>parmat</code>.</p> <p><code>wt = gfweight(genpoly, n)</code> returns the minimum distance of the <i>cyclic</i> code whose codeword length is <code>n</code> and whose generator polynomial is represented by <code>genpoly</code>. <code>genpoly</code> is a row vector that gives the coefficients of the generator polynomial in order of ascending powers.</p>
<b>Examples</b>	<p>The commands below illustrate three different ways to compute the minimum distance of a (7,4) cyclic code.</p> <pre>n = 7; % Generator polynomial of (7,4) cyclic code genpoly = cyclpoly(n, 4); [parmat, genmat] = cyclgen(n, genpoly); wts = [gfweight(genmat, 'gen'), gfweight(parmat, 'par'), ...        gfweight(genpoly, n)]  wts =        3      3      3</pre>
<b>See Also</b>	<code>hammgen</code> , <code>bchpoly</code> , <code>cyclpoly</code>



<b>Purpose</b>	Produce parity-check and generator matrices for Hamming code
<b>Syntax</b>	<pre> h = hammgen(m); h = hammgen(m, pol); [h, g] = hammgen(...); [h, g, n, k] = hammgen(...); </pre>
<b>Description</b>	<p>For all syntaxes, the codeword length is <math>n</math>. <math>n</math> has the form <math>2^m-1</math> for some positive integer <math>m</math> greater than or equal to 3. The message length, <math>k</math>, has the form <math>n-m</math>.</p> <p><code>h = hammgen(m)</code> produces an <math>m</math>-by-<math>n</math> parity-check matrix for a Hamming code having codeword length <math>n = 2^m-1</math>. <math>m</math> is a positive integer greater than or equal to 3. The message length of the code is <math>n-m</math>. The binary primitive polynomial used to produce the Hamming code is MATLAB's default primitive polynomial for <math>GF(2^m)</math>, represented by <code>gfprimdf(m)</code>.</p> <p><code>h = hammgen(m, pol)</code> produces an <math>m</math>-by-<math>n</math> parity-check matrix for a Hamming code having codeword length <math>n = 2^m-1</math>. <math>m</math> is a positive integer greater than or equal to 3. The message length of the code is <math>n-m</math>. <code>pol</code> is a row vector that gives the coefficients, in order of ascending powers, of the binary primitive polynomial for <math>GF(2^m)</math> that is used to produce the Hamming code. <code>hammgen</code> produces an error if <code>pol</code> represents a polynomial that is not, in fact, primitive.</p> <p><code>[h, g] = hammgen(...)</code> is the same as <code>h = hammgen(...)</code> except that it also produces the <math>k</math>-by-<math>n</math> generator matrix <code>g</code> that corresponds to the parity-check matrix <code>h</code>. <math>k</math>, the message length, equals <math>n-m</math>, or, <math>2^m-1-m</math>.</p> <p><code>[h, g, n, k] = hammgen(...)</code> is the same as <code>[h, g] = hammgen(...)</code> except that it also returns the codeword length <math>n</math> and the message length <math>k</math>.</p>

---

**Note** If your value of  $m$  is less than 25 and if your primitive polynomial is MATLAB's default primitive polynomial for  $GF(2^m)$ , then the syntax `hammgen(m)` is likely to be faster than the syntax `hammgen(m, pol)`.

---

<b>Examples</b>	The command below exhibits the parity-check and generator matrices for a Hamming code with codeword length $7 = 2^3-1$ and message length $4 = 7-3$ .
-----------------	---

# hammgen

---

```
[h, g, n, k] = hammgen(3)
```

h =

1	0	0	1	0	1	1
0	1	0	1	1	1	0
0	0	1	0	1	1	1

g =

1	1	0	1	0	0	0
0	1	1	0	1	0	0
1	1	1	0	0	1	0
1	0	1	0	0	0	1

n =

7

k =

4

The command below, which uses  $1 + x^2 + x^3$  as the primitive polynomial for  $\text{GF}(2^3)$ , shows that the parity-check matrix depends on the choice of primitive polynomial. Notice that h1 below is different from h in the example above.

```
h1 = hammgen(3, [1 0 1 1])
```

h1 =

1	0	0	1	1	1	0
0	1	0	0	1	1	1
0	0	1	1	1	0	1

<b>Algorithm</b>	Unlike <code>gftuple</code> which processes one $m$ -tuple at a time, <code>hammgen</code> generates the entire sequence from 0 to $2^m-1$ . The computation algorithm uses all previously computed values to produce the computation result.
<b>See Also</b>	<code>gftuple</code> , <code>gfrepconv</code> , <code>gfpri mck</code> , <code>gfpri mfd</code> , <code>gfpri mdf</code>

# hank2sys

---

**Purpose** Convert a Hankel matrix to a linear system model

**Syntax**

```
[num, den] = hank2sys(h, i ni , tol)
[num, den, sv] = hank2sys(h, i ni , tol)
[a, b, c, d] = hank2sys(h, i ni , tol)
[a, b, c, d, sv] = hank2sys(h, i ni , tol)
```

**Description** [num, den] = hank2sys(h, i ni , tol) converts a Hankel matrix h to a linear system transfer function with numerator num and denominator den. The vectors num and den list the coefficients of their respective polynomials in order of ascending exponents. i ni is the system impulse at time zero. If tol > 1, then tol is the order of the conversion. If tol < 1, then tol is the tolerance in selecting the conversion order based on the singular values. If you omit tol, then its default value is 0.01. This conversion uses the singular value decomposition method.

[num, den, sv] = hank2sys(h, i ni , tol) is the same as the syntax above, except that sv is a vector that lists the singular values of h.

[a, b, c, d] = hank2sys(h, i ni , tol) converts a Hankel matrix h to a corresponding linear system state-space model. a, b, c, and d are matrices. The input parameters are the same as in the first syntax above.

[a, b, c, d, sv] = hank2sys(h, i ni , tol) is the same as the syntax above, except that sv is a vector that lists the singular values of h.

**Examples**

```
h = hankel ([1 0 1]);
[num, den, sv] = hank2sys(h, 0, .01)
```

num =

```
0      1. 0000   -0. 0000    1. 0000
```

den =

```
1. 0000    0. 0000    0. 0000    0. 0000
```

`SV =`

1.6180

1.0000

0.6180

**See Also**

`hilb`, `ir`, `hankel`, `rcosflt`

**Purpose** Design a Hilbert transform IIR filter

**Syntax**

```
hilbiir;  
hilbiir(ts);  
hilbiir(ts, dly);  
hilbiir(ts, dly, bandwidth);  
hilbiir(ts, dly, bandwidth, tol);  
[num, den] = hilbiir(...);  
[num, den, sv] = hilbiir(...);  
[a, b, c, d] = hilbiir(...);  
[a, b, c, d, sv] = hilbiir(...);
```

**Description** The function `hilbiir` designs a Hilbert transform filter. The output is either:

- A plot of the filter's impulse response, or
- A quantitative characterization of the filter, using either a transfer function model or a state-space model

### Background Information

An ideal Hilbert transform filter has the transfer function  $H(s) = -j\text{sgn}(s)$ , where  $\text{sgn}(\cdot)$  is the signum function (`sign` in MATLAB). The impulse response of the Hilbert transform filter is

$$h(t) = \frac{1}{\pi t}$$

Since the Hilbert transform filter is a noncausal filter, the `hilbiir` function introduces a group delay, `dly`. A Hilbert transform filter with this delay has the impulse response

$$h(t) = \frac{1}{\pi(t - \text{dly})}$$

### Choosing a Group Delay Parameter

The filter design is an approximation. If you provide the filter's group delay as an input argument, then these two suggestions can help improve the accuracy of the results:

- Choose the sample time  $t_s$  and the filter's group delay  $dly$  so that  $dly$  is at least a few times larger than  $t_s$  and  $\text{rem}(dly, t_s) = t_s/2$ . For example, you can set  $t_s$  to  $2*dly/N$ , where  $N$  is a positive integer.
- At the point  $t = dly$ , the impulse response of the Hilbert transform filter can be interpreted as 0,  $-\text{Inf}$ , or  $\text{Inf}$ . If `hilbiir` encounters this point, then it sets the impulse response there to zero. To improve accuracy, avoid the point  $t = dly$ .

### Syntaxes for Plots

Each of these syntaxes produces a plot of the impulse response of the filter that the `hilbiir` function designs, as well as the impulse response of a corresponding ideal Hilbert transform filter.

`hilbiir` plots the impulse response of a fourth-order digital Hilbert transform filter with a 1-second group delay. The sample time is  $2/7$  seconds. In this particular design, the tolerance index is 0.05. The plot also displays the impulse response of the ideal Hilbert transform filter with a 1-second group delay.

`hilbiir(ts)` plots the impulse response of a fourth-order Hilbert transform filter with a sample time of  $t_s$  seconds and a group delay of  $t_s*7/2$  seconds. The tolerance index is 0.05. The plot also displays the impulse response of the ideal Hilbert transform filter having a sample time of  $t_s$  seconds and a group delay of  $t_s*7/2$  seconds.

`hilbiir(ts, dly)` is the same as the syntax above, except that the filter's group delay is  $dly$  for both the ideal filter and the filter that `hilbiir` designs. See "Choosing a Group Delay Parameter" above for guidelines on choosing  $dly$ .

`hilbiir(ts, dly, bandwdth)` is the same as the syntax above, except that `bandwdth` specifies the assumed bandwidth of the input signal and that the filter design might use a compensator for the input signal. If `bandwdth` = 0 or `bandwdth` >  $1/(2*t_s)$ , then `hilbiir` does not use a compensator.

`hilbiir(ts, dly, bandwdth, tol)` is the same as the syntax above, except that `tol` is the tolerance index. If `tol` < 1, then the order of the filter is determined by

$$\frac{\text{truncated-singular-value}}{\text{maximum-singular-value}} < \text{tol}$$

If  $\text{tol} > 1$ , then the order of the filter is  $\text{tol}$ .

## Syntaxes for Transfer Function and State-Space Quantities

Each of these syntaxes produces quantitative information about the filter that `hilbiir` designs, but does *not* produce a plot. The input arguments for these syntaxes (if you provide any) are the same as those described in the “Syntaxes for Plots” section above.

`[num, den] = hilbiir(...)` outputs the numerator and denominator of the IIR filter's transfer function.

`[num, den, sv] = hilbiir(...)` outputs the numerator and denominator of the IIR filter's transfer function, and the singular values of the Hankel matrix that `hilbiir` uses in the computation.

`[a, b, c, d] = hilbiir(...)` outputs the discrete-time state-space model of the designed Hilbert transform filter. `a`, `b`, `c`, and `d` are matrices.

`[a, b, c, d, sv] = hilbiir(...)` outputs the discrete-time state-space model of the designed Hilbert transform filter, and the singular values of the Hankel matrix that `hilbiir` uses in the computation.

## Algorithm

The `hilbiir` function calculates the impulse response of the ideal Hilbert transform filter response with a group delay. It fits the response curve using a singular-value decomposition method. See the book by Kailath listed below.

## Examples

At the MATLAB prompt, type `hilbiir` or `[num, den] = hilbiir` for an example using the function's default values.

## See Also

`grpdelay` (Signal Processing Toolbox)

## References

Kailath, Thomas. *Linear Systems*. Englewood Cliffs, N.J.: Prentice-Hall, 1980.



<b>Purpose</b>	Check if the input is a valid trellis structure
<b>Syntax</b>	<code>[i sok, status] = istrellis(s);</code>
<b>Description</b>	<p><code>[i sok, status] = istrellis(s)</code> checks if the input <code>s</code> is a valid trellis structure. If the input is a valid trellis structure, then <code>i sok</code> is 1 and <code>status</code> is an empty string. Otherwise, <code>i sok</code> is 0 and <code>status</code> is a string that indicates why <code>s</code> is not a valid trellis structure.</p> <p>A valid trellis structure is a MATLAB structure whose fields are as in the table below.</p>

**Table 3-16: Fields of a Valid Trellis Structure for a Rate  $k/n$  Code**

Field in Trellis Structure	Dimensions	Meaning
<code>numInputSymbol s</code>	Scalar	Number of input symbols to the encoder: $2^k$
<code>numOutputsymbol s</code>	Scalar	Number of output symbols from the encoder: $2^n$
<code>numStates</code>	Scalar	Number of states in the encoder
<code>nextStates</code>	<code>numStates-by-<math>2^k</math></code> matrix	Next states for all combinations of current state and current input
<code>outputs</code>	<code>numStates-by-<math>2^k</math></code> matrix	Outputs (in octal) for all combinations of current state and current input

In the `nextStates` matrix, each entry is an integer between 0 and `numStates-1`. The element in the  $s$ th row and  $u$ th column denotes the next state when the starting state is  $s-1$  and the input bits have decimal representation  $u-1$ . To convert the input bits to a decimal value, use the first input bit as the most significant bit (MSB). For example, the second column of the `nextStates` matrix stores the next states when the current set of input values is  $\{0, \dots, 0, 1\}$ .

To convert the state to a decimal value, use this rule: If  $k$  exceeds 1, then the shift register that receives the first input stream in the encoder provides the least significant bits in the state number, while the shift register that receives the last input stream in the encoder provides the most significant bits in the state number.

In the `outputs` matrix, the element in the  $s$ th row and  $u$ th column denotes the encoder's output when the starting state is  $s-1$  and the input bits have decimal representation  $u-1$ . To convert to decimal value, use the first output bit as the MSB.

## Examples

These commands assemble the fields into a very simple trellis structure, and then verify the validity of the trellis structure.

```
trellis.numInputSymbols = 2;  
trellis.numOutputSymbols = 2;  
trellis.numStates = 2;  
trellis.nextStates = [0 1; 0 1];  
trellis.outputs = [0 0; 1 1];  
[isok, status] = istrellis(trellis)
```

```
isok =
```

```
1
```

```
status =
```

```
''
```

## See Also

`poly2trellis`, `struct`, `convenc`, `vitdec`

<b>Purpose</b>	Optimize quantization parameters using the Lloyd algorithm
<b>Syntax</b>	<pre> [partition, codebook] = lloyds(trainingset, initcodebook); [partition, codebook] = lloyds(trainingset, length); [partition, codebook] = lloyds(trainingset, ..., tol); [partition, codebook] = lloyds(trainingset, ..., tol, plotflag); [partition, codebook, distor] = lloyds(...); [partition, codebook, distor, reldistor] = lloyds(...); </pre>
<b>Description</b>	<p><code>[partition, codebook] = lloyds(trainingset, initcodebook)</code> optimizes the scalar quantization parameters <code>partition</code> and <code>codebook</code> for the training data in the vector <code>trainingset</code>. <code>initcodebook</code>, a vector of length at least 2, is the initial guess of the codebook values. The output <code>codebook</code> is a vector of the same length as <code>initcodebook</code>. The output <code>partition</code> is a vector whose length is one less than the length of <code>codebook</code>.</p> <p>See either “Representing Quantization Parameters” on page 2-14 or the reference page for <code>quantiz</code> in this chapter, for a description of the formats of <code>partition</code> and <code>codebook</code>.</p> <hr/> <p><b>Note</b> <code>lloyds</code> optimizes for the data in <code>trainingset</code>. For best results, <code>trainingset</code> should be similar to the data that you plan to quantize.</p> <hr/> <p><code>[partition, codebook] = lloyds(trainingset, length)</code> is the same as the first syntax, except that the scalar argument <code>length</code> indicates the size of the vector <code>codebook</code>. This syntax does not include an initial codebook guess.</p> <p><code>[partition, codebook] = lloyds(trainingset, ..., tol)</code> is the same as the two syntaxes above, except that <code>tol</code> replaces <math>10^{-7}</math> in condition 1 of the algorithm description below.</p> <p><code>[partition, codebook] = lloyds(trainingset, ..., tol, plotflag)</code> is the same as the syntax above, except that it also plots the original signal, the optimized <code>partition</code>, and <code>codebook</code> in a figure. The value of <code>plotflag</code> is not important.</p>

`[partition, codebook, distort] = lloyds(...)` returns the final mean square distortion in the variable `distort`.

`[partition, codebook, distort, reldistort] = lloyds(...)` returns a value `reldistort` that is related to the algorithm's termination. In case 1 of "Algorithm" below, `reldistort` is the relative change in distortion between the last two iterations. In case 2, `reldistort` is the same as `distort`.

## Examples

The code below optimizes the quantization parameters for a sinusoidal transmission via a 3-bit channel. Since the typical data is sinusoidal, `trainingset` is a sampled sine wave. Since the channel can transmit 3 bits at a time, `lloyds` prepares a codebook of length  $2^3$ .

```
% Generate a complete period of a sinusoidal signal.
x = sin([0:1000]*pi/500);
[partition, codebook] = lloyds(x, 2^3)

partition =

    -0.8540    -0.5973    -0.3017     0.0031     0.3077     0.6023     0.8572

codebook =

Columns 1 through 7

    -0.9504    -0.7330    -0.4519    -0.1481     0.1558     0.4575     0.7372

Column 8

     0.9515
```

## Algorithm

`lloyds` uses an iterative process to try to minimize the mean square distortion. The optimization processing ends when either:

- 1 The relative change in distortion between iterations is less than  $10^{-7}$ , or
- 2 The distortion is less than `eps*max(trainingset)`, where `eps` is MATLAB's floating-point relative accuracy

## See Also

`compand`, `dpcmopt`, `quantiz`

**References**

S. P. Lloyd. "Least Squares Quantization in PCM." *IEEE Transactions on Information Theory*. Vol IT-28, March 1982, 129-137.

J. Max. "Quantizing for Minimum Distortion." *IRE Transactions on Information Theory*. Vol. IT-6, March 1960, 7-12.

# marcumq

---

**Purpose** Generalized Marcum Q function

**Syntax**  $Q = \text{marcumq}(a, b);$   
 $Q = \text{marcumq}(a, b, m);$

**Description**  $Q = \text{marcumq}(a, b)$  computes the Marcum Q function of  $a$  and  $b$ , defined by

$$Q(a, b) = \int_b^{\infty} x \exp\left(-\frac{x^2 + a^2}{2}\right) I_0(ax) dx$$

where  $a$  and  $b$  are nonnegative real numbers. In this expression,  $I_0$  is the modified Bessel function of the first kind of zero order.

$Q = \text{marcumq}(a, b, m)$  computes the generalized Marcum Q, defined by

$$Q_m(a, b) = \frac{1}{a^{m-1}} \int_b^{\infty} x^m \exp\left(-\frac{x^2 + a^2}{2}\right) I_{m-1}(ax) dx$$

where  $a$  and  $b$  are nonnegative real numbers, and  $m$  is a nonnegative integer. In this expression,  $I_{m-1}$  is the modified Bessel function of the first kind of order  $m-1$ .

**See Also** `bessel i`; `ncx2cdf` (Statistics Toolbox)

**References** Cantrell, P. E. and A. K. Ojha, "Comparison of Generalized Q-Function Algorithms." *IEEE Transactions on Information Theory*, vol. IT-33, July 1987, 591-596.

Marcum, J. I. "A Statistical Theory of Target Detection by Pulsed Radar: Mathematical Appendix." RAND Corporation, Santa Monica, CA, Research Memorandum RM-753, July 1, 1948. Reprinted in *IRE Transactions on Information Theory*, vol. IT-6, April 1960, 59-267.

McGee, W. F. "Another Recursive Method of Computing the Q Function." *IEEE Transactions on Information Theory*, vol. IT-16, July 1970, 500-501.

**Purpose** Map a digital signal to an analog signal

**Syntax**

```
modmap('method',...);
y = modmap(x, Fd, Fs, 'ask', M);
y = modmap(x, Fd, Fs, 'fsk', M, tone);
y = modmap(x, Fd, Fs, 'msk');
y = modmap(x, Fd, Fs, 'psk', M);
y = modmap(x, Fd, Fs, 'qask', M);
y = modmap(x, Fd, Fs, 'qask/arb', inphase, quadr);
y = modmap(x, Fd, Fs, 'qask/cir', numsig, amp, phs);
```

Optional Inputs	Input	Default Value
	tone	Fd
	amp	[ 1:length(numsig) ]
	phs	numsig*0

**Description** The digital modulation process consists of two steps: mapping the digital signal to an analog signal and modulating this analog signal. The function `modmap` performs the first step. You can perform the second step using `amod`, `amodce`, or your own custom modulator. The table below lists the digital modulation schemes that `modmap` supports.

Modulation Scheme	Value of 'method'
M-ary amplitude shift keying	'ask'
M-ary frequency shift keying	'fsk'
Minimum shift keying	'msk'
M-ary phase shift keying	'psk'
Quadrature amplitude shift keying	'qask', 'qask/cir', or 'qask/arb'

To Plot a Signal Constellation

`modmap('method',...)` creates a plot that characterizes the M-ary modulation method that `'method'` specifies. `'method'` is one of the entries in the

right-hand column of the table above. If '*method*' is a value other than '**fsk**' or '**msk**', then the plot shows the signal constellation; otherwise, it shows the spectrum.

For most methods, the input parameters that follow '*method*' in this syntax are the same as those that follow '*method*' in the corresponding mapping syntax. For more information about them, see the section “To Map a Digital Signal (Specific Syntax Information)” below.

However, if '*method*' is '**msk**', then the syntax is

```
modmap('msk', Fd)
```

where  $F_d$  is the sampling rate of the message signal.

## To Map a Digital Signal (General Information)

The generic syntax  $y = \text{modmap}(x, F_d, F_s, \dots)$  maps the digital message signal  $x$  onto an analog signal.  $x$  is a matrix of nonnegative integers. The sizes of  $x$  and  $y$  depend on the modulation method:

- **(ASK, FSK, MSK methods)** If  $x$  is a vector of length  $n$ , then  $y$  is a column vector of length  $n \cdot F_s / F_d$ . Otherwise, if  $x$  is  $n$ -by- $m$ , then  $y$  is  $(n \cdot F_s / F_d)$ -by- $m$  and each column of  $x$  is processed separately.
- **(PSK, QASK methods)** If  $x$  is a vector of length  $n$ , then  $y$  is an  $n \cdot F_s / F_d$ -by-2 matrix. Otherwise, if  $x$  is  $n$ -by- $m$ , then  $y$  is  $(n \cdot F_s / F_d)$ -by- $2m$  and each column of  $x$  is processed separately. The odd-numbered columns in  $y$  represent in-phase components and the even-numbered columns represent quadrature components.

The sampling rates in Hertz of  $x$  and  $y$ , respectively, are  $F_d$  and  $F_s$ . (Thus  $1/F_d$  represents the time interval between two consecutive samples in  $x$ , and similarly for  $y$ .) The ratio  $F_s/F_d$  must be a positive integer.

## To Map a Digital Signal (Specific Syntax Information)

$y = \text{modmap}(x, F_d, F_s, \text{'ask'}, M)$  maps to an  $M$ -ary amplitude shift keying signal constellation. Each entry of  $x$  must be in the range  $[0, M-1]$ . Each entry of  $y$  is in the range  $[-1, 1]$ .

$y = \text{modmap}(x, F_d, F_s, \text{'fsk'}, M, \text{tone})$  maps to frequencies in an  $M$ -ary frequency shift keying set. Each entry of  $x$  must be in the range  $[0, M-1]$ . The



optional argument `tone` is the separation between successive frequencies in the FSK set. The default value of `tone` is `Fd`.

`y = modmap(x, Fd, Fs, 'msk')` maps to frequencies in a minimum shift keying set. Each entry of `x` is either 0 or 1. The separation between the two frequencies is `Fd/2`.

`y = modmap(x, Fd, Fs, 'psk', M)` maps to an `M`-ary phase shift keying signal constellation. Each entry of `x` must be in the range `[0, M-1]`.

`y = modmap(x, Fd, Fs, 'qask', M)` maps to an `M`-ary quadrature amplitude shift keying square signal constellation. The table below shows the maximum value of the in-phase and quadrature components in `y`, for several small values of `M`.

M	Maximum of y	M	Maximum of y
2	1	32	5
4	1	64	7
8	3 (quadrature maximum is 1)	128	11
16	3	256	15

**Note** To see how symbols are mapped to the constellation points, generate a square constellation plot using `qaskenco(M)` or `modmap('qask', M)`.

`y = modmap(x, Fd, Fs, 'qask/arb', inphase, quadr)` maps to a quadrature amplitude shift keying signal constellation that you define using the vectors `inphase` and `quadr`. The signal constellation point for the  $k$ th message has in-phase component `inphase(k+1)` and quadrature component `quadr(k+1)`.

`y = modmap(x, Fd, Fs, 'qask/cir', numsig, amp, phs)` maps to a quadrature amplitude shift keying circular signal constellation. `numsig`, `amp`, and `phs` are vectors of the same length. The entries in `numsig` and `amp` must be positive. If  $k$  is an integer in the range `[1, length(numsig)]`, then `amp(k)` is the radius of

the  $k$ th circle,  $\text{numsig}(k)$  is the number of constellation points on the  $k$ th circle, and  $\text{phs}(k)$  is the phase of the first constellation point plotted on the  $k$ th circle. All points on the  $k$ th circle are evenly spaced. If you omit  $\text{phs}$ , then its default value is  $\text{numsig} \times 0$ . If you omit  $\text{amp}$ , then its default value is  $[1: \text{length}(\text{numsig})]$ .

---

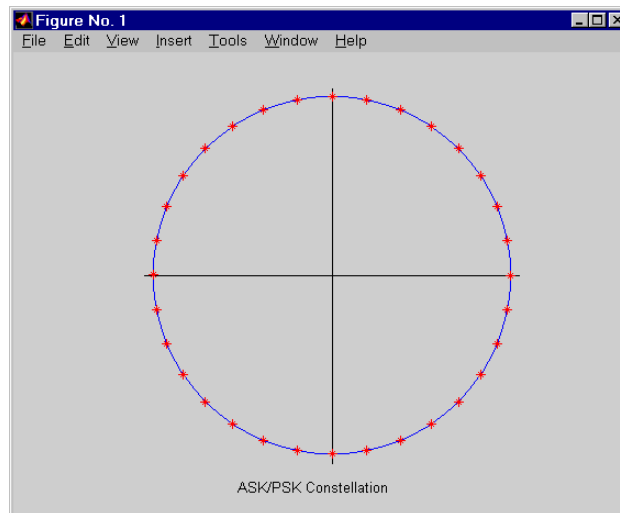
**Note** To see how symbols are mapped to the constellation points, generate a labeled circle constellation plot using `apkconst(numsig, amp, phs, 'n')`.

---

## Examples

The command below plots a phase shift keying (PSK) signal constellation with 32 points.

```
modmap('psk', 32);
```



The script below maps a digital signal using the 32-point PSK constellation. It then adds noise and computes the resulting error rate while demapping. Your results might vary because the example uses random numbers.

```
M = 32; Fd = 1; Fs = 3;  
x = randint(100, 1, M); % Original signal
```

```
y = modmap(x, Fd, Fs, 'psk', M); % Mapped signal, using 32-ary PSK
ynoi sy = y+.1*rand(100*Fs, 2); % Mapped signal with noise added
z = demodmap(ynoi sy, Fd, Fs, 'psk', M); % Demapped noisi y signal
s = symerr(x, z) % Number of errors after demapping noisi y signal
```

s =

8

## See Also

demodmap, dmod, dmodce, amod, amodce, apkconst

## oct2dec

---

<b>Purpose</b>	Convert octal numbers to decimal numbers				
<b>Syntax</b>	<code>d = oct2dec(c)</code>				
<b>Description</b>	<code>d = oct2dec(c)</code> converts an octal matrix <code>c</code> to a decimal matrix <code>d</code> , element by element. In both octal and decimal representations, the rightmost digit is the least significant.				
<b>Examples</b>	<p>The command below converts a 2-by-2 octal matrix.</p> <pre>d = oct2dec([12 144; 0 25])</pre> <p>d =</p> <table><tr><td>10</td><td>100</td></tr><tr><td>0</td><td>21</td></tr></table> <p>For instance, the octal number 144 is equivalent to the decimal number 100 because <math>144 \text{ (octal)} = 1*8^2 + 4*8^1 + 4*8^0 = 64 + 32 + 4 = 100</math>.</p>	10	100	0	21
10	100				
0	21				
<b>See Also</b>	<code>bi2de</code>				

<b>Purpose</b>	Convert convolutional code polynomials to trellis description
<b>Syntax</b>	<pre>trellis = poly2trellis(ConstraintLength, CodeGenerator); trellis = poly2trellis(ConstraintLength, CodeGenerator, ...     FeedbackConnection);</pre>
<b>Description</b>	<p>The <code>poly2trellis</code> function accepts a polynomial description of a convolutional encoder and returns the corresponding trellis structure description. The output of <code>poly2trellis</code> is suitable as an input to the <code>convenc</code> and <code>vitdec</code> functions, and as a mask parameter for the Convolutional Encoder, Viterbi Decoder, and APP Decoder blocks in the Communications Blockset.</p> <p><code>trellis = poly2trellis(ConstraintLength, CodeGenerator)</code> performs the conversion for a rate <math>k/n</math> feedforward encoder. <code>ConstraintLength</code> is a 1-by-<math>k</math> vector that specifies the delay for the encoder's <math>k</math> input bit streams. <code>CodeGenerator</code> is a <math>k</math>-by-<math>n</math> matrix of octal numbers that specifies the <math>n</math> output connections for each of the encoder's <math>k</math> input bit streams.</p> <p><code>trellis = poly2trellis(ConstraintLength, CodeGenerator, ... FeedbackConnection)</code> is the same as the syntax above, except that it applies to a feedback, not feedforward, encoder. <code>FeedbackConnection</code> is a 1-by-<math>k</math> vector of octal numbers that specifies the feedback connections for the encoder's <math>k</math> input bit streams.</p> <p>For both syntaxes, the output is a MATLAB structure whose fields are as in the table below.</p>

**Table 3-17: Fields of the Output Structure `trellis` for a Rate  $k/n$  Code**

Field in <code>trellis</code> Structure	Dimensions	Meaning
<code>numInputSymbols</code>	Scalar	Number of input symbols to the encoder: $2^k$
<code>numOutputSymbols</code>	Scalar	Number of output symbols from the encoder: $2^n$
<code>numStates</code>	Scalar	Number of states in the encoder

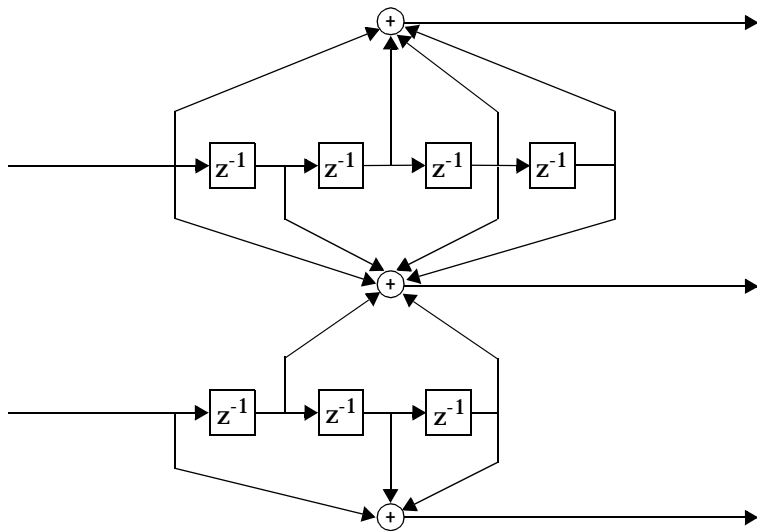
Table 3-17: Fields of the Output Structure trellis for a Rate k/n Code (Continued)

Field in trellis Structure	Dimensions	Meaning
nextStates	numStates-by- $2^k$ matrix	Next states for all combinations of current state and current input.
outputs	numStates-by- $2^k$ matrix	Outputs (in octal) for all combinations of current state and current input

For more about this structure, see the reference page for the `istrellis` function.

Examples

Consider the rate 2/3 feedforward convolutional encoder depicted in the figure below. The reference page for the `convenc` function includes an example that uses this encoder.



For this encoder, the `ConstraintLength` vector is [5,4] and the `CodeGenerator` matrix is [27,33,0; 0,5,13]. The output below reveals part of the corresponding trellis structure description of this encoder.

```
trellis = poly2trellis([5 4],[27 33 0; 0 5 13])
```

```
trellis =

    numInputSymbols: 4
    numOutputSymbols: 8
    numStates: 128
    nextStates: [128x4 double]
    outputs: [128x4 double]
```

The scalar field `trellis.numInputSymbols` has the value 4 because the combination of two input bit streams can produce four different input symbols. Similarly, `trellis.numOutputSymbols` is 8 because the three output bit streams can produce eight different output symbols.

The scalar field `trellis.numStates` is 128 (that is,  $2^7$ ) because each of the encoder's seven memory registers can have one of two binary values.

To get details about the matrix fields `trellis.nextStates` and `trellis.outputs`, inquire specifically about them. As an example, the command below displays the first five rows of the 128-by-4 matrix `trellis.nextStates`.

```
trellis.nextStates(1:5,:)
```

```
ans =

    0    64     8    72
    0    64     8    72
    1    65     9    73
    1    65     9    73
    2    66    10    74
```

This first row indicates that if the encoder starts in the zeroth state and receives input bits of 00, 01, 10, or 11, respectively, then the next state will be the 0th, 64th, 8th, or 72nd state, respectively. The 64th state means that the bottom-left memory register in the diagram contains the value 1, while the other six memory registers contain zeros.

## See Also

`istrellis`, `convenc`, `vitdec`

# qaskdeco

---

**Purpose** Demap a message from a QASK square signal constellation

**Syntax** `msg = qaskdeco(i nphase, quadr, M);`  
`msg = qaskdeco(i nphase, quadr, M, mi nmax);`

**Description** `msg = qaskdeco(i nphase, quadr, M)` demaps the message signal `msg` from the M-ary quadrature amplitude shift keying (QASK) square signal constellation points given in the vectors `i nphase` and `quadr`. Here `i nphase` lists the in-phase components of the points and `quadr` lists the corresponding quadrature components. `M` must be a power of 2. `qaskdeco` uses the default minimum/maximum value of the in-phase component and quadrature component. The defaults corresponding to small values of `M` are in the table on the reference page for the function `qaskenco`.

---

**Note** To see how symbols are mapped to the constellation points, generate a constellation plot using `qaskenco(M)`.

---

`msg = qaskdeco(i nphase, quadr, M, mi nmax)` is the same as the syntax above, except that `mi nmax` specifies the minimum and maximum in-phase and quadrature component values. `mi nmax` is a 2-by-2 matrix of the form shown below.

$$\text{minmax} = \begin{bmatrix} \text{in-phase minimum} & \text{in-phase maximum} \\ \text{quadrature minimum} & \text{quadrature maximum} \end{bmatrix}$$

**Examples** The commands below show that `qaskdeco` and `qaskenco` are inverse operations.

```
msg = [0 3 5 3 2 5]'; M = 8;  
[i nphase, quadr] = qaskenco(msg, M); % Map the message.  
newmsg = qaskdeco(i nphase, quadr, M) % Demap to recover data.  
  
newmsg =  
  
0  
3
```



5  
3  
2  
5

**See Also**      qaskenco, decode, demodmap

**Purpose** Map a message to a QASK square signal constellation

**Syntax**

```
qaskenco(M)
qaskenco(msg, M)
[inphase, quadr] = qaskenco(M)
[inphase, quadr] = qaskenco(msg, M)
```

**Description** qaskenco(M) plots the square signal constellation for M-ary quadrature amplitude shift keying (QASK) modulation, labeling the Mpoints with numbers in the range [0, M-1]. M must be a power of 2. If M is a perfect square, then qaskenco labels the constellation points so as to implement Gray code.

qaskenco(msg, M) is the same as the syntax above, except that only those points with labels in the vector msg are plotted. The elements in msg must be integers in the range [0, M-1].

[inphase, quadr] = qaskenco(M) returns vectors inphase and quadr that represent the coordinates of the points in the signal constellation for M-ary QASK modulation. inphase gives the in-phase component of each point and quadr gives the quadrature component of each point. M must be a power of 2.

[inphase, quadr] = qaskenco(msg, M) is the same as the syntax above, except that inphase and quadr represent only those constellation points with labels in the vector msg. (These labels are the same number labels that appear in the plot that the command qaskenco(msg, M) produces.) The elements in msg must be integers in the range [0, M-1].

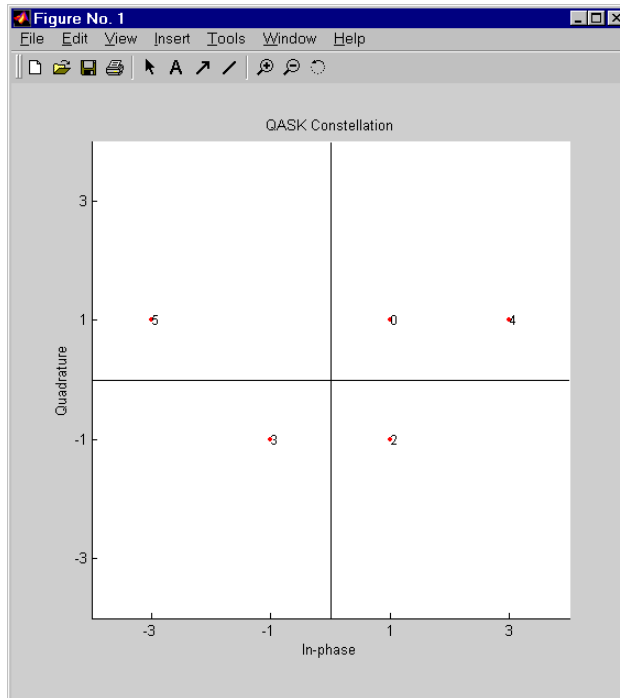
The table below shows the maximum value of inphase and quadr, for several small values of M

M	Maximum of inphase and quadr	M	Maximum of inphase and quadr
2	1	32	5
4	1	64	7
8	3 (maximum of quadr is 1)	128	11
16	3	256	15

## Examples

The command below displays that part of the 8-ary QASK square constellation that corresponds to the points in the digital message signal [0 3 4 3 2 5].

```
qaskenco([0 3 4 3 2 5], 8)
```



The commands below capture the same information in vectors `inphase` and `quadr` instead of in a plot.

```
[inphase, quadr] = qaskenco([0 3 5 3 2 5], 8);  
inphase'
```

```
ans =
```

```
1    -1    -3    -1    1    -3
```

```
quadr'
```

## qaskenco

---

```
ans =
```

```
      1      -1      1      -1      -1      1
```

The command below captures in `inphase` and `quadr` the coordinates of all eight points in the 8-ary QASK square constellation.

```
[ inphase2, quad2] = qaskenco(8);
```

### See Also

`encode`, `modmap`, `qaskdeco`

<b>Purpose</b>	Produce a quantization index and a quantized output value
<b>Syntax</b>	<pre> index = quantiz(sig, partition); [index, quants] = quantiz(sig, partition, codebook); [index, quants, distort] = quantiz(sig, partition, codebook); </pre>
<b>Description</b>	<p><code>index = quantiz(sig, partition)</code> returns the quantization levels in the real vector <code>signal sig</code> using the parameter <code>partition</code>. <code>partition</code> is a real vector whose entries are in strictly ascending order. If <code>partition</code> has length <math>n</math>, then <code>index</code> is a column vector whose <math>k</math>th entry is:</p> <ul style="list-style-type: none"> <li>• 0 if <math>\text{sig}(k) \leq \text{partition}(1)</math></li> <li>• <math>m</math> if <math>\text{partition}(m) &lt; \text{sig}(k) \leq \text{partition}(m+1)</math></li> <li>• <math>n</math> if <math>\text{partition}(n) &lt; \text{sig}(k)</math></li> </ul> <p><code>[index, quants] = quantiz(sig, partition, codebook)</code> is the same as the syntax above, except that <code>codebook</code> prescribes a value for each partition in the quantization and <code>quants</code> contains the quantization of <code>sig</code> based on the quantization levels and prescribed values. <code>codebook</code> is a vector whose length exceeds the length of <code>partition</code> by one. <code>quants</code> is a row vector whose length is the same as the length of <code>sig</code>. <code>quants</code> is related to <code>codebook</code> and <code>index</code> by</p> <pre>quants(ii) = codebook(index(ii)+1);</pre> <p>where <code>ii</code> is an integer between 1 and <code>length(sig)</code>.</p> <p><code>[index, quants, distort] = quantiz(sig, partition, codebook)</code> is the same as the syntax above, except that <code>distort</code> estimates the mean square distortion of this quantization data set.</p>
<b>Examples</b>	<p>The command below rounds several numbers between 1 and 100 up to the nearest multiple of ten. <code>quants</code> contains the rounded numbers, and <code>index</code> tells which quantization level each number is in.</p> <pre> [index, quants] = quantiz([3 34 84 40 23], 10:10:90, 10:10:100)  index =      0     3 </pre>

# quantiz

---

8  
3  
2

quants =

10      40      90      40      30

**See Also**

lloyd, dpcmenco, dpcmdeco

<b>Purpose</b>	Generate bit error patterns
<b>Syntax</b>	<pre>out = randerr(m); out = randerr(m, n); out = randerr(m, n, errors); out = randerr(m, n, errors, state);</pre>
<b>Description</b>	<p>For all syntaxes, <code>randerr</code> treats each row of <code>out</code> independently.</p> <p><code>out = randerr(m)</code> generates an <math>m</math>-by-<math>m</math> binary matrix, each row of which has exactly one nonzero entry in a random position. Each allowable configuration has an equal probability.</p> <p><code>out = randerr(m, n)</code> generates an <math>m</math>-by-<math>n</math> binary matrix, each row of which has exactly one nonzero entry in a random position. Each allowable configuration has an equal probability.</p> <p><code>out = randerr(m, n, errors)</code> generates an <math>m</math>-by-<math>n</math> binary matrix, where <code>errors</code> determines how many nonzero entries are in each row:</p> <ul style="list-style-type: none"><li>• If <code>errors</code> is a scalar, then it is the number of nonzero entries in each row.</li><li>• If <code>errors</code> is a row vector, then it lists the possible number of nonzero entries in each row.</li><li>• If <code>errors</code> is a matrix having two rows, then the first row lists the possible number of nonzero entries in each row and the second row lists the probabilities that correspond to the possible error counts.</li></ul> <p>Once <code>randerr</code> determines the <i>number</i> of nonzero entries in a given row, each configuration of that number of nonzero entries has equal probability.</p> <p><code>out = randerr(m, n, prob, state)</code> is the same as the syntax above, except that it first resets the state of MATLAB's uniform random number generator <code>rand</code> to the integer <code>state</code>.</p>
<b>Examples</b>	<p>To generate an 8-by-7 binary matrix, each row of which is equally likely to have either zero or two nonzero entries, use the command below.</p> <pre>out = randerr(8, 7, [0 2])</pre>

## randerr

---

```
out =
```

0	0	0	0	0	0	0
0	0	0	0	0	0	0
0	0	1	0	0	0	1
1	0	1	0	0	0	0
0	0	0	0	0	0	0
0	0	0	0	0	0	0
0	0	0	0	1	1	0
1	0	1	0	0	0	0

To alter the scenario above by making it three times as likely that a row has two nonzero entries, use the command below instead. Notice that the second row of the error parameter sums to one.

```
out2 = randerr(8, 7, [0 2; .25 .75])
```

```
out =
```

0	0	0	0	0	0	0
1	0	0	0	0	0	1
1	0	0	0	0	0	1
0	0	0	1	0	1	0
0	0	0	0	0	0	0
0	1	0	0	0	0	1
0	0	0	0	0	0	0
1	0	0	0	1	0	0

### See Also

rand, randint, randsrc



<b>Purpose</b>	Generate matrix of uniformly distributed random integers
<b>Syntax</b>	<pre>out = randint out = randint(m); out = randint(m, n); out = randint(m, n, range); out = randint(m, n, range, state);</pre>
<b>Description</b>	<p><code>out = randint</code> generates a random scalar that is either zero or one, with equal probability.</p> <p><code>out = randint(m)</code> generates an <math>m</math>-by-<math>m</math> binary matrix, each of whose entries independently takes the value zero with probability <math>1/2</math>.</p> <p><code>out = randint(m, n)</code> generates an <math>m</math>-by-<math>n</math> binary matrix, each of whose entries independently takes the value zero with probability <math>1/2</math>.</p> <p><code>out = randint(m, n, range)</code> generates an <math>m</math>-by-<math>n</math> integer matrix. If <code>range</code> is zero, then <code>out</code> is a zero matrix. Otherwise, the entries are uniformly distributed and independently chosen from the range:</p> <ul style="list-style-type: none"><li>• <code>[0, range-1]</code> if <code>range</code> is a positive integer</li><li>• <code>[range+1, 0]</code> if <code>range</code> is a negative integer</li><li>• Between <code>min</code> and <code>max</code>, inclusive, if <code>range = [min, max]</code> or <code>[max, min]</code></li></ul> <p><code>out = randint(m, n, range, state)</code> is the same as the syntax above, except that it first resets the state of MATLAB's uniform random number generator <code>rand</code> to the integer state.</p>
<b>Examples</b>	<p>To generate a 10-by-10 matrix whose elements are uniformly distributed in the range from 0 to 7, you can use either of the following commands.</p> <pre>out = randint(10, 10, [0, 7]); out = randint(10, 10, 8);</pre>
<b>See Also</b>	<code>rand</code> , <code>randsrc</code> , <code>randerr</code>

# randsrc

---

**Purpose** Generate random matrix using prescribed alphabet

**Syntax**

```
out = randsrc;  
out = randsrc(m);  
out = randsrc(m, n);  
out = randsrc(m, n, alphabet);  
out = randsrc(m, n, [alphabet; prob]);  
out = randsrc(m, n, ..., state);
```

**Description**

`out = randsrc` generates a random scalar that is either -1 or 1, with equal probability.

`out = randsrc(m)` generates an m-by-m matrix, each of whose entries independently takes the value -1 with probability 1/2, and 1 with probability 1/2.

`out = randsrc(m, n)` generates an m-by-n matrix, each of whose entries independently takes the value -1 with probability 1/2, and 1 with probability 1/2.

`out = randsrc(m, n, alphabet)` generates an m-by-n matrix, each of whose entries is independently chosen from the entries in the row vector `alphabet`. Each entry in `alphabet` occurs in `out` with equal probability. Duplicate values in `alphabet` are ignored.

`out = randsrc(m, n, [alphabet; prob])` generates an m-by-n matrix, each of whose entries is independently chosen from the entries in the row vector `alphabet`. Duplicate values in `alphabet` are ignored. The row vector `prob` lists corresponding probabilities, so that the symbol `alphabet(k)` occurs with probability `prob(k)`, where `k` is any integer between one and the number of columns of `alphabet`. The elements of `prob` must add up to one.

`out = randsrc(m, n, ..., state);` is the same as the two preceding syntaxes, except that it first resets the state of MATLAB's uniform random number generator `rand` to the integer state.

**Examples**

To generate a 10-by-10 matrix whose elements are uniformly distributed among members of the set {-3,-1,1,3}, you can use either of these commands.

```
out = randsrc(10, 10, [-3 -1 1 3]);  
out = randsrc(10, 10, [-3 -1 1 3; .25 .25 .25 .25]);
```

To skew the probability distribution so that -1 and 1 each occur with probability .3, while -3 and 3 each occur with probability .2, use this command.

```
out = randsrc(10, 10, [-3 -1 1 3; .2 .3 .3 .2]);
```

**See Also**

rand, randi nt, randerr

**Purpose** Design a raised cosine FIR filter

**Syntax**

```
b = rcosfir(R, n_T, rate, T);
b = rcosfir(R, n_T, rate, T, filter_type);
rcosfir(...);
rcosfir(..., color);
[b, sample_time] = rcosfir(...);
```

Optional Inputs	Input	Default Value
	n_T	[-3, 3]
	rate	5
	T	1

**Description** The time response of the raised cosine filter has the form

$$h(t) = \frac{\sin(\frac{\pi t}{T}) \cos(\frac{\pi R t}{T})}{(\frac{\pi t}{T}) \left(1 - \frac{4R^2 t^2}{T^2}\right)}$$

`b = rcosfir(R, n_T, rate, T)` designs a raised cosine filter and returns a vector `b` of length  $(n\_T(2) - n\_T(1)) * rate + 1$ . The filter's rolloff factor is `R`, where  $0 \leq R \leq 1$ . `T` is the duration of each bit in seconds. `n_T` is a length-two vector that indicates the number of symbol periods before and after the peak response. `rate` is the number of points in each input symbol period of length `T`. The input sample rate is `T` samples per second, while the output sample rate is `T*rate` samples per second.

The order of the FIR filter is

$$(n\_T(2) - n\_T(1)) * rate$$

The arguments `n_T`, `rate`, and `T` are optional inputs whose default values are [-3,3], 5, and 1, respectively.

`b = rcosfir(R, n_T, rate, T, filter_type)` designs a square-root raised cosine filter if `filter_type` is `'sqrt'`. If `filter_type` is `'normal'` then this syntax is the same as the previous one.

The impulse response of a square root raised cosine filter is

$$h(t) = 4r \frac{\cos\left((1+r)\pi\frac{t}{T}\right) + \frac{\sin\left((1-r)\pi\frac{t}{T}\right)}{4r\frac{t}{T}}}{\pi\sqrt{T}\left(\left(4r\frac{t}{T}\right)^2 - 1\right)}$$

`rcosfir(...)` produces plots of the time and frequency responses of the raised cosine filter.

`rcosfir(..., color)` uses the string `color` to determine the plotting color. The choices for `color` are the same as those listed for the `plot` function.

`[b, sample_time] = rcosfir(...)` returns the FIR filter and its sample time.

## Examples

The commands below compare different rolloff factors.

```
rcosfir(0);
subplot(211); hold on;
subplot(212); hold on;
rcosfir(.5, [], [], [], [], 'r-');
rcosfir(1, [], [], [], [], 'g-');
```

## See Also

`rcosflt`, `rcosfir`, `firrcos` (Signal Processing Toolbox)

## References

Korn, Israel. *Digital Communications*. New York: Van Nostrand Reinhold, 1985.

**Purpose** Filter the input signal using a raised cosine filter

**Syntax**

```
y = rcosflt(x, Fd, Fs);  
y = rcosflt(x, Fd, Fs, 'filter_type', r, delay, tol);  
y = rcosflt(x, Fd, Fs, 'filter_type/Fs', r, delay, tol);  
y = rcosflt(x, Fd, Fs, 'filter_type/filter', num, den);  
y = rcosflt(x, Fd, Fs, 'filter_type/filter', num, den, delay);  
y = rcosflt(x, Fd, Fs, 'filter_type/filter/Fs', num, den, ...);  
[y, t] = rcosflt(...);
```

Optional Inputs	Input	Default Value
	<i>filter_type</i>	<b>fir/normal</b>
	<i>r</i>	0.5
	<i>delay</i>	3
	<i>tol</i>	0.01
	<i>den</i>	1

**Description** The function `rcosflt` passes an input signal through a raised cosine filter. You can either let `rcosflt` design a raised cosine filter automatically or you can specify the raised cosine filter yourself using input arguments.

**Designing the Filter Automatically**

`y = rcosflt(x, Fd, Fs)` designs a raised cosine FIR filter and then filters the input signal `x` using it. The sample frequency for the digital input signal `x` is `Fd`, and the sample frequency for the output signal `y` is `Fs`. The ratio `Fs/Fd` must be an integer. In the course of filtering, `rcosflt` upsamples the data by a factor of `Fs/Fd`, by inserting zeros between samples. The order of the filter is `1+2*delay*Fs/Fd`, where `delay` is 3 by default. If `x` is a vector, then the sizes of `x` and `y` are related by this equation.

$$\text{length}(y) = (\text{length}(x) + 2 * \text{delay}) * Fs/Fd$$

Otherwise, `y` is a matrix, each of whose columns is the result of filtering the corresponding column of `x`.

`y = rcosflt(x, Fd, Fs, 'filter_type', r, delay, tol)` designs a raised cosine FIR or IIR filter and then filters the input signal `x` using it. The ratio `Fs/Fd` must be an integer. `r` is the rolloff factor for the filter, a real number in the range `[0, 1]`. `delay` is the filter's group delay, measured in input samples. The actual group delay in the filter design is `delay/Fd` seconds. The input `tol` is the tolerance in the IIR filter design. FIR filter design does not use `tol`.

The characteristics of `x`, `Fd`, `Fs`, and `y` are as in the first syntax.

The fourth input argument, `'filter_type'`, is a string that determines the type of filter that `rcosflt` should design. Use one of the values in the table below.

Table 3-18: Values of `filter_type` to Determine the Type of Filter

Type of Filter	Value of opt
FIR raised cosine filter	<b>fir</b> or <b>fir/normal</b>
IIR raised cosine filter	<b>iir</b> or <b>iir/normal</b>
Square-root FIR raised cosine filter	<b>fir/sqrt</b>
Square-root IIR raised cosine filter	<b>iir/sqrt</b>

`y = rcosflt(x, Fd, Fs, 'filter_type/Fs', r, delay, tol)` is the same as the previous syntax, except that it assumes that `x` has sample frequency `Fs`. This syntax does not upsample `x` any further. If `x` is a vector, then the relative sizes of `x` and `y` are related by this equation.

$$\text{length}(y) = \text{length}(x) + (2 * \text{delay} * Fs/Fd)$$

As before, if `x` is a nonvector matrix, then `y` is a matrix each of whose columns is the result of filtering the corresponding column of `x`.

Specifying the Filter Using Input Arguments

`y = rcosflt(x, Fd, Fs, 'filter_type/filter', num, den)` filters the input signal `x` using a filter whose transfer function numerator and denominator are given in `num` and `den`, respectively. If `filter_type` includes **fir**, then omit `den`. This syntax uses the same arguments `x`, `Fd`, `Fs`, and `filter_type` as explained in the first and second syntaxes above.

`y = rcosflt(x, Fd, Fs, 'filter_type/filter', num, den, delay)` uses `delay` in the same way that the `rcosine` function uses it. This syntax assumes that the filter described by `num`, `den`, and `delay` was designed using `rcosine`.

As before, if `x` is a nonvector matrix, then `y` is a matrix each of whose columns is the result of filtering the corresponding column of `x`.

`y = rcosflt(x, Fd, Fs, 'filter_type/filter/Fs', num, den, ...)` is the same as the earlier syntaxes, except that it assumes that `x` has sample frequency `Fs` instead of `Fd`. This syntax does not upsample `x` any further. If `x` is a vector, then the relative sizes of `x` and `y` are related by this equation.

$$\text{length}(y) = \text{length}(x) + (2 * \text{delay} * Fs/Fd)$$

### Additional Output

`[y, t] = rcosflt(...)` outputs `t`, a vector that contains the sampling time points of `y`.

### See Also

`rcosfir`, `rcosfir`, `rcosine`, `rcosdemo`, `grpdelay` (Signal Processing Toolbox)

### References

Korn, Israel. *Digital Communications*. New York: Van Nostrand Reinhold, 1985.



**Purpose** Design a raised cosine IIR filter

**Syntax**

```
[num, den] = rcosiir(R, T_delay, rate, T, tol);
[num, den] = rcosiir(R, T_delay, rate, T, tol, filter_type);
rcosiir(...);
rcosiir(..., color);
[num, den, sample_time] = rcosiir(...);
```

Optional Inputs	Input	Default Value
	T_delay	3
	rate	5
	T	1
	tol	0.01

**Description** The time response of the raised cosine filter has the form

$$h(t) = \frac{\sin\left(\frac{\pi t}{T}\right) \cos\left(\frac{\pi R t}{T}\right)}{\frac{\pi t}{T} \sqrt{1 - \frac{4 R^2 t^2}{T^2}}}$$

[num, den] = rcosiir(R, T\_delay, rate, T, tol) designs an IIR approximation of an FIR raised cosine filter, and returns the numerator and denominator of the IIR filter. The filter's rolloff factor is R, where  $0 \leq R \leq 1$ . T is the symbol period in seconds. The filter's group delay is T\_delay symbol periods. rate is the number of sample points in each interval of duration T. The input sample rate is T samples per second, while the output sample rate is T\*rate samples per second. If tol is an integer greater than one, then it becomes the order of the IIR filter; if tol is less than 1, then it indicates the relative tolerance for rcosiir to use when selecting the order based on the singular values.

The arguments T\_delay, rate, T, and tol are optional inputs whose default values are 3, 5, 1, and 0.01, respectively.

`[num,den] = rcosiiir(R,T_delay,rate,T,tol,filter_type)` designs a square-root raised cosine filter if *filter\_type* is '**sqrt**'. If *filter\_type* is '**normal**' then this syntax is the same as the previous one.

`rcosiiir(...)` plots the time and frequency responses of the raised cosine filter.

`rcosiiir(...,color)` uses the string *color* to determine the plotting color. The choices for *color* are the same as those listed for the `plot` function.

`[num,den,sample_time] = rcosiiir(...)` returns the transfer function and the sample time of the IIR filter.

### Examples

The script below compares different values of *T\_delay*.

```
rcosiiir(0,10);  
subplot(211); hold on;  
subplot(212); hold on;  
col = ['r-'; 'g-'; 'b-'; 'm-'; 'c-'; 'w-'];  
R = [8, 6, 4, 3, 2, 1];  
for i = R  
    rcosiiir(0,i,[],[],[],[],col(find(R==i),:));  
end;
```

This example shows how the filter's frequency response more closely approximates that of the ideal raised cosine filter as *T\_delay* increases.

### See Also

`rcosfir`, `rcosflt`, `rcosine`, `grpdelay` (Signal Processing Toolbox)

### References

Kailath, Thomas. *Linear Systems*. Englewood Cliffs, N.J.: Prentice-Hall, 1980.  
Korn, Israel. *Digital Communications*. New York: Van Nostrand Reinhold, 1985.

**Purpose** Design a raised cosine filter

**Syntax**

```
num = rcosine(Fd, Fs);
[num, den] = rcosine(Fd, Fs, type_flag);
[num, den] = rcosine(Fd, Fs, type_flag, r);
[num, den] = rcosine(Fd, Fs, type_flag, r, delay);
[num, den] = rcosine(Fd, Fs, type_flag, r, delay, tol);
```

**Description** `num = rcosine(Fd, Fs)` designs a finite impulse response (FIR) raised cosine filter and returns its transfer function. The input digital signal has frequency `Fd`. The sampling frequency for the filter is `Fs`. The ratio `Fs/Fd` must be a positive integer. The default rolloff factor is `.5`. The filter's group delay, which is the time between the input to the filter and the filter's peak response, is three input samples. Equivalently, the group delay is  $3/Fd$  seconds.

`[num, den] = rcosine(Fd, Fs, type_flag)` designs a raised cosine filter using directions in the string variable `type_flag`. Filter types are listed in the table below, along with the corresponding values of `type_flag`.

**Table 3-19: Types of Filter and Corresponding Values of type\_flag**

Type of Filter	Value of type_flag
Finite impulse response (FIR)	'default' or 'fir/normal'
Infinite impulse response (IIR)	'iir' or 'iir/normal'
Square-root raised cosine FIR	'sqrt' or 'fir/sqrt'
Square-root raised cosine IIR	'iir/sqrt'

The default tolerance value in IIR filter design is 0.01.

`[num, den] = rcosine(Fd, Fs, type_flag, r)` specifies the rolloff factor, `r`. The rolloff factor is a real number in the range `[0, 1]`.

`[num, den] = rcosine(Fd, Fs, type_flag, r, delay)` specifies the filter's group delay, measured in input samples. `delay` is a positive integer. The actual group delay in the filter design is `delay/Fd` seconds.

## rcosine

---

[num, den] = rcosine(Fd, Fs, *type\_flag*, r, delay, tol) specifies the tolerance in the IIR filter design. FIR filter design does not use tol.

**See Also** rcosfir, rcosiir, rcosflt, rcosdemo, grpdelay (Signal Processing Toolbox)

**References** Korn, Israel. *Digital Communications*. New York: Van Nostrand Reinhold, 1985.

Purpose

Reed-Solomon decoder

Syntax

```
msg = rsdeco(code, n, k);
msg = rsdeco(code, n, k, format);
msg = rsdeco(code, field, ...);
[msg, err] = rsdeco(...);
[msg, err, ccode] = rsdeco(...);
[msg, err, ccode, cerr] = rsdeco(...);
```

Description

For All Syntaxes

The encoding counterpart for this function is rsenco.

In all cases, the codeword length  $n$  must have the form  $2^m-1$  where  $m$  is an integer greater than or equal to 3.

The matrix code, which contains the code words to be decoded, can have one of several formats. The table below shows the formats for `msg`, how the optional argument `format` should reflect the format of `msg`, and how the format of the output code depends on these choices. If `format` is not specified as input, then its default value is **binary**.

Table 3-20: Information Formats for Reed-Solomon Decoding

Format of code	Value of format Argument	Format of msg
Binary matrix with $m$ columns	'binary'	Binary matrix with $m$ columns
Example: code = [0 0 0; 0 1 1; 0 1 1; 1 1 0; 1 0 1; 1 0 0; 0 1 1]		
Binary column vector	'binary'	Binary column vector
Example: code = [0 0 0, 0 1 1, 0 1 1, 1 1 0, 1 0 1, 1 0 0, 0 1 1]';		

Table 3-20: Information Formats for Reed-Solomon Decoding (Continued)

Format of code	Value of format Argument	Format of msg
Matrix of integers in the range $[0, 2^m-1]$ , with n columns	' decimal '	Matrix of integers in the range $[0, 2^m-1]$ , with k columns
Example: code = [ 0, 6, 6, 3, 5, 1, 6]		
Matrix of integers in the range $[-1, 2^m-2]$ , with n columns	' power'	Matrix of integers in the range $[-1, 2^m-2]$ , with k columns
Example: code = [- 1, 5, 5, 2, 4, 0, 5]		

For Specific Syntaxes

`msg = rsdeco(code, n, k)` decodes `code` using the Reed-Solomon decoding method. `n` is the codeword length and `k` is the message length. `code` has either of the two binary formats described in Table 3-20, Information Formats for Reed-Solomon Decoding.

`msg = rsdeco(code, n, k, format)` is the same as the syntax above, except that `format` specifies the format of `code`. Table 3-20, Information Formats for Reed-Solomon Decoding, lists the possible values for `format`, as well as the corresponding shape and contents of `code`.

`msg = rsdeco(code, field, ...)` is a faster variation of the syntaxes above. `field` is a matrix that lists all elements of  $GF(2^m)$  in the format described in “List of All Elements of a Galois Field” on page 2-91. The size of `field` determines `n`.

`[msg, err] = rsdeco(...)` outputs the number `err`, which specifies the number of errors that occurred in the decoding.

`[msg, err, ccode] = rsdeco(...)` outputs `ccode`, a corrected version of `code`. The format of `ccode` matches the format of `code` in the input.

`[msg, err, ccode, cerr] = rsdeco(...)` outputs the number `cerr`, which specifies the number of errors found in the `ccode` column.

**Examples**

This example creates and decodes a noisy code. Although some codewords contain errors, the decoded message contains no errors.

```
L = 1000; % Number of bits in the computation
m = 4;
n = 2^m - 1; % Codeword length
k = n - 4; % Message word length
rand('state', 9876); % Initialize random number generator.
msg = randint(L, 1); % L bits of data
field = gftuple([-1 : n-1]', m); % List of elements in GF(2^m)
[code, added] = rsenco(msg, field, k); % Encode the data.
msg = [msg; zeros(added, 1)]; % Pad msg for later comparison.

% Add burst errors of length m to the code.
noi = rand(length(code)/m, 1) < .03; % Three percent noise
noi = (noi*ones(1, m))'; noi = noi(:);
code_noi = rem(code + noi, 2);

% Decode the noisy code.
[dec, err, ccode, err_c] = rsdeco(code_noi, field, k);
err_c = reshape(err_c, n, length(err_c)/n)';
% Number of code symbols that contain at least one error
num_err_codesyms = sum(err_c(:, 1) > 0)
% Number of bit errors after decoding
num_err_decbits = sum(abs(dec - msg))

num_err_codesyms =

    36

num_err_decbits =

    0
```

**See Also**

rsenco, rsdecode, rspoly, encode, decode

# rsdecode

---

**Purpose** Reed-Solomon decoding using the exponential format

**Syntax**

```
msg = rsdecode(code, k);  
msg = rsdecode(code, k, m);  
msg = rsdecode(code, k, fi el d);  
[msg, err] = rsdecode(...);  
[msg, err, ccode] = rsdecode(...);
```

**Description** **For All Syntaxes**

The encoding counterpart for this function is `rsencode`.

`rsdecode` uses the exponential format to represent elements of  $\text{GF}(2^m)$ . For example, an entry of 2 represents the element  $\alpha^2$ , where  $\alpha$  is a primitive element of  $\text{GF}(2^m)$ . If `fi el d` is not used as an input argument, then the exponential format is relative to a root of MATLAB's default primitive polynomial for  $\text{GF}(2^m)$ . If `fi el d` is used as an input argument, then its format and the formats in `msg` and `code` are all relative to the same primitive element of  $\text{GF}(2^m)$ . See "Representing Elements of Galois Fields" on page 2-90 for more information about these formats.

Since  $\text{GF}(2^m)$  has  $2^m$  elements, each codeword represents  $2^m(2^m-1)$  bits of information. Each decoded message represents  $2^m*k$  bits of information.

**For Specific Syntaxes**

`msg = rsdecode(code, k)` decodes `code` using the Reed-Solomon method. `k` is the message length. The codeword length  $n$  must have the form  $2^m-1$  for some integer  $m$  greater than or equal to 3. `code` is a matrix with  $n$  columns. Each row of `code` represents one codeword. Each entry of `code` represents an element of  $\text{GF}(2^m)$  in exponential format. `msg` is a matrix with `k` columns. Each row of `msg` represents one message. Each entry of `msg` is the exponential format of an element of  $\text{GF}(2^m)$ .

`msg = rsdecode(code, k, m)` is the same as the first syntax when the matrix `code` has  $2^m-1$  columns. This syntax is faster than the first.

`msg = rsdecode(code, k, fi el d)` is the same as the first syntax, except that `fi el d` is a matrix that lists the elements of  $\text{GF}(2^m)$  in the format described in



“List of All Elements of a Galois Field” on page 2-91. This syntax is faster than the first two.

`[msg, err] = rsdecode(...)` returns a column vector `err` that gives information about error correction. A nonnegative integer in `err(r)` indicates the number of errors corrected in the  $r$ th codeword; a negative integer indicates that there are more errors in the  $r$ th codeword than can be corrected.

`[msg, err, ccode] = rsdecode(...)` returns the corrected code in `ccode`.

## Examples

The script below continues the example from the reference page for `rsencode`. After corrupting some symbols from the code, it tries to recover the message.

```
m = 3; n = 2^m-1; % Codeword length is 7.
field = gftuple([-1:2^m-2]', m, 2); % List of elements in GF(2^m)
msg = [5 0 1; 2 3 4];
k = size(msg, 2); % Message length = number of columns of msg
genpoly = rspoly(n, k, field); % Generator polynomial
code = rsencode(msg, genpoly, n, field);
% Change up to three of the code symbols.
noisycode = code;
noisycode(1, 2) = randint(1, 1, [-1, n-1]);
noisycode(2, 1) = randint(1, 1, [-1, n-1]);
noisycode(2, 5) = randint(1, 1, [-1, n-1]);
% Try to decode.
[newmsg, err, ccode] = rsdecode(noisycode, k, field);
if ccode==code
    disp('All errors were corrected.')
end
if newmsg==msg
    disp('The message was recovered perfectly.')
end
```

Unless one of the random integers was zero, `err` is the matrix `[1; 2]`, which reflects the fact that we put one error in the first row of `noisycode` and two errors in the second row. Since this code’s error-correction capability is  $\text{floor}((n-k)/2)$ , or 2, all errors are corrected in this example.

## See Also

`rsencode`, `encode`, `decode`, `rsdeco`

# rsdecof

---

<b>Purpose</b>	Decode an ASCII file that was encoded using Reed-Solomon code
<b>Syntax</b>	<pre>rsdecof(file_in, file_out); rsdecof(file_in, file_out, err_cor);</pre>
<b>Description</b>	<p>This function is the inverse process of the function <code>rsencof</code> in that it decodes a file that <code>rsencof</code> encoded.</p> <p><code>rsdecof(file_in, file_out)</code> decodes the ASCII file <code>file_in</code> that was previously created by the function <code>rsencof</code> using an error-correction capability of 5. The decoded message is written to <code>file_out</code>. Both <code>file_in</code> and <code>file_out</code> are string variables.</p> <hr/> <p><b>Note</b> If the number of characters in <code>file_in</code> is not an integer multiple of 127, then the function appends <code>char(4)</code> symbols to the data it must decode. If you encode and then decode a file using <code>rsencof</code> and <code>rsdecof</code>, respectively, then the decoded file might have <code>char(4)</code> symbols at the end that the original file does not have.</p> <hr/> <p><code>rsdecof(file_in, file_out, err_cor)</code> is the same as the first syntax, except that <code>err_cor</code> specifies the error-correction capability for each block of 127 codeword characters. The message length is <math>127 - 2 * \text{err\_cor}</math>. The value in <code>err_cor</code> must match the value used in <code>rsencof</code> when <code>file_in</code> was created.</p>
<b>Examples</b>	An example is on the reference page for <code>rsencof</code> .
<b>See Also</b>	<code>rsencof</code> , <code>encode</code> , <code>decode</code> , <code>rsenco</code> , <code>rsdeco</code>

**Purpose** Reed-Solomon encoder

**Syntax**

```
code = rsenco(msg, n, k);  
code = rsenco(msg, n, k, format);  
code = rsenco(msg, n, k, format, genpoly);  
code = rsenco(msg, field, ...);  
[code, added] = rsenco(...);
```

**Description** For All Syntaxes

The decoding counterpart for this function is rsdeco.

In all cases, the codeword length  $n$  must have the form  $2^m - 1$  where  $m$  is an integer greater than or equal to 3.

The matrix `msg`, which contains the messages to be encoded, can have one of several formats. Table 3-21, Information Formats for Reed-Solomon Encoding, shows which formats are allowed for `msg`, how the optional argument `format` should reflect the format of `msg`, and how the format of the output code depends on these choices. If `format` is not specified as input, then its default value is **'binary'**.

**Table 3-21: Information Formats for Reed-Solomon Encoding**

Format of msg	Value of format Argument	Format of code
Binary matrix with $m$ columns	<b>'binary'</b>	Binary matrix with $m$ columns
Example: <code>msg = [1 1 0; 1 0 1; 1 0 0; 0 1 1; 1 1 0; 1 0 1; 1 0 0; 0 1 1]</code>		
Binary column vector	<b>'binary'</b>	Binary column vector
Example: <code>msg = [1 1 0, 1 0 1, 1 0 0, 0 1 1, 1 1 0, 1 0 1, 1 0 0, 0 1 1]'</code>		

Table 3-21: Information Formats for Reed-Solomon Encoding (Continued)

Format of msg	Value of format Argument	Format of code
Matrix of integers in the range $[0, 2^m-1]$ , with k columns	' decimal '	Matrix of integers in the range $[0, 2^m-1]$ , with n columns
Example: msg = [3, 5, 1, 6; 3, 5, 1, 6]		
Matrix of integers in the range $[-1, 2^m-2]$ , with k columns	' power '	Matrix of integers in the range $[-1, 2^m-2]$ , with n columns
Example: msg = [2, 4, 0, 5; 2, 4, 0, 5]		

For Specific Syntaxes

`code = rsenco(msg, n, k)` encodes `msg` using the Reed-Solomon encoding method. `k` is the message length. `msg` has either of the two binary formats described in Table 3-21, Information Formats for Reed-Solomon Encoding. The generator polynomial for the code is the output of the function `rspoly`.

`code = rsenco(msg, n, k, format)` is the same as the syntax above, except that `format` specifies the format of `msg`. Table 3-21, Information Formats for Reed-Solomon Encoding, lists the possible values for `format`, as well as the corresponding shape and contents of `msg`.

`code = rsenco(msg, n, k, format, genpoly)` is the same as the syntax above, except that `genpoly` is a row vector that gives the coefficients, in order of ascending powers, of the generator polynomial for the code. Each coefficient is an element of  $GF(2^m)$  expressed in exponential format. For a description of exponential format, see “Exponential Format” on page 2-90.

`code = rsenco(msg, field, ...)` is a faster variation of the syntaxes above. `field` is a matrix that lists all elements of  $GF(2^m)$  in the format described in “List of All Elements of a Galois Field” on page 2-91. The size of `field` determines `n`.

[code, added] = rsenco(. . .) returns the additional variable added. added is the number of zeros that were placed at the end of the message matrix before encoding, in order for the matrix to have the appropriate shape.

**Algorithm**

rsenco invokes the function rsencode, which processes data in power format. If msg has decimal or binary format, then rsenco converts it to the power format, passes it to rsencode, and converts the code back to the original format of msg. Binary data has the longest processing time.

**See Also**

rsdeco, encode, decode, rsencode

**Purpose** Reed-Solomon encoding using the exponential format

**Syntax**

```
code = rsencode(msg, genpoly, n);  
code = rsencode(msg, genpoly, n, m);  
code = rsencode(msg, genpoly, n, field);
```

**Description** For All Syntaxes

The decoding counterpart for this function is `rsdecode`.

`rsencode` uses the exponential format to represent elements of  $\text{GF}(2^m)$ . For example, an entry of 2 represents the element  $\alpha^2$ , where  $\alpha$  is a primitive element of  $\text{GF}(2^m)$ . If `field` is not used as an input argument, then the exponential format is relative to a root of MATLAB's default primitive polynomial for  $\text{GF}(2^m)$ . If `field` is used as an input argument, then its format and the formats in `msg` and `code` are all relative to the same primitive element of  $\text{GF}(2^m)$ . See "Representing Elements of Galois Fields" on page 2-90 for more information about these formats.

Since  $\text{GF}(2^m)$  has  $2^m$  elements, each codeword represents  $2^m(2^m-1)$  bits of information. Each decoded message represents  $2^{m*k}$  bits of information.

**For Specific Syntaxes**

`code = rsencode(msg, genpoly, n)` encodes the message `msg` using the Reed-Solomon coding method. `n`, the codeword length, must have the form  $2^m-1$  for some integer  $m$  greater than or equal to 3. If the message length is  $k$ , then `msg` is a matrix having  $k$  columns. Each entry of `msg` represents an element of  $\text{GF}(2^m)$  in exponential format. Each row of `msg` is treated as a separate message. Each row of `code` represents a codeword, and each entry is the exponential format of an element of  $\text{GF}(2^m)$ . The last  $k$  columns of `code` are just `msg`; that is, the parity bits are at the beginning of each codeword. `genpoly` is a row vector that gives the coefficients, in order of ascending powers, of the generator polynomial. Each coefficient is specified in exponential format.

`code = rsencode(msg, genpoly, n, m)` is the same as `code = rsencode(msg, genpoly, 2^m-1)` when `m` is an integer greater than or equal to 3. Specifying `m` as a fourth input argument speeds the execution.

`code = rsencode(msg, genpoly, n, field)` is the same as the first syntax, except that `field` is a matrix that lists the elements of  $\text{GF}(2^m)$  in the format described in “List of All Elements of a Galois Field” on page 2-91. This syntax is faster than the first one.

## Examples

The commands below use the third syntax of `rsencode` to encode two messages.

```
m = 3; n = 2^m-1; % Codeword length is 7.
field = gftuple([-1:2^m-2]', m, 2); % List of elements in GF(2^m)
msg = [5 0 1; 2 3 4];
k = size(msg, 2); % Message length = number of columns of msg
genpoly = rspoly(n, k, field); % Generator polynomial
code = rsencode(msg, genpoly, n, field);
```

The reference page for `rsdecode` continues this example by corrupting the code and then decoding it.

## See Also

`rsdecode`, `encode`, `decode`, `rspoly`, `rsdeco`

# rsencof

---

**Purpose** Encode an ASCII file using Reed-Solomon code

**Syntax** `rsencof(file_in, file_out);`  
`rsencof(file_in, file_out, err_cor);`

**Description** `rsencof(file_in, file_out)` encodes the ASCII file `file_in` using (127, 117) Reed-Solomon code. The error-correction capability of this code is 5 for each block of 127 codeword characters. This function writes the encoded text to the file `file_out`. Both `file_in` and `file_out` are string variables.

`rsencof(file_in, file_out, err_cor)` is the same as the first syntax, except that `err_cor` specifies the error correction capability for each block of 127 codeword characters. The message length is  $127 - 2 * \text{err\_cor}$ .

---

**Note** If the number of characters in `file_in` is not an integer multiple of  $127 - 2 * \text{err\_cor}$ , then the function appends `char(4)` symbols to `file_out`.

---

**Examples** The file `matlabroot/toolbox/comm/comm/oct2dec.m` contains text help for the `oct2dec` function in this toolbox. The commands below encode the file using `rsencof` and then decode it using `rsdecof`.

```
file_in = [matlabroot ' /toolbox/comm/comm/oct2dec.m' ];  
file_out = 'encodedfile'; % Or use another filename  
rsencof(file_in, file_out) % Encode the file.
```

```
file_in = file_out;  
file_out = 'decodedfile'; % Or use another filename  
rsdecof(file_in, file_out) % Decode the file.
```

To see the original file and the decoded file in the MATLAB workspace, use the commands below (or similar ones if you modified the filenames above).

```
type oct2dec.m  
type decodedfile
```

**See Also** `rsdecof`, `encode`, `decode`, `rsenco`, `rsdeco`



**Purpose** Produce Reed-Solomon code generator polynomial

**Syntax**

```
genpoly = rspoly(n, k);
genpoly = rspoly(n, k, m);
genpoly = rspoly(n, k, field);
[genpoly, t] = rspoly(...);
```

**Description** `genpoly = rspoly(n, k)` finds the generator polynomial of a Reed-Solomon code with codeword length  $n$  and message length  $k$ . `genpoly` is a row vector that represents the coefficients of the generator polynomial in order of ascending powers. Each coefficient is an element of  $GF(2^m)$  represented in exponential format, as described in the section “Representing Elements of Galois Fields” on page 2-90.

`genpoly = rspoly(n, k, m)` is the same as `genpoly = rspoly(2m-1, k)`, but faster. If  $n$  does not equal  $2^m-1$ , then an error results.

`genpoly = rspoly(n, k, field)` is the same as the first syntax listed, except that `field` indirectly specifies the primitive element for  $GF(2^m)$  relative to which the coefficients in `genpoly` are expressed. `field` is a matrix that lists the elements of  $GF(2^m)$  in the format described in “List of All Elements of a Galois Field” on page 2-91. Both `field` and `genpoly` use exponential formats relative to the same primitive element. This syntax is faster than the first syntax listed.

`[genpoly, t] = rspoly(...)` returns in `t` the error-correction capability of the Reed-Solomon code.

**Examples** The command below shows that the (15, 11) Reed-Solomon code generator polynomial is  $\alpha^{10} + \alpha^3 X + \alpha^6 X^2 + \alpha^{13} X^3 + X^4$ .

```
genpoly = rspoly(15, 11, 4)

genpoly =

    10     3     6    13     0
```

The syntax below uses `field` as the third input argument in `rspoly` and obtains the same result.

# rspoly

---

```
m = 4;
field = gftuple([-1:2^m-2]', m, 2);
genpoly2 = rspoly(15, 11, field)

genpoly2 =

    10     3     6    13     0
```

## See Also

encode, decode, rsenco, rsdeco

**Purpose** Generate a scatter plot

**Syntax**

```
scatterplot(x);
scatterplot(x, n);
scatterplot(x, n, offset);
scatterplot(x, n, offset, plotstring);
scatterplot(x, n, offset, plotstring, h);
h = scatterplot(...);
```

**Description** `scatterplot(x)` produces a scatter plot for the signal `x`. The interpretation of `x` depends on its shape and complexity:

- If `x` is a real two-column matrix, then `scatterplot` interprets the first column as in-phase components and the second column as quadrature components.
- If `x` is a complex vector, then `scatterplot` interprets the real part as in-phase components and the imaginary part as quadrature components.
- If `x` is a real vector, then `scatterplot` interprets it as a real signal.

`scatterplot(x, n)` is the same as the first syntax, except that the function plots every `n`th value of the signal, starting from the first value. That is, the function decimates `x` by a factor of `n` before plotting.

`scatterplot(x, n, offset)` is the same as the first syntax, except that the function plots every `n`th value of the signal, starting from the `(offset+1)`st value in `x`.

`scatterplot(x, n, offset, plotstring)` is the same as the syntax above, except that `plotstring` determines the plotting symbol, line type, and color for the plot. `plotstring` is a string whose format and meaning are the same as in the `plot` function.

`scatterplot(x, n, offset, plotstring, h)` is the same as the syntax above, except that the scatter plot is in the figure whose handle is `h`, rather than a new figure. `h` must be a handle to a figure that `scatterplot` previously generated. To plot multiple signals in the same figure, use `hold on`.

`h = scatterplot(...)` is the same as the earlier syntaxes, except that `h` is the handle to the figure that contains the scatter plot.

# scatterplot

---

## Examples

See “Example: Scatter Plots” on page 2-12 or the example on the reference page for `demodmap`. Both examples illustrate how to plot multiple signals in a single scatter plot.

For an online demonstration, use `scattereyedemo`.

## See Also

`eyediagram`, `plot`, `hold`

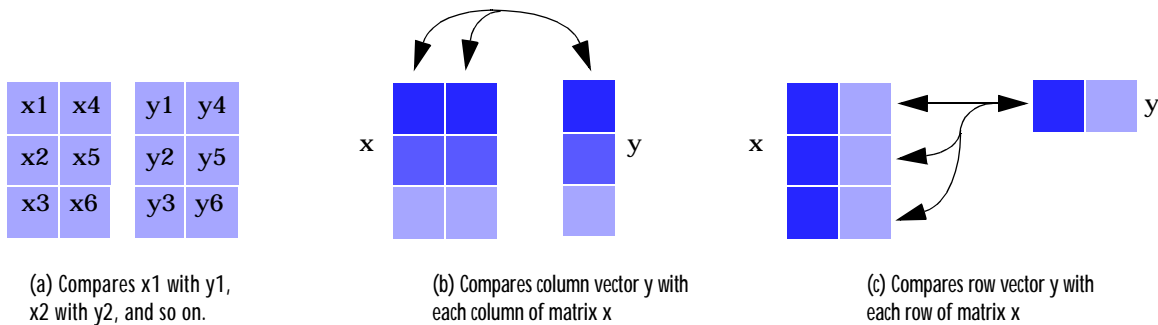
**Purpose** Compute number of symbol errors and symbol error rate

**Syntax**

```
[number, ratio] = symerr(x, y);
[number, ratio] = symerr(x, y, flag);
[number, ratio, loc] = symerr(...)
```

**Description** For All Syntaxes

The `symerr` function compares binary representations of elements in `x` with those in `y`. The schematics below illustrate how the shapes of `x` and `y` determine which elements `symerr` compares.



The output number is a scalar or vector that indicates the number of elements that differ. The size of number is determined by the optional `flag` and by the dimensions of `x` and `y`. The output `ratio` equals number divided by the total number of elements in the *smaller* input.

**For Specific Syntaxes**

`[number, ratio] = symerr(x, y)` compares the elements in `x` and `y`. The sizes of `x` and `y` determine which elements are compared:

- If `x` and `y` are matrices of the same dimensions, then `symerr` compares `x` and `y` element-by-element. number is a scalar. See schematic (a) in the figure.
- If one is a row (respectively, column) vector and the other is a two-dimensional matrix, then `symerr` compares the vector element-by-element with *each row (resp., column)* of the matrix. The length of the vector must equal the number of columns (resp., rows) in the matrix.

number is a column (resp., row) vector whose *m*th entry indicates the number of elements that differ when comparing the vector with the *m*th row (resp., column) of the matrix. See schematics (b) and (c) in the figure.

[number, ratio] = symerr(x, y, flag) is similar to the previous syntax, except that flag can override the defaults that govern which elements symerr compares and how symerr computes the outputs. The values of flag are 'overall', 'column-wise', and 'row-wise'. The table below describes the differences that result from various combinations of inputs. In all cases, ratio is number divided by the total number of elements in y.

Table 3-22: Comparing a Two-Dimensional Matrix x with Another Input y

Shape of y	flag	Type of Comparison	number
Two-dimensional matrix	'overall' (default)	Element-by-element	Total number of symbol errors
	'column-wise'	<i>m</i> th column of x vs. <i>m</i> th column of y	Row vector whose entries count symbol errors in each column
	'row-wise'	<i>m</i> th row of x vs. <i>m</i> th row of y	Column vector whose entries count symbol errors in each row
Column vector	'overall' (default)	y vs. each column of x	Total number of symbol errors
	'column-wise'	y vs. each column of x	Row vector whose entries count symbol errors in each column of x
Row vector	'overall' (default)	y vs. each row of x	Total number of symbol errors
	'row-wise'	y vs. each row of x	Column vector whose entries count symbol errors in each row of x

[number, ratio, loc] = symerr(...) returns a binary matrix loc that indicates which elements of x and y differ. An element of loc is zero if the corresponding comparison yields no discrepancy, and one otherwise.

Examples

On the reference page for bi terr, the last example uses symerr.

The command below illustrates how `symerr` works when one argument is a vector and the other is a matrix. It compares the vector `[1, 2, 3]'` to the columns

$$\begin{bmatrix} 1 \\ 3 \\ 3 \end{bmatrix}, \begin{bmatrix} 1 \\ 2 \\ 3 \end{bmatrix}, \begin{bmatrix} 3 \\ 2 \\ 8 \end{bmatrix}, \text{ and } \begin{bmatrix} 1 \\ 2 \\ 3 \end{bmatrix}$$

of the matrix.

```
num = symerr([1 2 3]', [1 1 3 1; 3 2 2 2; 3 3 8 3])
```

```
num =
```

```
1      0      2      0
```

As another example, the command below illustrates the use of `flag` to override the default row-by-row comparison. Notice that `number` and `ratio` are scalars.

```
format rat; [number, ratio, loc] = symerr([1 2; 3 4], ...  
[1 3], 'overall')
```

```
number =
```

```
3
```

```
ratio =
```

```
3/4
```

```
loc =
```

```
0      1  
1      1
```

## See Also

`biterr`

# syndtable

---

<b>Purpose</b>	Produce syndrome decoding table
<b>Syntax</b>	<code>t = syndtable(parmat);</code>
<b>Description</b>	<p><code>t = syndtable(parmat)</code> returns a decoding table for an error-correcting binary code having codeword length <math>n</math> and message length <math>k</math>. <code>parmat</code> is an <math>(n-k)</math>-by-<math>n</math> parity-check matrix for the code. <code>t</code> is a <math>2^{n-k}</math>-by-<math>n</math> binary matrix. The <math>r</math>th row of <code>t</code> is an error pattern for a received binary codeword whose syndrome has decimal integer value <math>r-1</math>. (The syndrome of a received codeword is its product with the transpose of the parity-check matrix.) In other words, the rows of <code>t</code> represent the coset leaders from the code's standard array.</p> <p>When converting between binary and decimal values, the leftmost column is interpreted as the <i>most</i> significant digit. This differs from the default convention in the <code>bi2de</code> and <code>de2bi</code> commands.</p>
<b>Examples</b>	An example is in the section "Decoding Table" on page 2-34.
<b>See Also</b>	<code>decode</code> , <code>hammgen</code> , <code>gfcosets</code>
<b>References</b>	Clark, George C. Jr. and J. Bibb Cain. <i>Error-Correction Coding for Digital Communications</i> . New York: Plenum Press, 1981.



**Purpose** Convert a vector into a matrix

**Syntax**

```
mat = vec2mat(vec, matcol);  
mat = vec2mat(vec, matcol, padding);  
[mat, padded] = vec2mat(...);
```

**Description** `mat = vec2mat(vec, matcol)` converts the vector `vec` into a matrix with `matcol` columns, creating one row at a time. If the length of `vec` is not multiple of `matcol`, then extra zeros are placed in the last row of `mat`. The matrix `mat` has `ceil(length(vec)/matcol)` rows.

`mat = vec2mat(vec, matcol, padding)` is the same as the first syntax, except that the extra entries placed in the last row of `mat` are not necessarily zeros. The extra entries are taken from the matrix `padding`, in order. If `padding` has fewer entries than are needed, then the last entry is used repeatedly.

`[mat, padded] = vec2mat(...)` returns an integer `padded` that indicates how many extra entries were placed in the last row of `mat`.

---

**Note** `vec2mat` is similar to the built-in MATLAB function `reshape`. However, given a vector input, `reshape` creates a matrix one *column* at a time instead of one row at a time. Also, `reshape` requires the input and output matrices to have the same number of entries, whereas `vec2mat` places extra entries in the output matrix if necessary.

---

## Examples

```
vec = [1 2 3 4 5];  
[mat, padded] = vec2mat(vec, 3)
```

`mat =`

```
     1     2     3  
     4     5     0
```

`padded =`

```
     1
```

# vec2mat

---

```
[mat2, padded2] = vec2mat(vec, 4)
```

```
mat2 =
```

1	2	3	4
5	0	0	0

```
padded2 =
```

```
3
```

```
mat3 = vec2mat(vec, 4, [10 9 8; 7 6 5; 4 3 2])
```

```
mat3 =
```

1	2	3	4
5	10	7	4

## See Also

[reshape](#)

Purpose

Convolutionally decode binary data using the Viterbi algorithm

Syntax

```
decoded = vitdec(code, trellis, tblen, opmode, dectype);
decoded = vitdec(code, trellis, tblen, opmode, 'soft', nsdec);
decoded = vitdec(..., 'cont', ..., initmetric, initstates, initinputs);
[decoded finalmetric finalstates finalinputs] = ...
    vitdec(..., 'cont', ...);
```

Description

`decoded = vitdec(code, trellis, tblen, opmode, dectype)` decodes the vector `code` using the Viterbi algorithm. The MATLAB structure `trellis` specifies the convolutional encoder that produced `code`; the format of `trellis` is described in “Trellis Description of a Convolutional Encoder” on page 2-46 and the reference page for the `istrellis` function. `code` contains one or more symbols, each of which consists of  $\log_2(\text{trellis.numOutputSymbols})$  bits. Each symbol in the vector `decoded` consists of  $\log_2(\text{trellis.numInputSymbols})$  bits. `tblen` is a positive integer scalar that specifies the traceback depth.

The string `opmode` indicates the decoder’s operation mode and its assumptions about the corresponding encoder’s operation. Choices are in the table below.

Table 3-23: Values of `opmode` Input

Value	Meaning
'trunc'	The encoder is assumed to have started at the all-zeros state. The decoder traces back from the state with the best metric.
'term'	The encoder is assumed to have both started and ended at the all-zeros state. The decoder traces back from the all-zeros state.
'cont'	The encoder is assumed to have started at the all-zeros state. The decoder traces back from the state with the best metric. A delay equal to <code>tblen</code> symbols elapses before the first decoded symbol appears in the output.

The string *dectype* indicates the type of decision that the decoder makes, and influences the type of data the decoder expects in code. Choices are in the table below.

Table 3-24: Values of *dectype* Input

Value	Meaning
'unquant'	code contains real input values, where 1 represents a logical zero and -1 represents a logical one.
'hard'	code contains binary input values.
'soft'	For soft-decision decoding, use the syntax below. Note that nsdec is required for soft-decision decoding.

Syntax for Soft Decision Decoding

`decoded = vitdec(code, trellis, tblen, opmode, 'soft', nsdec)` decodes the vector code using soft-decision decoding. code consists of integers between 0 and  $2^{nsdec}-1$ , where 0 represents the most confident 0 and  $2^{nsdec}-1$  represents the most confident 1.

Additional Syntaxes for Continuous Operation Mode

`decoded = vitdec(..., 'cont', ..., initmetric, initstates, initinputs)` is the same as the earlier syntaxes, except that the decoder starts with its state metrics, traceback states, and traceback inputs specified by *initmetric*, *initstates*, and *initinputs*, respectively. Each real number in *initmetric* represents the starting state metric of the corresponding state. *initstates* and *initinputs* jointly specify the initial traceback memory of the decoder; both are `trellis.numStates-by-tblen` matrices. *initstates* consists of integers between 0 and `trellis.numStates-1`. If the encoder schematic has more than one input stream, then the shift register that receives the first input stream provides the least significant bits in *initstates*, while the shift register that receives the last input stream provides the most significant bits in *initstates*. The vector *initinputs* consists of integers between 0 and `trellis.numInputSymbols-1`. To use default values for all of the last three arguments, specify them as `[], [], []`.

`[decoded, final metric, final states, final inputs] = ...`  
`vitdec(..., 'cont', ...)` is the same as the earlier syntaxes, except that the final three output arguments return the state metrics, traceback states, and traceback inputs, respectively, at the end of the decoding process. `final metric` is a vector with `trellis.numStates` elements which correspond to the final state metrics. `final states` and `final inputs` are both matrices of size `trellis.numStates-by-tblen`. The elements of `final states` have the same format as those of `initstates`.

## Examples

The example below encodes random data and adds noise. Then it decodes the noisy code three times to illustrate the three decision types that `vitdec` supports. Notice that for unquantized and soft decisions, the output of `convenc` does not have the same data type that `vitdec` expects for the input code, so it is necessary to manipulate `ncode` before invoking `vitdec`.

```
trell = poly2trellis(3,[6 7]); % Define trellis.
msg = randint(100,1,2,123); % Random data
code = convenc(msg,trell); % Encode.
ncode = rem(code + randerr(200,1,[0 1;.95 .05]),2); % Add noise.
tblen = 3; % Traceback length
% Use hard decisions.
decoded1 = vitdec(ncode,trell,tblen,'cont','hard');
% Use unquantized decisions.
ucode = 1-2*ncode; % +1 & -1 represent zero & one, respectively.
decoded2 = vitdec(ucode,trell,tblen,'cont','unquant');
% Use soft decisions.
% To prepare for soft-decision decoding, map to decision values.
[x,qcode] = quantiz(1-2*ncode,[-.75 -.5 -.25 0 .25 .5 .75],...
[7 6 5 4 3 2 1 0]); % Values in qcode are between 0 and 2^3-1.
decoded3 = vitdec(qcode,trell,tblen,'cont','soft',3);

% Compute bit error rates, using the fact that the decoder
% output is delayed by tblen symbols.
[n1,r1] = biterr(decoded1(tblen+1:end),msg(1:end-tblen));
[n2,r2] = biterr(decoded2(tblen+1:end),msg(1:end-tblen));
[n3,r3] = biterr(decoded3(tblen+1:end),msg(1:end-tblen));
disp(['The bit error rates are: ',num2str([r1 r2 r3])])
```

The bit error rates are:      0.020619      0.020619      0.020619

The example below illustrates how to use the final state and initial state arguments when invoking `vitdec` repeatedly. Notice that `[decoded4; decoded5]` is the same as `decoded6`.

```
trel = poly2trellis(3,[6 7]);
code = convenc(randint(100,1,2,123),trel);
% Decode part of code, recording final state for later use.
[decoded4,f1,f2,f3] = vitdec(code(1:100),trel,3,'cont','hard');
% Decode the rest of code, using state input arguments.
decoded5 = vitdec(code(101:200),trel,3,'cont','hard',f1,f2,f3);
% Decode the entire code in one step.
decoded6 = vitdec(code,trel,3,'cont','hard');
isequal(decoded6,[decoded4; decoded5])

ans =

     1
```

**See Also** `convenc`, `poly2trellis`, `istrellis`

**References** Gitlin, Richard D., Jeremiah F. Hayes, and Stephen B. Weinstein. *Data Communications Principles*. New York: Plenum, 1992.

<b>Purpose</b>	Generate white Gaussian noise
<b>Syntax</b>	<pre>y = wgn(m, n, p); y = wgn(m, n, p, imp); y = wgn(m, n, p, imp, state); y = wgn(..., powertype); y = wgn(..., outputtype);</pre>
<b>Description</b>	<p><code>y = wgn(m, n, p)</code> generates an <i>m</i>-by-<i>n</i> matrix of white Gaussian noise. <i>p</i> specifies the power of <i>y</i> in decibels. The default load impedance is 1 Ohm.</p> <p><code>y = wgn(m, n, p, imp)</code> is the same as the previous syntax, except that <i>imp</i> specifies the load impedance in Ohms.</p> <p><code>y = wgn(m, n, p, imp, state)</code> is the same as the previous syntax, except that <code>wgn</code> first resets the state of MATLAB's normal random number generator <code>randn</code> to the integer state.</p> <p><code>y = wgn(..., powertype)</code> is the same as the previous syntaxes, except that the string <i>powertype</i> specifies the units of <i>p</i>. Choices for <i>powertype</i> are <b>'dB'</b>, <b>'dBm'</b>, and <b>'linear'</b>.</p> <p><code>y = wgn(..., outputtype)</code> is the same as the previous syntaxes, except that the string <i>outputtype</i> specifies whether the noise is real or complex. Choices for <i>outputtype</i> are <b>'real'</b> and <b>'complex'</b>. If <i>outputtype</i> is <b>'complex'</b>, then the real and imaginary parts of <i>y</i> each have a noise power of <i>p</i>/2.</p>
<b>Examples</b>	<p>To generate a column vector of length 100 containing real white Gaussian noise of power 0 dB, use this command:</p> <pre>y1 = wgn(100, 1, 0);</pre> <p>To generate a column vector of length 100 containing complex white Gaussian noise, each component of which has a noise power of 0 dB, use this command:</p> <pre>y2 = wgn(100, 1, 0, 'complex');</pre>
<b>See Also</b>	<code>randn</code> , <code>awgn</code>





## A

- addition in Galois fields 2-97
- admod 3-12
- ademodce 3-16
- A-law companders 2-22
- amod 3-20
- amodce 3-25
- amplitude modulation (AM) sample code
  - for basic example 2-61
  - using demodulation offsets 3-14
  - using filters 2-63
  - using Hilbert filter 3-24
  - using Hilbert filter in baseband simulation 3-27
  - using single and double sidebands 3-22
- amplitude shift keying (ASK) sample code
  - for mapping 2-68
  - to plot waveforms 3-81
  - using passband simulation 3-61
- analog
  - signals, representing 2-59
- analog-to-digital conversion 2-14
- apkconst 3-28
- arbitrary signal constellations 2-72
- arithmetic
  - in Galois fields 2-97
  - sample code
    - using extension fields 2-98
    - using prime fields 2-97
- awgn 3-32

## B

- baseband
  - modulated signal 2-60
  - simulation 2-58

- BCH coding
  - discussion of 2-38
  - generator polynomial for 2-33
  - sample code 3-34
    - for tracking errors 3-71
    - using various coding methods 3-93
  - summary of tools for 2-26
- bchdeco 3-34
- bchenco 3-36
- bchpoly 3-37
- bi2de 3-41
- binary
  - codes 2-25
  - matrix format for messages and codewords 2-27
    - in Reed-Solomon codes 2-28
  - sample code 3-92
  - numbers, order of digits and 2-28
  - vector format for messages and codewords 2-26
    - sample code 3-92
- binary-to-decimal conversion 3-41
- bipolar random numbers 2-4
- bit error rates 2-7
- bi terr 3-43
- bits
  - random 2-5
- block coding 2-24
  - functions 2-26
  - techniques 2-25
    - See also* specific coding techniques
- Bose-Chaudhuri-Hocquenghem (BCH) coding
  - discussion of 2-38
  - generator polynomial for 2-33
  - sample code 3-34
    - for tracking errors 3-71

- using various coding methods 3-93
  - summary of tools for 2-26
- burst errors 2-39

## C

- carrier frequency ( $F_c$ ) 2-58
  - relative to sampling rate 2-58
- carrier signal 2-58
  - initial phase
    - for analog modulation 2-62
    - for digital modulation 2-77
- circle signal constellations 2-71
  - default values for 2-72
  - plotting 2-72
- code generator matrices
  - converting to parity-check matrices 2-41
    - sample code 2-32
  - finding 2-41
  - representing 2-30
- code generator polynomials
  - finding 2-40
  - representing 2-33
- codebooks
  - optimizing 2-18
    - for DPCM 2-21
    - sample code 2-18
    - sample code for DPCM 2-21
  - representing 2-15
- codewords
  - definition 2-26
  - representing 2-26
- coding
  - block 2-24
    - functions 2-26
    - techniques 2-25
  - convolutional 2-43

- examples 2-52
  - features of the toolbox 2-43
  - sample code 2-50
  - using polynomial description 2-43
  - using trellis description 2-46
- source 2-14
  - features of the toolbox 2-14
- compand 3-49
- companders 2-22
  - sample code 2-22
- complex envelope 2-60
- compression
  - of data 2-14
- compressors 2-22
  - sample code 2-22
- conjugate elements of Galois fields 3-103
- constellations, signal 2-70
  - arbitrary 2-72
  - circle 2-71
    - default values for 2-72
    - plotting 2-72
  - hexagonal, sample code 2-73
  - plots
    - interpreting 2-69
  - square 2-71
    - plotting 2-71
  - triangular, sample code 2-73
- constraint length
  - of convolutional code 2-44
- convenc 3-51
- converting
  - analog to digital 2-14
  - binary to decimal 3-41
  - decimal to binary 3-67
  - exponential to polynomial format for Galois field elements 2-94
    - sample code 2-95

- generator matrices to parity-check matrices 2-41
  - sample code 2-32
- polynomial representations 3-125
- polynomial to exponential format for Galois field elements 2-96
  - sample code 2-96
- sampling rates 3-85
- vectors to matrices 3-199
- convolutional coding 2-43
  - examples 2-52
  - features of the toolbox 2-43
  - sample code 2-50
  - using polynomial description 2-43
    - sample code 2-46
  - using trellis description 2-46
- correction vector 2-34
- correlation techniques
  - in demodulation 2-66
- cosets, cyclotomic 3-103
- Costas phase-locked loop
  - for analog modulation 2-62
  - for digital modulation 2-77
- cycl gen 3-53
- cyclic coding
  - discussion of 2-37
  - generator polynomial for 2-33
  - sample code 2-37, 3-93
    - using various coding methods 3-93
  - summary of tools for 2-26
- cyclotomic cosets 3-103
- cycl pol y 3-55

## D

- data compression 2-14
- ddemod 3-57

- ddemodce 3-62
- de2bi 3-67
- decimal format for messages and codewords 2-28
  - in Reed-Solomon coding 2-29
  - sample code 3-92
- decision timing
  - and eye diagrams 2-8
  - sample code
    - for eye diagrams 2-10
    - for scatter plots 2-12
- decode 3-69
- decoding tables
  - representing 2-34
- default
  - Galois field parameters, in error-control coding 2-34
  - primitive polynomials 2-93
- delta modulation 2-19
  - sample code 2-20
  - See also* differential pulse code modulation
- demodmap 3-73
- demodulation
  - definition of 2-56
  - digital 2-66
    - functions for 2-67
  - features of the toolbox 2-58
  - noncoherent 2-77
  - See also* modulation
- diagrams
  - eye 2-8
    - sample code 2-9
  - scatter 2-11
    - sample code 2-12
- differential pulse code modulation (DPCM) 2-19
  - parameters, optimizing 2-21
    - sample code 2-21
  - parameters, representing 2-14

- sample code 2-20
- digital
  - modulation 2-66
    - functions for 2-67
  - signals, representing 2-67
- displaying polynomials 2-100
- distortion
  - from DPCM, reducing 2-21
  - from quantization, reducing 2-18
- division in Galois fields 2-97
- dmod 3-78
- dmodce 3-82
- DPCM 2-19
  - parameters, optimizing 2-21
    - sample code 2-21
  - parameters, representing 2-14
  - sample code 2-20
- dpcmdeco 3-86
- dpcmenco 3-87
- dpcmopt 3-88

## E

- encode 3-89
- error
  - analysis, features of the toolbox 2-3
  - integers 2-5
  - patterns 2-5
  - predictive 2-19
  - rates
    - bit versus symbol 2-8
    - sample code 2-7
- error-control coding
  - base 2 only 2-25
  - features of the toolbox 2-25
  - methods supported in toolbox 2-25
  - terminology and notation 2-26

- with default Galois field parameters 2-34
- error-correction capability
  - of BCH codes 2-41
  - of Hamming codes 2-34
  - of Reed-Solomon codes 2-41
- error-rate analysis 2-7
- expanders 2-22
  - sample code 2-22
- exponential format
  - for Galois field elements 2-90
  - in Reed-Solomon coding 2-29
- eye diagrams 2-8
  - sample code 2-9
- eyedi agram 3-95

## F

- features of the toolbox
  - error-analysis 2-3
  - error-control coding 2-25
  - modulation/demodulation 2-58
  - source coding 2-14
- feedback connection polynomials 2-45
- fields, finite 2-89
  - See also* Galois fields
- filtering data over Galois fields 3-110
- filters
  - designing and applying raised cosine 2-83
  - designing Hilbert transform 2-80
  - designing raised cosine 2-87
    - using after analog demodulation 2-62
      - choosing cutoff frequency 2-63
      - resulting time lag 2-64
    - using after digital demodulation 2-77
  - using raised cosine 2-81
  - using square-root raised cosine 2-85

finite fields 2-89  
     *See also* Galois fields  
 font when using `gfpretty` 3-118  
 format of Galois field elements 2-90  
     simplifying and converting to exponential  
         format 2-96  
         sample code 2-96  
     simplifying and converting to polynomial  
         format 2-94  
         sample code 2-95  
 formatting polynomials 2-100  
 formatting, of signals 2-14  
 frequency modulation (FM)  
     sample code 3-18  
 frequency shift keying (FSK) 2-66  
     sample code 3-85

## G

Galois fields 2-89  
     in error-control coding 2-33  
         avoiding explicit reference to 2-34  
     representing elements of 2-90  
 Gaussian elimination  
     in `gflineq` 3-113  
     in `gfrank` 3-124  
 Gaussian noise, generating 2-3  
`gen2par` 3-97  
 generator matrices  
     converting to parity-check matrices 2-41  
         sample code 2-32  
     finding 2-41  
     representing 2-30  
 generator polynomials  
     finding 2-40  
     for convolutional code 2-44  
     representing 2-33

`gfadd` 3-99  
`gfconv` 3-101  
`gfcosets` 3-103  
`gfdeconv` 3-105  
`gfdi v` 3-108  
`gffilter` 3-110  
`gflineq` 3-112  
`gfminpol` 3-114  
`gfmul` 3-116  
`gfplus` 3-117  
`gfpretty` 3-118  
`gfprimck` 3-120  
`gfpri mdf` 3-121  
`gfpri mfd` 3-122  
`gfrank` 3-124  
`gfrepconv` 3-125  
`gfroots` 3-126  
`gfsub` 3-128  
`gftrunc` 3-130  
`gftuple` 3-131  
`gfweight` 3-134

## H

`hamngen` 3-135  
 Hamming coding 2-39  
     for single-error-correction 2-34  
     sample code 2-34  
         using various coding methods 3-93  
         using various formats 3-92  
     summary of tools for 2-26  
 Hamming weight 3-134  
`hank2sys` 3-138  
 hard-decision decoding 2-50  
 Hilbert filters  
     designing 2-80  
     for amplitude modulation 2-62

- sample code 3-24
- for baseband amplitude modulation
  - sample code 3-27
- hilbair 3-140

**I**

- initial phases of carrier signal
  - for analog modulation 2-62
  - for digital modulation 2-77
- in-phase/quadrature coordinates 2-69
- inverses, multiplicative in Galois fields 3-108
- irreducible polynomials 2-101
  - sample code 2-101
- istrellis 3-143

**L**

- length of polynomial representations
  - minimizing 2-100
- linear block coding 2-35
  - sample code 2-36
- linear equations over Galois fields 3-112
  - and filtering 3-110
- linear predictors 2-19
  - optimizing 2-21
    - sample code 2-21
  - representing 2-19
- list of elements of Galois fields 2-91
  - generating 2-96
    - sample code 2-92
  - in error-control coding, avoiding explicit
    - reference to 2-34
- Lloyd algorithm, for optimizing quantization
  - parameters 2-18
- lloyds 3-145

**M**

- mapped signals
  - representing 2-67
    - for PSK and QASK 2-69
- mapping
  - functions for 2-67
  - without modulating 2-76
- marcumq 3-148
- matrices over Galois fields
  - rank 3-124
- messages
  - definition 2-26
  - representing
    - for coding functions 2-26
    - for modulation functions 2-67
- minimal polynomials of Galois field elements
  - 2-101
    - sample code 2-101
- minimum distance 3-134
- minimum shift keying (MSK) 2-66
- modmap 3-149
- modulated signals, representing 2-68
- modulation 2-56
  - definition of 2-56
  - delta 2-19
    - sample code 2-20
    - See also* differential pulse code modulation
  - differential pulse code. *See* differential pulse code modulation
  - digital 2-66
    - functions for 2-67
  - features of the toolbox 2-58
  - methods supported in toolbox 2-57
  - of several signals 2-60
    - sample code 3-18
  - terminology 2-58
  - without mapping 2-76

- mu-law companders 2-22
  - sample code 2-22
- multiple roots of polynomials over Galois fields 3-126
- multiplication in Galois fields 2-97
- multiplicative inverses in Galois fields 3-108

## N

- nonbinary codes 2-25
  - Reed-Solomon 2-39
- noncausality of filters 2-78
- noncoherent demodulation 2-77
- Nyquist sampling theorem 2-58

## O

- oct2dec 3-154
- optimizing
  - DPCM parameters 2-21
    - sample code 2-21
  - quantization parameters 2-18
    - sample code 2-18
- order of digits in binary numbers 2-28
- order, predictive 2-19

## P

- parity-check matrices
  - finding 2-41
  - representing 2-31
- partitions
  - optimizing 2-18
    - for DPCM 2-21
    - sample code 2-18
    - sample code for DPCM 2-21
  - representing 2-14

- passband
  - simulation 2-58
- phase modulation (PM)
  - sample code 2-64
- phase shift keying (PSK) sample code
  - for basic example 2-74
  - for demapping 3-76
  - for mapping 2-69
  - for plotting signal constellation 3-152
- phase-locked loop, Costas
  - for analog modulation 2-62
  - for digital modulation 2-77
- points, decision
  - and eye diagrams 2-8
  - sample code
    - for eye diagrams 2-10
    - for scatter plots 2-12
- poly2trellis 3-155
- polynomial
  - description of an encoder
    - sample code 2-46
  - description of encoder 2-43
  - format for Galois field elements 2-91
- polynomials
  - displaying formatted 2-100
  - generator, finding 2-40
- polynomials over Galois fields 2-99
  - adding 2-100
  - dividing 2-100
  - irreducible 2-101
  - minimal 2-101
    - sample code 2-101
  - multiplying 2-100
- primitive 2-101
  - default 2-93
  - definition 2-89
  - finding 3-122

- representation choices for 3-125
- roots of 2-102
  - sample code 2-102
- subtracting 2-100
- truncating 2-100
- predictive
  - error 2-19
  - order 2-19
- predictive quantization 2-19
  - features of the toolbox 2-14
  - parameters, optimizing 2-21
    - sample code 2-21
  - parameters, representing 2-14
  - sample code 2-20
- predictors 2-19
  - linear 2-19
  - optimizing 2-21
    - sample code 2-21
  - representing 2-19
- primitive
  - element, definition 2-89
  - polynomials 2-101
    - consistent use of 2-93
    - default 2-93
    - definition 2-89
    - finding 3-122
    - in error-control coding, avoiding explicit reference to 2-34
    - sample code 2-101
- punctured convolutional code 2-54
- Q**
- QAM
  - representing signals for 2-60
  - sample code 2-60
- QASK sample code
  - using eye diagram 2-9
  - using scatter plot and square constellation 2-12
- qaskdeco 3-158
- qaskenco 3-160
- quadrature amplitude modulation (QAM)
  - representing signals for 2-60
  - sample code 2-60
- quadrature amplitude shift keying (QASK) sample code
  - using eye diagram 2-9
  - using scatter plot and square constellation 2-12
- quantiz 3-163
- quantization 2-14
  - coding 2-17
  - DPCM parameters, optimizing 2-21
    - sample code 2-21
  - features of the toolbox 2-14
  - parameters, optimizing 2-18
    - sample code 2-18
  - parameters, representing 2-14
  - predictive 2-19
    - sample code 2-20
  - sample code 2-15
  - vector 2-14
- R**
- raised cosine filters
  - designing and applying 2-83
  - designing but not applying 2-87
  - filtering with 2-81
  - square-root 2-85
- randerr 3-165
- randint 3-167
- random
  - bipolar symbols 2-4
  - bits 2-5



- in error patterns 2-5
- integers 2-5
- signals 2-3
  - features of the toolbox 2-3
- symbols 2-4
- `randsrc` 3-168
- rank of matrices over Galois fields 3-124
- `rcosfir` 3-170
- `rcosflt` 3-172
- `rcosfir` 3-175
- `rcosine` 3-177
- redundancy, reducing 2-14
- Reed-Solomon coding
  - discussion of 2-39
  - generator polynomial for 2-33
  - summary of tools for 2-26
- references
  - convolutional coding 2-55
  - error-control coding 2-42
  - Galois fields 2-103
  - modulation/demodulation 2-77
- representation of polynomials, shortening 2-100
- representing
  - analog signals 2-59
  - codebooks 2-15
  - codewords 2-26
  - decoding tables 2-34
  - digital signals 2-67
  - Galois field elements 2-90
    - in a list 2-91
    - in exponential format 2-90
    - in polynomial format 2-91
  - generator matrices 2-30
  - generator polynomials 2-33
  - mapped signals 2-67
    - for PSK and QASK 2-69
  - messages

- for coding functions 2-26
  - for modulation functions 2-67
- modulated signals 2-68
- parity-check matrices 2-31
- partitions 2-14
- predictors 2-19
- quantization parameters 2-14
- signal constellations 2-70
- roots of polynomials over Galois fields 2-102
  - sample code 2-102
- `rsdeco` 3-179
- `rsdecode` 3-182
- `rsdecof` 3-184
- `rsenco` 3-185
- `rsencode` 3-188
- `rsencof` 3-190
- `rspoly` 3-191

## S

- sampling rate 2-58
  - change during mapping 2-66
  - individual 2-70
  - of signals, relative 2-67
  - relative to carrier frequency 2-58
  - significance of 2-70
- scalar quantization
  - coding 2-17
  - features of the toolbox 2-14
  - parameters, representing 2-14
  - sample code 2-15
- scatter plots 2-11
  - sample code 2-12
  - using modulation 2-74
- `scatterplot` 3-193
- signal constellations 2-70
  - arbitrary 2-72

- circle 2-71
  - default values for 2-72
  - plotting 2-72
- hexagonal, sample code 2-73
- plots
  - interpreting 2-69
- square 2-71
  - plotting 2-71
- triangular, sample code 2-73
- signal formatting 2-14
  - features of the toolbox 2-14
- Signal Processing Toolbox
  - for filter design 2-62
- simplifying
  - exponential format of Galois field elements 2-96
    - sample code 2-96
  - polynomial format of Galois field elements 2-94
    - sample code 2-95
- soft-decision decoding 2-51
  - sample code 2-51
- source coding 2-14
  - features of the toolbox 2-14
- square signal constellations 2-71
  - plotting 2-71
- subtraction in Galois fields 2-97
- symbol error rates 2-7
- symerr 3-195
- syndrome 2-34
- syndtable 3-198

## T

- terminology
  - modulation/demodulation 2-58
- timing, decision

- and eye diagrams 2-8
- sample code
  - for eye diagrams 2-10
  - for scatter plots 2-12
- training data
  - for optimizing DPCM quantization parameters 2-21
  - for optimizing quantization parameters 2-18
- trellis
  - description of encoder 2-46
  - structure 2-47
    - sample code 2-49
- truncating polynomials over Galois fields 2-100

## V

- vec2mat 3-199
- vector quantization 2-14
- vi tdec 3-201

## W

- weight, Hamming 3-134
- wgn 3-205
- white Gaussian noise, generating 2-3