Lecture 23
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## Differential Encoding

Suppose we have a signal that is slowly varying. For instance, if we were looking at a video frame by frame we would see that only a few pixels are changing between subsequent frames. In this case, rather than encoding the signal as is, we would first sample it and look at the difference signal and encode this instead:


This $d_{n}$ would then need to be quantized, creating an estimate $\left(\hat{d_{n}}\right)$ which would contain some quantization noise $\left(q_{n}\right)$

$$
\begin{gathered}
Q\left(d_{n}\right)=\hat{d}_{n} \\
\hat{d}_{n}=d_{n}+q_{n}
\end{gathered}
$$

What we are truly interested in recovering are the values of $x$. Unfortunately, the quantization error will get accumulated in the value of $x$. The reason being, that the operation forming $d_{n}$ is of the following matrix form

$$
\left(\begin{array}{cccccc}
1 & 1 & 0 & 0 & 0 & 0 \\
0 & -1 & 1 & 0 & 0 & 0 \\
0 & 0 & -1 & 1 & 0 & 0 \\
\vdots & \vdots & 0 & \ddots & \ddots & 0 \\
0 & 0 & 0 & 0 & -1 & 1
\end{array}\right)
$$

Inverting this matrix we obtain the accumulator matrix

$$
\left(\begin{array}{cccccc}
1 & 1 & 0 & 0 & 0 & 0 \\
1 & 1 & 1 & 0 & 0 & 0 \\
1 & 1 & 1 & 1 & 0 & 0 \\
\vdots & \vdots & \vdots & \vdots & \ddots & 0 \\
1 & 1 & 1 & 1 & \ldots & 1
\end{array}\right)
$$

If you make an error in the $d_{n}$ the noise in $x_{n}$ will be accumulated as follows

$$
\hat{x}_{n}=x_{0}+\sum_{i=0}^{n} q_{n}
$$

The quantization noise $q_{n}$ can be positive or negative and in the long term we expect it to add up to zero, but what happens with high probability is that the range of the error becomes too great for us to handle. Therefore, we adopt the following strategy

$$
d_{n}=x_{n}-\hat{x}_{n-1}
$$

We can then implement this technique using the following encoder


The corresponding decoder will look as follows


This encoder/decoder scheme shown is governed by the relationship introduced previously. Rearranged it looks as follows

$$
\hat{x_{n}}=\hat{d_{n}}+\hat{x}_{n-1}
$$

Walking through an example using this scheme we would have the following sequence

$$
\begin{align*}
\hat{x_{0}} & =x_{0}  \tag{1}\\
\hat{x_{1}} & =\hat{d_{1}}+\hat{x_{0}}  \tag{2}\\
& =d_{1}+q_{1}+\hat{x_{0}}  \tag{3}\\
& =x_{1}+q_{1}  \tag{4}\\
\hat{x_{2}} & =\hat{d_{2}}+\hat{x_{1}}  \tag{5}\\
& =d_{2}+q_{2}+\hat{x_{1}}  \tag{6}\\
& =x_{2}+q_{2} \tag{7}
\end{align*}
$$

We see that in general

$$
\begin{aligned}
\hat{x_{n}} & =\hat{d}_{n}+\hat{x}_{n-1} \\
& =d_{n}+q_{n}+\hat{x}_{n-1} \\
& =x_{n}+q_{n}
\end{aligned}
$$

Notice that while the quantization noise $q_{n}$ was originally accumulated in $x_{n}$, by adopting the aforementioned strategy we have made each estimate of $x_{n}$ dependent on only its own respective quantization error. This type of method is known as a Differential Pulse Coded Modulation Scheme (DPCM).

## Differential Pulse Coded Modulation Scheme

Generalizing the method described above, we see that we use a previous predicted value of $x_{n}$, denoted by $p_{n}$, to construct the difference sequence which is then quantized and used to predict the current value of $x_{n}$. Furthermore, the predicted $x_{n}$ is operated on by a Predictor $(P)$ which provides an estimate of the previous $x_{n}$ in order to recursively find the difference signal. This is shown in the respective encoder and decoder figures below


Formalizing this procedure, we write

$$
\begin{aligned}
p_{n} & =f\left(\hat{x}_{n-1}, x_{n-2}, x_{n-3}, \ldots, \hat{x_{0}}\right) \\
d_{n} & =x_{n}-p_{n} \\
& =x_{n}-f\left(\hat{x}_{n-1}, x_{n-2}, x_{n-3}, \ldots, \hat{x_{0}}\right)
\end{aligned}
$$

We are hopeful that the values of $d_{n}$ are much smaller than the values of $x_{n}$, given that it is a difference signal for a slowly varying signal. So if we were looking at the energy in $d_{n}$ we would guess it to be much smaller than the energy in $x_{n}$. When we are dealing with the energy of a signal it is analogous to its variance, which we define as follows
So if we were to optimize something in this DPCM scheme, we would want to find $f$ such that we minimize the energy in $d_{n}$

Find $f\left(\hat{x}_{n-1}, x_{n-2}, x_{n-3}^{\hat{n}}, \ldots, \hat{x_{0}}\right)$ such that $\sigma_{d}{ }^{2}$ is minimized.

$$
\sigma_{d}^{2}=E\left[\left(x_{n}-f\left(\hat{x}_{n-1}, x_{n-2}, x_{n-3}^{\hat{n}}, \ldots, \hat{x_{0}}\right)\right)^{2}\right]
$$

Since it is necessary to know previous estimates of $x_{n}$ in order to calculate $f$, but we also need to know $f$ in order to calculate previous estimates, we find ourselves in a roundabout and difficult situation. Therefore, we make the following assumption. Suppose,

$$
p_{n}=f\left(x_{n-1}, x_{n-2}, x_{n-3}, \ldots, x_{0}\right)
$$

In other words, we design the predictor assuming there is no quantization. This is called a Fine Quantization assumption. In many cases this makes sense because the estimate of $x_{n}$ is very close to $x_{n}$, as the difference signal is very small, and the quantization error is even smaller. Now we have

$$
\sigma_{d}^{2}=E\left[\left(x_{n}-f\left(x_{n-1}, x_{n-2}, x_{n-3}, \ldots, x_{0}\right)^{2}\right]\right.
$$

Considering the limitations on $f$, we note that $f$ can be any nonlinear function. But we often don't know a way to implement many nonlinear functions. So we assume further that the predictor $f$ is linear and seek the best linear function. By definition, a linear function expressing $f\left(x_{n-1}, x_{n-2}, x_{n-3}, \ldots, x_{0}\right)$ will merely be a weighted sum of previous values of $x_{n}$. Assuming we are looking back through a "window" at the first $N$ samples of a signal, the predictor can then be expressed as follows

$$
p_{n}=\sum_{i=1}^{N} a_{i} x_{n-i}
$$

Now the variance of $d_{n}$ becomes

$$
\sigma_{d}^{2}=E\left[\left(x_{n}-\sum_{i=1}^{N} a_{i} x_{n-i}\right)^{2}\right]
$$

Now we need to find the coefficients $a_{i}$ which will minimize this variance. To do this we take the derivative with respect to $a_{i}$ and set it equal to zero.

$$
\begin{aligned}
\frac{\partial \sigma_{d}^{2}}{\partial a_{1}} & =E\left[-2\left(x_{n}-\sum_{i=1}^{N} a_{i} x_{n-i}\right) x_{n-1}\right] \\
& =-2 E\left[x_{n} x_{n-1}-\sum_{i=1}^{N} a_{i} x_{n-i} x_{n-1}\right]=0 \\
& \vdots \\
\frac{\partial \sigma_{d}^{2}}{\partial a_{j}} & =2 E\left[x_{n} x_{n-j}-\sum_{i=1}^{N} a_{i} x_{n-i} x_{n-j}\right]=0
\end{aligned}
$$

By setting the derivative above to zero and rearranging, we see that the coefficients we are looking for are dependent upon second order statistics

$$
E\left[x_{n} x_{n-j}\right]=\sum_{i=1}^{N} a_{i} E\left[x_{n-i} x_{n-j}\right]
$$

We make the assumption that $x$ is stationary, wich means that the correlation between two values of $x_{n}$ is only a function of the lag between them. Formally, we denote this by the fact that the expectation of the product of two values of $x_{n}$ separated in time by $k$ samples (also known as the autocorrelation) is purely a function of the time difference, or lag, $k$

$$
E\left[x_{n} x_{n+k}\right]=R_{x x}(k)
$$

The relationship governing the coefficients $a_{i}$ can then be rewritten using this notation as

$$
R_{x x}(j)=\sum_{i=1}^{N} a_{i} R_{x x}(i-j) \text { for } 1 \leq j \leq N
$$

Thus, looking at the autocorrelation functions for each $j$ we have

$$
\begin{gathered}
R_{x x}(1)=\sum_{i=1}^{N} a_{i} R_{x x}(i-1) \\
R_{x x}(2)=\sum_{i=1}^{N} a_{i} R_{x x}(i-2) \\
\vdots \\
R_{x x}(N)=\sum_{i=1}^{N} a_{i} R_{x x}(i-N)
\end{gathered}
$$

Now that we have $N$ equations and $N$ unknowns we can solve for the coefficients $a_{i}$ in matrix form

$$
\left(\begin{array}{cccc}
R_{x x}(0) & R_{x x}(1) & \ldots & R_{x x}(N-1) \\
R_{x x}(1) & R_{x x}(0) & & \vdots \\
\vdots & & \ddots & \vdots \\
R_{x x}(N-1) & & \ldots & R_{x x}(0)
\end{array}\right)\left(\begin{array}{c}
a_{1} \\
a_{2} \\
\vdots \\
a_{n}
\end{array}\right)=\left(\begin{array}{c}
R_{x x}(1) \\
R_{x x}(2) \\
\vdots \\
R_{x x}(N)
\end{array}\right)
$$

Rewriting this more compactly where $R$ is a matrix and $a$ and $P$ are column vectors we have

$$
\begin{gathered}
R a=P \\
a=R^{-1} P
\end{gathered}
$$

Thus, we see in order to determine the coefficients for a linear predictor for DPCM we must invert the autocorrelation matrix and multiply by the autocorrelation vector $P$. Once we have determined the coefficients we can design the predictor necessary for our DPCM scheme.

Now suppose that our signal is rapidly changing as shown below


We see that since the signal varies significantly over short intervals the difference signal $d_{n}$ would be very large and DPCM might not be a very good scheme. Suppose instead that we passed this rapidly changing signal through both a low pass and high pass filter as shown below


Low Pass


High Pass

Let us create two signals based off the values of $x_{n}$ to emulate this low pass and high pass response. Let $y_{n}$ represent an averaging operation that smoothes out the response of $x_{n}$ (low pass) and let $z_{n}$ represent a difference operation which emulates the high frequency variation of $x_{n}$ (high pass).

$$
\begin{aligned}
& y_{n}=\frac{x_{n}+x_{n-1}}{2} \\
& z_{n}=\frac{x_{n}-x_{n-1}}{2}
\end{aligned}
$$

Applying this method for each $x_{n}$ we would send or store two values ( $y_{n}$ and $z_{n}$ ). This is unnecessary. Instead, what we can do is apply the following strategy

$$
\begin{aligned}
& y_{2 n}=\frac{x_{2 n}+x_{2 n-1}}{2} \\
& z_{2 n}=\frac{x_{2 n}-x_{2 n-1}}{2}
\end{aligned}
$$

Then we can recover both even and odd values of $x_{n}$ as follows

$$
\begin{gathered}
y_{2 n}+z_{2 n}=x_{2 n} \\
y_{2 n}-z_{2 n}=x_{2 n-1}
\end{gathered}
$$

This process of splitting signal components into multiple portions is called decimation and can be extrapolated out further until a point at which you perform bit allocation. This method is called sub-band coding. We note that DPCM is well suited for the $y_{n}$ low pass components whereas another technique would likely suite the $z_{n}$ high pass components more effectively.

## Distributed Storage

Today's storage systems often utilize distributed storage. In such a system, data will be stored in many separate locations on servers. Suppose we want to do data compression in such a system. What we would like to do is query a portion of our data set - this could be 1 page out of an entire document for instance. Rather than having to sift through the entire data set to find 1 page, we would like to be able to go directly there. In other words, we would like such a system to be query efficient. For this reason suppose we divide data out into sections, such as by page, before compressing it in an effort to preserve information about where the data came from. Such a partitioning of a data set is shown below

$$
(101|010| 011|\ldots| \ldots)
$$

Next we define query efficiency as the \# of bits we need to process to get back a single bit. Suppose that the partitioned bit stream shown above has length $N$ and we have partitioned it in chunks of $m=3$ bits. In this instance, the query efficiency would be m .

Suppose we have a binary vector of length n represented by x which we can compress to $H(x)+1$. Our compression would then be

$$
\begin{aligned}
& \text { Compression Rate }=\frac{H\left(x_{1}^{n}\right)+1}{n} \\
& =\frac{H\left(x_{1}^{n}\right)}{n}+\frac{1}{n} \\
& =\frac{n H(x)}{n}+\frac{1}{n} \\
& =H(x)+\frac{1}{n}
\end{aligned}
$$

Now, in the case of our previous example regarding query efficiency. The compression rate will be

$$
\begin{aligned}
\text { Rate } & =H(x)+\frac{1}{m} \\
& =H(x)+\frac{1}{\text { query efficiency }}
\end{aligned}
$$

If we were able to compress the file as a whole the query efficiency would be huge and the Compression Rate would be

$$
\begin{aligned}
& \text { Rate }=H(x)+\frac{1}{N} \\
& \text { Difference in rate from optimal } \stackrel{=}{=} \frac{1}{\text { Query Efficiency }}
\end{aligned}
$$

Thus we see there is a trade off between Compression Rate and Query Efficiency. Let us end the lecture by propose a research problem were we find a better relation between query efficiency and redundancy.

